

# **HUMPHREY IMPOUNDMENT SEMI-ANNUAL 1<sup>ST</sup> HALF 2025 INTERIM MEASURE REPORT**

**TRADEPOINT ATLANTIC  
SPARROWS POINT, MARYLAND**

Prepared for:



**TRADEPOINT ATLANTIC**  
6995 Bethlehem Blvd  
Sparrows Point, MD 21219

Prepared by:



**ARM GROUP LLC**  
9175 Guilford Road  
Suite 310  
Columbia, Maryland 21046  
  
ARM Project No. 21010214

Respectfully Submitted:

A handwritten signature in black ink, appearing to read "Charlotte Norris".

Charlotte Norris  
Staff Geologist

A handwritten signature in black ink, appearing to read "Kaye Guille".

Kaye Guille, P.E., PMP  
Senior Engineer

Revision 0 - July 15, 2025

**TABLE OF CONTENTS**

|            |                                                     |          |
|------------|-----------------------------------------------------|----------|
| <b>1.0</b> | <b>Introduction.....</b>                            | <b>1</b> |
| 1.1.       | Site Description and History .....                  | 1        |
| <b>2.0</b> | <b>Interim Measure Results .....</b>                | <b>3</b> |
| 2.1.       | Groundwater Sampling.....                           | 3        |
| 2.1.1.     | 1 <sup>st</sup> Half 2025 Groundwater Sampling..... | 3        |
| 2.1.2.     | Groundwater Sampling Trends .....                   | 4        |
| 2.1.3.     | B24 Groundwater Sampling.....                       | 4        |
| 2.2.       | Surface Water Sampling.....                         | 5        |
| 2.3.       | Groundwater Elevation.....                          | 5        |
| 2.4.       | NAPL Observations .....                             | 5        |
| <b>3.0</b> | <b>Summary and Conclusions.....</b>                 | <b>7</b> |

**List of Figures (Following Text)**

|          |                                                               |
|----------|---------------------------------------------------------------|
| Figure 1 | Parcel B14 Site Location                                      |
| Figure 2 | Monitoring Well and Surface Water Sampling Locations          |
| Figure 3 | Groundwater PAL Exceedances – 1 <sup>st</sup> Half 2025       |
| Figure 4 | Historic Groundwater Results for Benzene                      |
| Figure 5 | Historic Groundwater Results for Naphthalene                  |
| Figure 6 | Groundwater Elevation Contour Map – 1 <sup>st</sup> Half 2025 |

**List of Tables (Following Text)**

|         |                                                                          |
|---------|--------------------------------------------------------------------------|
| Table 1 | Summary of Organics Detected in Groundwater – Parcel B14 1st Half 2025   |
| Table 2 | Historic Groundwater Sampling Results                                    |
| Table 3 | Summary of Organics Detected in Groundwater – Parcel B24 1st Half 2025   |
| Table 4 | Summary of Organics Detected in Surface Water – Parcel B14 1st Half 2025 |
| Table 5 | Historic Surface Water Sampling Results                                  |
| Table 6 | Parcel B14 Groundwater Elevations                                        |
| Table 7 | Parcel B14 NAPL Gauging                                                  |

**List of Appendices (Following Text)**

|            |                       |
|------------|-----------------------|
| Appendix A | Laboratory Reports    |
| Appendix B | Mann-Kendall Analysis |

## 1.0 INTRODUCTION

ARM Group LLC (ARM), on behalf of Tradepoint Atlantic (TPA), has prepared this Semi-Annual 1<sup>st</sup> Half 2025 Interim Measures (IM) Report for a portion of TPA property (formerly Sparrows Point Terminal, LLC) that has been designated as the Humphrey Impoundment located in Area B: Parcel B14 (the Site). Parcel B14 is comprised of 60.3 acres of the approximately 3,100-acre former steel making facility (**Figure 1**). The majority of Parcel B14 is occupied by the Humphrey Impoundment, which is approximately 43 acres in size. The Humphrey Impoundment has been identified as a Special Study Area due to the wastes which were historically managed within the area, and potential environmental releases which could have occurred due to its construction (slag base and sides). *The Work Plan for the Humphrey Impoundment* (Revision 2 dated March 29, 2022) was submitted to the Maryland Department of the Environment (MDE) and United States Environmental Protection Agency (USEPA), hereby referred to as the Agencies. The IM Work Plan for the Humphrey Impoundment was proposed as a component of the final remedy for Parcel B14 for the impoundment and the wastes contained within the impoundment. The IM Work Plan proposed groundwater and surface water sampling and gauging to monitor the Site conditions as the Humphrey Impoundment is filled, in preparation for development.

This 1<sup>st</sup> Half 2025 IM Report summarizes groundwater and surface water sampling and gauging results from January through June 2025. The approved IM proposed gauging and sampling of groundwater and surface water around the perimeter and in the vicinity of the Humphrey Impoundment to assess the interaction of the impoundment with groundwater and surface water and to monitor the effectiveness of the proposed IM at reducing migration of contaminants to groundwater and the potential need for additional contingent groundwater remedies. These gauging and sampling activities include the adjacent areas of Parcel B24 and portions of the Tin Mill Canal (TMC).

### 1.1. SITE DESCRIPTION AND HISTORY

In preparation for development, filling activities for the Humphrey Impoundment commenced in January 2022 and have continued through 2025. The impoundment is being filled in quadrants. Filling of the impoundment is now approximately 95% complete and paving is approximately 50% complete. Future use for the impoundment will include a Stormwater Management Regional Pond and a parking lot.

From the late 1800s until 2012, the production and manufacturing of steel was conducted at Sparrows Point. Iron and steel production operations and processes at Sparrows Point included raw material handling, coke production, sinter production, iron production, steel production, and semi-finished and finished product preparation. In 1970, Sparrows Point was the largest steel facility in the United States, producing hot and cold rolled sheets, coated materials, pipes, plates, and rod and wire. The steel making operations at Sparrows Point ceased in fall 2012.

The majority of Parcel B14, as shown in **Figure 1**, is occupied by the Humphrey Impoundment, which is approximately 43 acres in size. As stated in the Description of Current Conditions (DCC) Report prepared by Rust Environment and Infrastructure, dated January 1998, the USEPA identified the Humphrey Impoundment as a potential concern due to the wastes which were historically managed within the impoundment, and potential environmental releases which could have occurred due to its construction (slag base and sides).

Between 1950 and 1970, Humphrey Creek existed as open water (the impoundment did not yet exist) and received wastewater from various steel processing areas including the Hot Strip Mill, Cold Sheet Mill, Tin Mill, and Rod & Wire Mill. Following construction of the TMC (ca. 1969), from 1970 to 1985 the Humphrey Impoundment was used as a dewatering area for on-site sludges and slurry materials generated from the Basic Oxygen Furnace and various on-site water treatment plants. Materials that were dewatered within the impoundment included: Blast Oxygen Furnace slurry; Blast Furnace G, H, J, K, and L thickener sludges; Humphrey Creek Wastewater Treatment Plant (HCWWTP) sludge; Sinter Plant slurry; Open Hearth (No.4) slurry; waste oil pit sludge and non-recoverable waste oil residue; and pre-limer clarifier sludge. Since 1985, the impoundment was used for sludge/slurry dewatering in emergency scenarios only (i.e., when upsets had occurred in the on-site water treatment systems). The MDE was notified prior to these emergency uses. According to the DCC Report, all the wastes that were placed inside the impoundment were determined to be non-hazardous.



## 2.0 INTERIM MEASURE RESULTS

### 2.1. GROUNDWATER SAMPLING

In December 2021, perimeter monitoring wells were installed at Humphrey Impoundment. Starting in January 2022, multiple rounds of monthly groundwater sampling data were collected in accordance with the *Humphrey Impoundment IM Work Plan* (Revision 2 dated March 29, 2022). On November 7, 2022, the Agencies approved a reduction in sampling frequency from monthly to quarterly. On October 30, 2023, the Agencies approved a reduction in gauging frequency from monthly to quarterly. On September 13, 2024, the Agencies approved an additional reduction in groundwater gauging, groundwater sampling, and surface water sampling frequency from quarterly to semi-annually.

As shown on **Figure 2**, groundwater samples were originally collected from 18 monitoring wells around the perimeter of the impoundment. These wells are HI10-MWS, HI11-MWS, HI12-MWS, HI13-MWS, HI14-MWS, HI15-MWS, HI16-MWS, HI17-MWS, HI18-MWS, HI19-MWS, HI20-MWS (replaced with HI20R2-MWS), HI21-MWS, HI22-MWS, TM04-PZM006, TM08R-PZM007, HI02-PZM006 (replaced with HI02R-PZM006), HI04-PZM006 (now abandoned), and HI07-PZM005 (replaced with HI07R-PZM005).

Groundwater samples were analyzed for volatile organic compounds (VOCs) via the USEPA method 8260 and semi-volatile organic compounds (SVOCs) via USEPA method 8270 during the groundwater monitoring events. Groundwater analytical results for the monitoring well sampling are included in **Table 1** and **Table 2**. Project Action Limit (PAL) exceedances for each of the sampling rounds are shown on **Figure 3**.

The groundwater PALs are based on the USEPA Maximum Contaminant Level or the USEPA Regional Screening Level for tap water. The laboratory reports are included in **Appendix A**.

#### 2.1.1. 1<sup>st</sup> Half 2025 Groundwater Sampling

During the 1<sup>st</sup> Half 2025 sampling event (conducted between April 8 and April 15<sup>th</sup>), 16 monitoring wells were sampled. No sample was collected from HI17-MWS (dry), and HI04-PZM006 is abandoned.

Benzene exceedances of the groundwater PAL (5 micrograms per liter, or µg/L) were observed in two baseline monitoring wells: HI19-MWS (12 µg/L) and TM04-PZM006 (64 µg/L). Measured benzene concentrations are consistent with historical values observed at these locations.

Naphthalene exceedances of the groundwater PAL (0.12 µg/L) were observed at 14 of 16 baseline monitoring wells sampled, with the maximum concentration observed in HI19-MWS (52 µg/L). Naphthalene concentrations are consistent with historical values. Historically, the maximum naphthalene concentrations for this study area have been detected at TM04-PZM006. However, naphthalene concentrations have been fluctuating at TM04-PZM006 over the last several quarters:

13 µg/L during the 2<sup>nd</sup> quarter 2024, 130 µg/L during the 2<sup>nd</sup> Half 2024, and 0.16 µg/L during the most recent sampling event. Concentrations at TM04-PZM006 will continue to be monitored for trend changes.

Benz[a]anthracene exceedances of the groundwater PAL (0.03 µg/L) were observed in four of the 16 monitoring wells samples, with the maximum concentration observed in HI12-MWS (0.09 µg/L). Benz[a]anthracene concentrations are consistent with historical values observed at these locations. There were also exceedances for bis(2-Ethylhexyl)phthalate (PAL of 6 µg/L) in one well (concentration of 8.2 µg/L in HI15-MWS) and pentachlorophenol (PAL of 1 µg/L) in two wells (concentrations of 8.0 J µg/L and 6.3 J µg/L in HI07R-MWS and HI18-MWS respectively).

### 2.1.2. Groundwater Sampling Trends

Benzene and naphthalene historic groundwater sampling results are compiled in **Table 2**. Trend plots for benzene and naphthalene (for wells where PAL exceedances have been identified in one or more monitoring events) are included as **Figure 4** and **Figure 5**, respectively. Overall, benzene and naphthalene concentrations in the wells do not appear to be changing as the impoundment is being filled.

As shown on **Figure 4**, monitoring well location TM04-PZM006 continues to be the main area of benzene groundwater impact. Historically, groundwater sampling results for this area are available from December 2001, July 2004, October 2017, and October 2021. A Mann-Kendall analysis of the benzene and naphthalene groundwater concentrations was conducted via the ChemStat 6.5 statistical software from 2001 – April 2025. Monitoring well TM04-PZM006 was evaluated with a 95% confidence factor and identified statistically significant decreasing trends in the benzene and naphthalene concentrations. The analysis output file is included in **Appendix B**.

### 2.1.3. B24 Groundwater Sampling

On April 8<sup>th</sup> and 15<sup>th</sup>, 2025, ARM sampled all three shallow monitoring wells on Parcel B24: B24-001-MWS, B24-002-MWS and TS03-DDP002. These three monitoring wells are located to the west of the impoundment and were included in the monitoring program as sentinel wells to identify potential migration towards Bear Creek.

The three monitoring wells were sampled for VOCs via USEPA Method 8260 and SVOCs by USEPA Method 8270 SIM. All results from the 1<sup>st</sup> Half 2025 sampling event are included in **Table 3**. The laboratory reports are included in **Appendix A**. All groundwater PAL exceedances are included in **Figure 3**.

Benzene did not exceed the groundwater PAL in the three Parcel B24 sentinel monitoring wells. B24-002-MWS had three VOC exceedances: a concentration of 4.5 µg/L of bromodichloromethane (PAL of 0.13 µg/L), 7.9 µg/L of bromoform (PAL of 3.3 µg/L), and 2.9 µg/L of chloroform (PAL of 0.22 µg/L). For SVOCs, naphthalene exceeded the groundwater PAL (0.12 µg/L) in B24-001-MWS, with an observed concentration of 3.6 µg/L. Groundwater

concentrations at the Parcel B24 sentinel monitoring wells have remained low since sampling began in May 2022 (following well installation in April 2022).

## 2.2. SURFACE WATER SAMPLING

Surface water samples were collected in April 2025 from four locations: TMC-Outlet (End of Canal), TMC-TM04 (adjacent to TM04-PZM006), TMC-Bend (from the bend in the TMC in the southeastern corner of B14, adjacent to HI15-MWS), and TMC-Rail Bridge (Mid Canal) (refer to **Figure 2**). All surface water samples were analyzed for VOCs via USEPA Method 8260 and SVOCs by USEPA Method 8270 SIM.

For comparison purposes, the USEPA Region 4 Surface Water Screening Values (for freshwater and chronic) (Region 4 Ecological Risk Assessment Supplemental Guidance, March 2018) were included in **Table 4** for parameters with detections. The laboratory reports are included in **Appendix A**.

For the 1<sup>st</sup> Half of 2025, benzene and naphthalene did not exceed the freshwater screening levels (160 µg/L and 21 µg/L, respectively). Benzene concentrations ranged from 0.17 J µg/L to 0.42 µg/L for benzene, while all naphthalene results were non-detect. Benzene and naphthalene historic surface water sampling results are compiled in **Table 5**.

Benzo[g,h,i]perylene did exceed its freshwater screening level at TMC-Outlet, with a concentration of 0.03 J µg/L versus the screening level of 0.012 µg/L. There were no other exceedances of surface water screening levels in the 1<sup>st</sup> Half of 2025.

## 2.3. GROUNDWATER ELEVATION

To assess potential impacts from the IM on groundwater flow direction, TPA is conducting groundwater gauging throughout the duration of the development/remedial activities. On October 30, 2023, the Agencies approved a reduction in gauging frequency from monthly to quarterly. Gauging results are summarized in **Table 6**. Gauging results from the 1<sup>st</sup> Half of 2025 have been used to create groundwater elevation contour maps, included as **Figure 6**. The following wells are not included in the gauging: HI17-MWS (dry) and HI20R2-MWS (not surveyed).

During groundwater monitoring events, each groundwater point was checked for the presence of NAPL using an oil-water interface probe. Apart from monitoring well HI17-MWS and HI20R-MWS, discussed below, NAPL was not identified in any of the monitoring wells.

## 2.4. NAPL OBSERVATIONS

During groundwater gauging on April 18, 2022, approximately 0.02 ft of NAPL was identified in HI17-MWS. MDE was notified of the NAPL detection on April 19, 2022. HI17-MWS was gauged at least weekly following this detection, with NAPL thickness ranging from the initial detection of 0.02 ft to not measurable. Three enhanced fluid recovery events were conducted at HI17-MWS

between April and May 2022; minimal product was recovered. In July 2022, a 10-foot-long trench was advanced to groundwater within the impoundment, approximately 20 feet west of HI17-MWS. No NAPL was observed within the trench for several weeks prior to backfilling.

Since July 2022, absorbent socks have been utilized in HI17-MWS, with an absorbent sock placed into the well for one month and then removed for one month to evaluate the amount of potential NAPL recharge into the well. A summary of NAPL observations during the 1<sup>st</sup> Half of 2025 is provided in **Table 7**. NAPL thickness has not increased at this location, with only trace amounts observed. Absorbent socks had either trace product or were ~1/3 saturated with product at removal. Absorbent socks will continue to be used at this location to remove NAPL from the well.

Monitoring well HI20-MWS was abandoned in March 2023 due to anticipated development activities on Sub-Parcel B8-1. Monitoring well HI20R-MWS was installed in January 2024, approximately 53 feet south of the previous well location due to construction of the Parcel B8-1 building pad; the HI20R-MWS location is within Parcel B14/Humphrey Impoundment (refer to **Figure 2**). During well installation, a black tar-like substance was observed at several depths in the boring. Monitoring well HI20R-MWS was developed on January 25, 2024, and no NAPL was identified. On February 20, 2024, the monitoring well was gauged before the Q1 2024 sampling event; 0.73 ft of NAPL was observed in the well. An absorbent sock was deployed in the well. However, subsequent gauging at HI20R-MWS have identified less than 0.02 ft of NAPL; the initial NAPL observation appears to have been an erroneous measurement due to the viscous nature of the NAPL. A summary of NAPL observations during the 1<sup>st</sup> Half of 2025 is provided in **Table 7**; 0.02 ft of NAPL was observed in January 2025, with all other months having trace amounts of NAPL or being dry. Absorbent socks had either trace product or were ~1/4 saturated with product at removal. This location will continue to be gauged, and absorbent socks deployed as needed.

Monitoring well HI20R2-MWS was installed in March 2024 directly next to the abandoned HI20-MWS well. Due to its proximity to HI20R-MWS, an absorbent sock was deployed in the well. No NAPL has been detected during monthly monitoring since its installation.

### 3.0 SUMMARY AND CONCLUSIONS

The IM Work Plan for the Humphrey Impoundment (Revision 2 dated March 29, 2022) was proposed as a component of the final remedy for the waste material contained in the Parcel B14 impoundment. The IM Work Plan proposed groundwater and surface water gauging and sampling to monitor the site conditions as the Humphrey Impoundment is filled, in preparation for development.

Filling of the impoundment is approximately 95% complete, with paving approximately 50% complete. Since commencement of the filling of HI, there have been no significant changes in groundwater concentrations, surface water concentrations, or groundwater elevations, and no seasonal trends have been identified. TMC surface-water elevation continues to be largely controlled at the HCWWTP intake. Groundwater gauging, groundwater sampling, and surface water sampling will continue to be conducted on a semi-annual basis.

Monitoring wells HI17-MWS and HI20R-MWS, where NAPL/sheen has been observed, will continue to be gauged monthly, with absorbent socks deployed as needed.

The next IM Report will be submitted following completion of groundwater and surface water activities for the second half of 2025.

---

---

## FIGURES

---

---



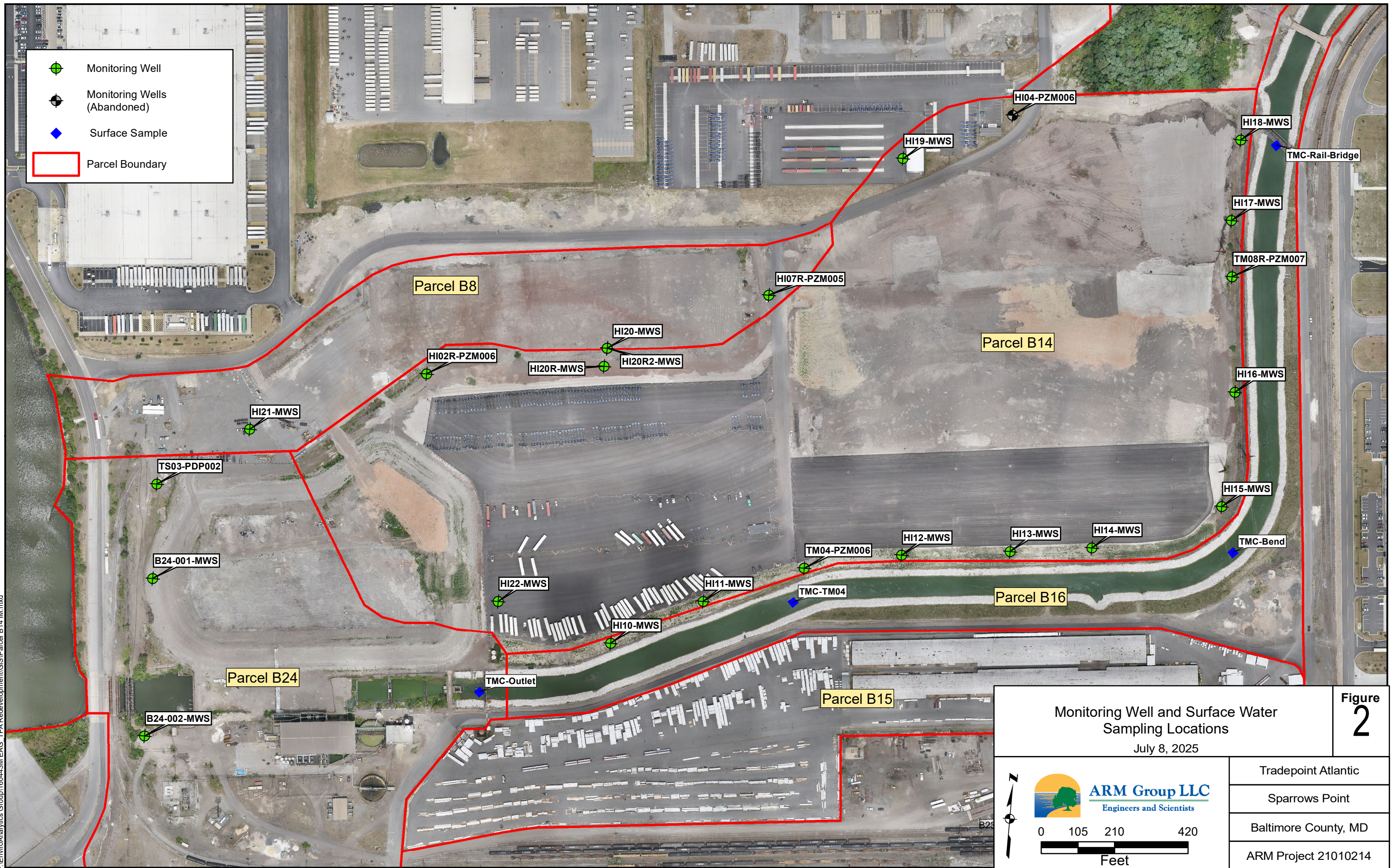


P:\EnviroAnalytics Group\150298M EAG Sparrows Point Area A\GIS\Site-wide Figures.mxd

|                                                                               |  |                          |  |
|-------------------------------------------------------------------------------|--|--------------------------|--|
| Tradepoint Atlantic Property<br>Parcel B14 Site Location<br>February 20, 2023 |  | Figure<br>1              |  |
| <br><br>                                                                      |  | Tradepoint Atlantic      |  |
|                                                                               |  | Sparrows Point           |  |
|                                                                               |  | Baltimore County, MD     |  |
|                                                                               |  | ARM Project No. 21010214 |  |



P:\EnviroAnalytics Group\160443M EAG TPA Redevelopment\GIS\Parcel B14 IM.mxd



Monitoring Well and Surface Water  
Sampling Locations

July 8, 2025



**ARM Group LLC**  
Engineers and Scientists

Tradepoint Atlantic

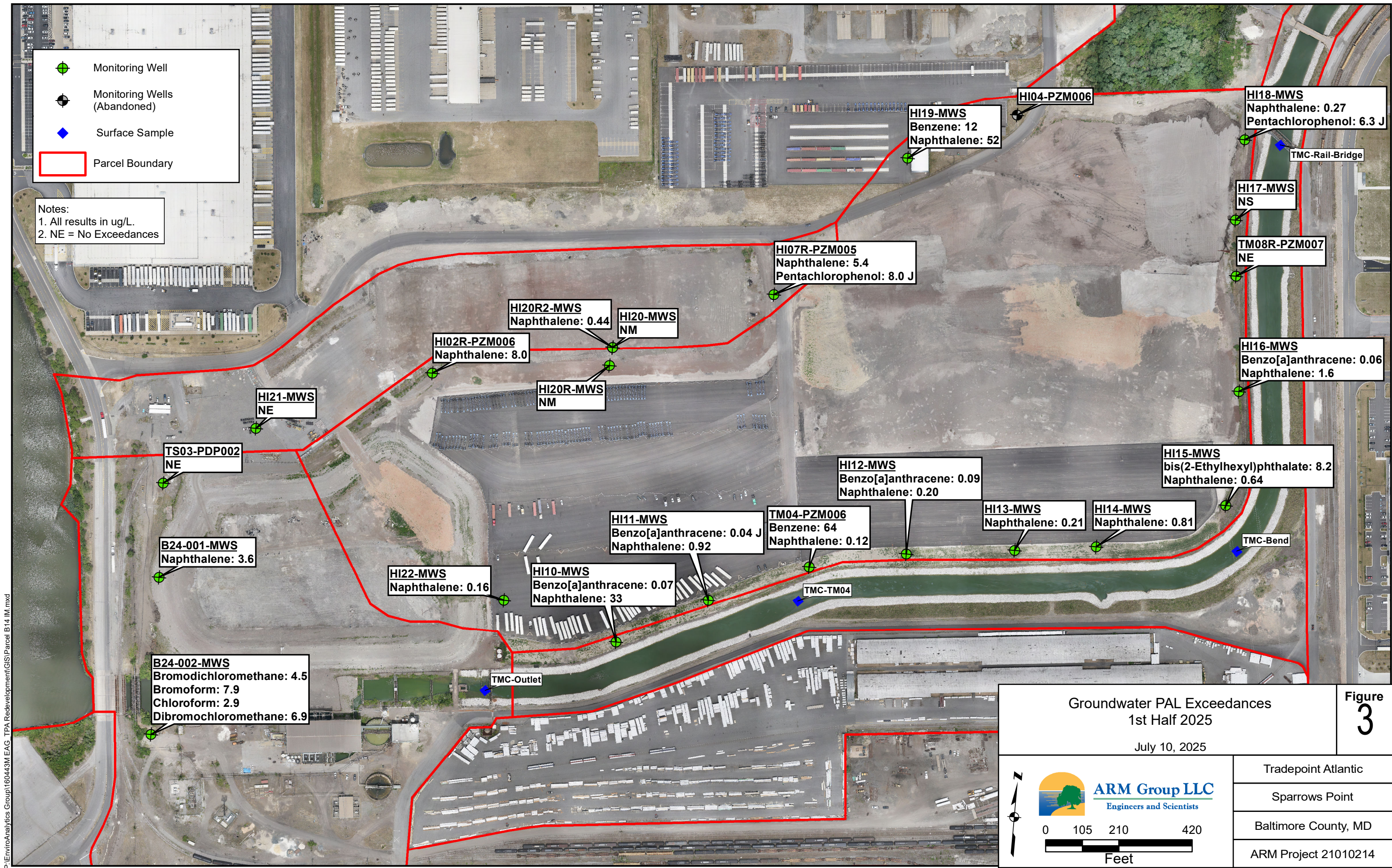
Sparrows Point

Baltimore County, MD

ARM Project 21010214

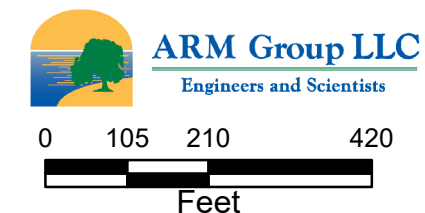
Figure  
**2**





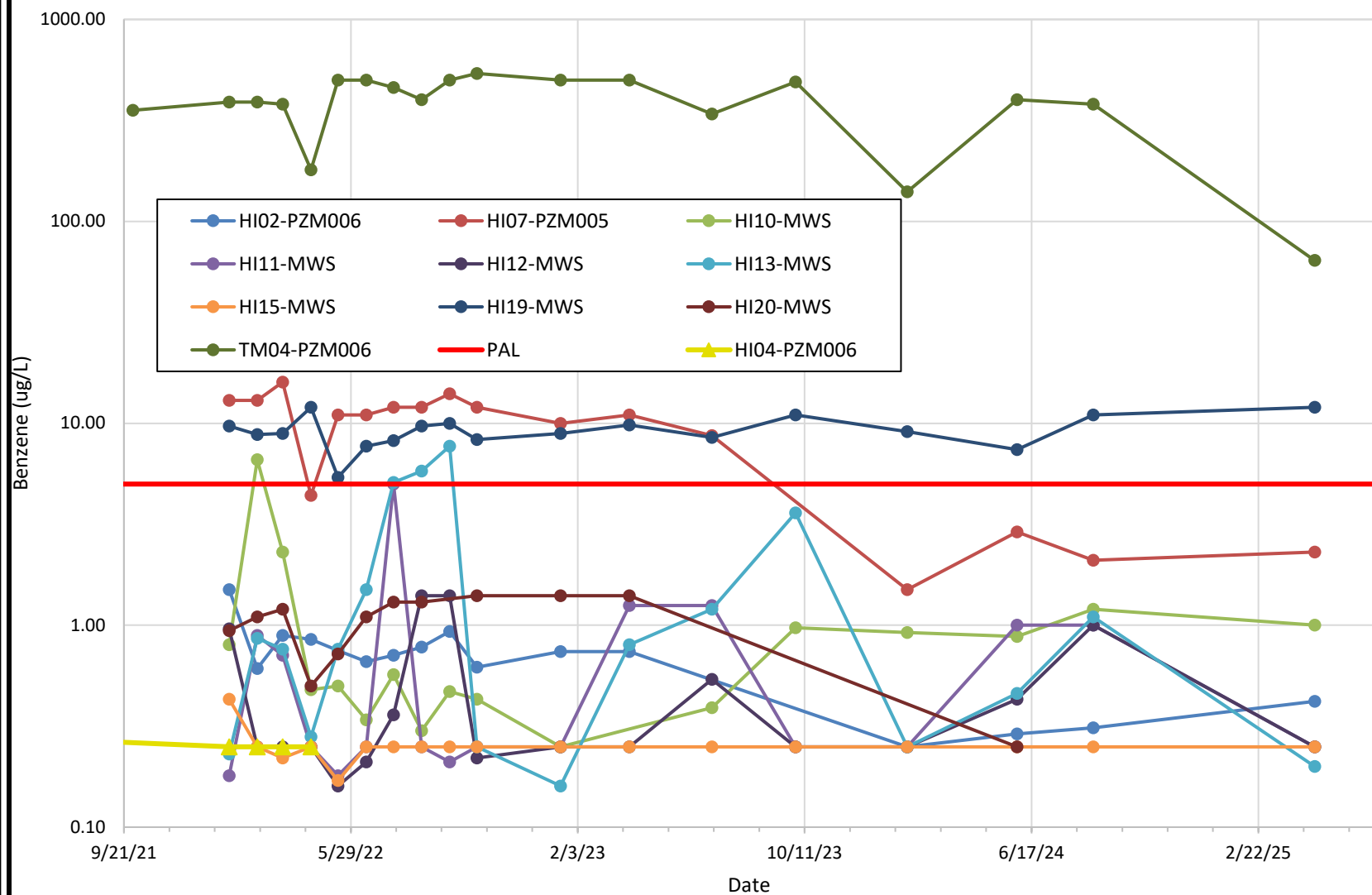
Groundwater PAL Exceedances  
1st Half 2025  
July 10, 2025

Figure  
3



|                      |
|----------------------|
| Tradepoint Atlantic  |
| Sparrows Point       |
| Baltimore County, MD |
| ARM Project 21010214 |





**ARM Group LLC**  
Engineers and Scientists

Parcel B14  
Tradeport Atlantic

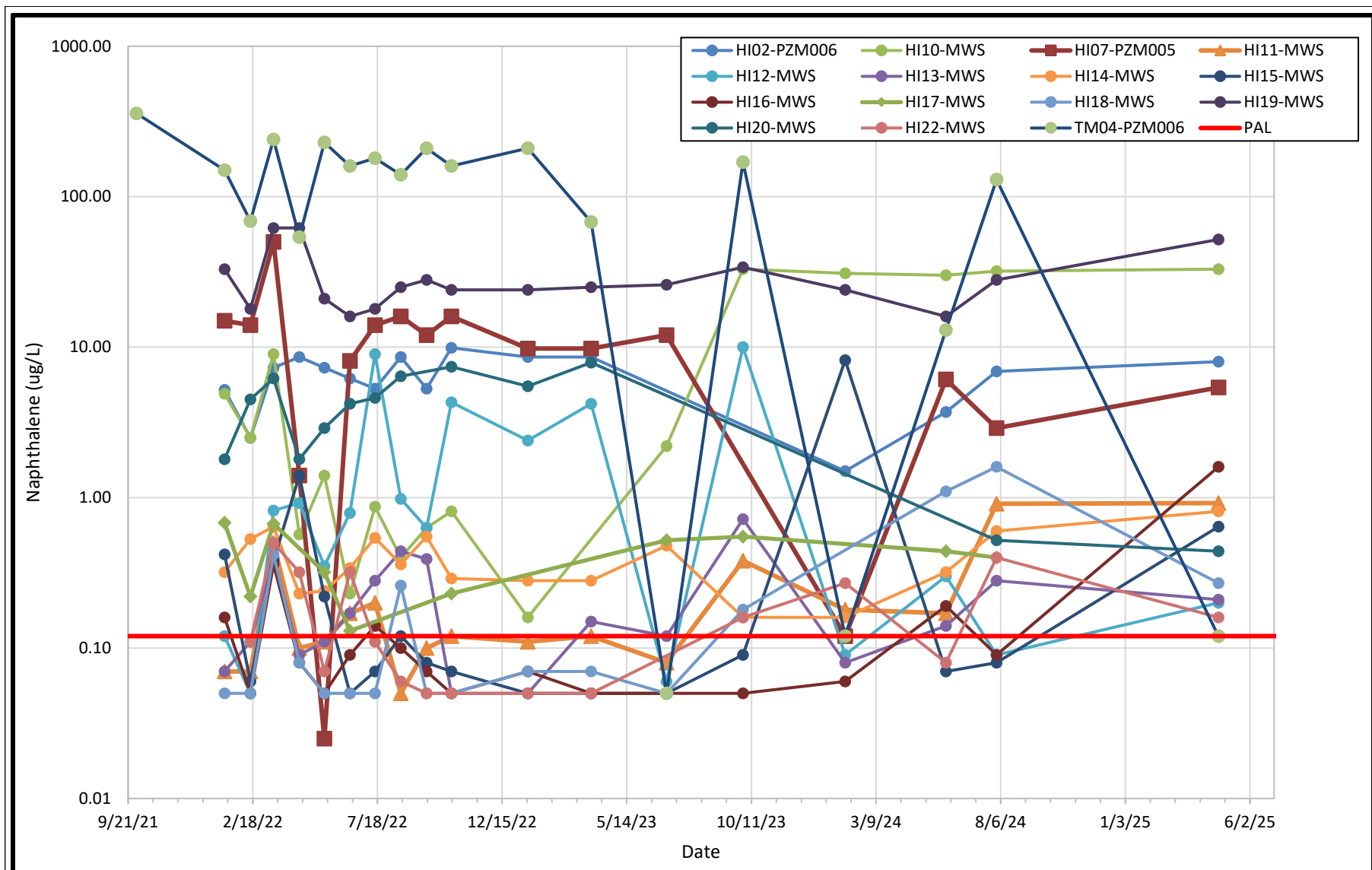
Sparrows Point,  
Maryland

## B14 Historic Groundwater Results for Benzene

July 2, 2025

**Figure  
4**





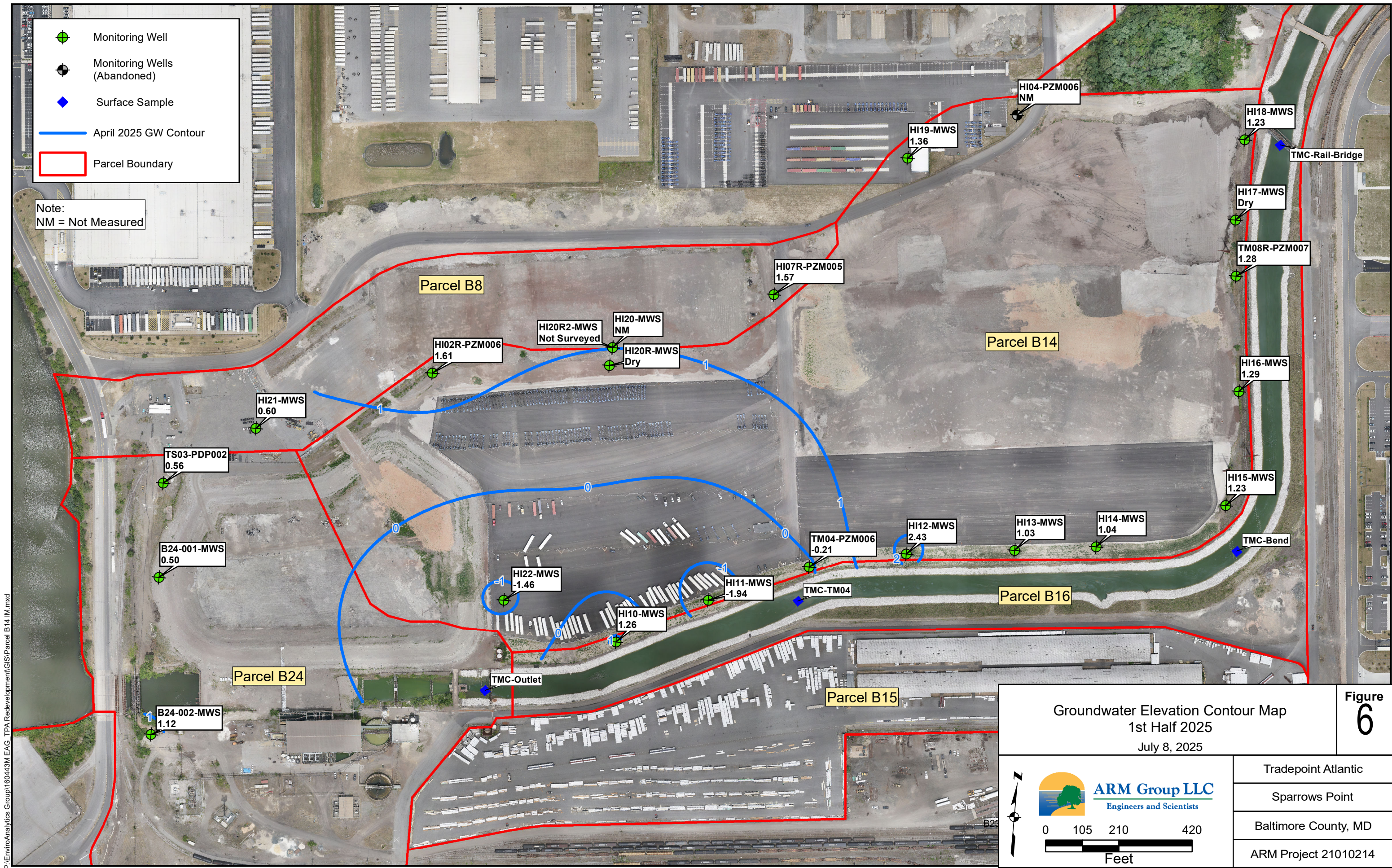
**ARM Group LLC**  
Engineers and Scientists

Parcel B14  
Tradeport Atlantic  
Sparrows Point,  
Maryland

**B14 Historic Groundwater  
Results for Naphthalene**  
July 2, 2025

**Figure  
5**





Groundwater Elevation Contour Map  
1st Half 2025  
July 8, 2025

Figure  
6

**ARM Group LLC**  
Engineers and Scientists

0 105 210 420  
Feet

|                      |
|----------------------|
| Tradepoint Atlantic  |
| Sparrows Point       |
| Baltimore County, MD |
| ARM Project 21010214 |



---

---

## TABLES

---

---



Table 1 - Parcel B14 Groundwater Sampling (1st Half 2025)  
Summary of Organics Detected in Groundwater

| Parameter                    |      | Units | PAL   | HI02R-PZM006<br>4/8/2025 | HI07R-PZM005<br>4/9/2025 | HI10-MWS<br>4/14/2025 | HI11-MWS<br>4/14/2025 | HI12-MWS<br>4/14/2025 | HI13-MWS<br>4/15/2025 | HI14-MWS<br>4/10/2025 | HI15-MWS<br>4/10/2025 |
|------------------------------|------|-------|-------|--------------------------|--------------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| SVOCs                        |      |       |       |                          |                          |                       |                       |                       |                       |                       |                       |
| 1,1-Biphenyl                 | SVOC | µg/L  | 0.83  | 0.24 J                   | 0.26 J                   | 0.32 J                | 2 U                   | 2 U                   | 2 U                   | 2 U                   | 2 U                   |
| 2,4-Dimethylphenol           | SVOC | µg/L  | 360   | 5.5                      | 5 U                      | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   |
| 2-Methylnaphthalene          | SVOC | µg/L  | 36    | 0.90                     | 0.85                     | 1.9                   | 0.04 J                | 0.1 U                 | 0.1 U                 | 0.05 J                | 0.04 J                |
| 2-Methylphenol               | SVOC | µg/L  | 930   | 5 U                      | 5 U                      | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   |
| 3&4-Methylphenol(m&p Cresol) | SVOC | µg/L  | 930   | 5 U                      | 3.6 J                    | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   |
| Acenaphthene                 | SVOC | µg/L  | 530   | 1.4                      | 0.55                     | 2.2                   | 0.07 J                | 0.1 U                 | 0.1 U                 | 1.6                   | 0.1 U                 |
| Acenaphthylene               | SVOC | µg/L  | 530   | 0.05 J                   | 0.36                     | 0.57                  | 0.05 J                | 0.04 J                | 0.1 U                 | 0.05 J                | 0.1 U                 |
| Anthracene                   | SVOC | µg/L  | 1800  | 0.97                     | 0.28                     | 0.40                  | 0.12                  | 0.09 J                | 0.04 J                | 0.19                  | 0.06 J                |
| Benzo[a]anthracene           | SVOC | µg/L  | 0.03  | 0.05 U                   | 0.05 U                   | 0.07                  | 0.04 J                | 0.09                  | 0.05 U                | 0.05 U                | 0.05 U                |
| Benzo[a]pyrene               | SVOC | µg/L  | 0.20  | 0.1 U                    | 0.1 U                    | 0.1 U                 | 0.1 U                 | 0.03 J                | 0.1 U                 | 0.1 U                 | 0.1 U                 |
| Benzo[b]fluoranthene         | SVOC | µg/L  | 0.25  | 0.05 U                   | 0.05 U                   | 0.03 J                | 0.05 U                | 0.07                  | 0.05 U                | 0.05 U                | 0.05 U                |
| Benzo[g,h,i]perylene         | SVOC | µg/L  | -     | 0.1 U                    | 0.1 U                    | 0.1 U                 | 0.1 U                 | 0.04 J                | 0.1 U                 | 0.1 U                 | 0.1 U                 |
| bis(2-Ethylhexyl)phthalate   | SVOC | µg/L  | 6.0   | 3 U                      | 3 U                      | 3 U                   | 3 U                   | 3 U                   | 3 U                   | 3 U                   | 8.2                   |
| Carbazole                    | SVOC | µg/L  | -     | 0.42 J                   | 0.92 J                   | 4.2                   | 2 U                   | 2 U                   | 2 U                   | 2 U                   | 2 U                   |
| Chrysene                     | SVOC | µg/L  | 25    | 0.1 U                    | 0.05 J                   | 0.04 J                | 0.1 U                 | 0.04 J                | 0.1 U                 | 0.1 U                 | 0.1 U                 |
| Fluoranthene                 | SVOC | µg/L  | 800   | 0.15                     | 0.89                     | 0.72                  | 0.07 J                | 0.10 J                | 0.1 U                 | 0.18                  | 0.1 U                 |
| Fluorene                     | SVOC | µg/L  | 290   | 1.0                      | 0.76                     | 1.8                   | 0.04 J                | 0.1 U                 | 0.1 U                 | 0.82                  | 0.03 J                |
| Indeno[1,2,3-c,d]pyrene      | SVOC | µg/L  | 0.25  | 0.1 U                    | 0.1 U                    | 0.1 U                 | 0.1 U                 | 0.04 J                | 0.1 U                 | 0.1 U                 | 1.1 J                 |
| Naphthalene                  | SVOC | µg/L  | 0.12  | 8.0                      | 5.4                      | 33                    | 0.92                  | 0.20                  | 0.21                  | 0.81                  | 0.64                  |
| Pentachlorophenol            | SVOC | µg/L  | 1.0   | 10 U                     | 8.0 J                    | 10 U                  | 10 U                  | 10 U                  | 10 U                  | 10 U                  | 10 U                  |
| Phenanthrene                 | SVOC | µg/L  | -     | 0.96                     | 1.5                      | 2.8                   | 0.06                  | 0.07                  | 0.05 U                | 0.85                  | 0.05 U                |
| Phenol                       | SVOC | µg/L  | 5800  | 5 U                      | 5 U                      | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   | 5 U                   |
| Pyrene                       | SVOC | µg/L  | 120   | 0.10                     | 0.56                     | 0.53                  | 0.06 J                | 0.08 J                | 0.1 U                 | 0.11                  | 0.1 U                 |
| VOCs                         |      |       |       |                          |                          |                       |                       |                       |                       |                       |                       |
| 1,1-Dichloroethane           | VOC  | µg/L  | 2.7   | 0.75 U                   | 0.26 J                   | 0.75 U                | 0.75 U                | 0.75 U                | 0.75 U                | 0.75 U                | 0.75 U                |
| Acetone                      | VOC  | µg/L  | 14000 | 3.1 J                    | 4.6 J                    | 2.9 J                 | 5.2                   | 4.5 J                 | 3.7 J                 | 2.2 J                 | 3.2 J                 |
| Benzene                      | VOC  | µg/L  | 5.0   | 0.42 J                   | 2.3                      | 1.0                   | 0.5 U                 | 0.5 U                 | 0.20 J                | 0.5 U                 | 0.5 U                 |
| Ethylbenzene                 | VOC  | µg/L  | 700   | 0.5 U                    | 0.5 U                    | 0.5 U                 | 0.5 U                 | 0.5 U                 | 0.5 U                 | 0.5 U                 | 0.5 U                 |
| m&p-Xylene                   | VOC  | µg/L  | -     | 1 U                      | 0.88 J                   | 0.36 J                | 1 U                   | 1 U                   | 1 U                   | 1 U                   | 1 U                   |
| o-Xylene                     | VOC  | µg/L  | -     | 1 U                      | 1 U                      | 1 U                   | 1 U                   | 1 U                   | 1 U                   | 1 U                   | 1 U                   |
| Toluene                      | VOC  | µg/L  | 1000  | 0.75 U                   | 1.5                      | 0.46 J                | 0.75 U                | 0.75 U                | 0.75 U                | 0.75 U                | 0.75 U                |
| Vinyl chloride               | VOC  | µg/L  | 2.0   | 1 U                      | 1 U                      | 1 U                   | 1 U                   | 1 U                   | 1 U                   | 1 U                   | 1 U                   |
| Xylenes                      | VOC  | µg/L  | 10000 | 1 U                      | 0.88 J                   | 0.36 J                | 1 U                   | 1 U                   | 1 U                   | 1 U                   | 1 U                   |

Detections in bold

Exceedances of EPA Region 4 Surface Water Screening Levels (freshwater, chronic) in red

U: This analyte was not detected in the sample. The numeric value represents the sample quantitation/detection limit.

J: The positive result reported for this analyte is a quantitative estimate.

JB: The positive result reported for this analyte is a quantitative estimate and was detected above the reporting limits in the associated method blank.

B: The analyte was detected above the reporting limits in the associated method blank.

µg/L: micrograms per liter

PAL: Project Action Limit



Table 1 - Parcel B14 Groundwater Sampling (1st Half 2025)  
Summary of Organics Detected in Groundwater

| Parameter                    |      | Units | PAL   | HI16-MWS<br>4/10/2025 | HI18-MWS<br>4/9/2025 | HI19-MWS<br>4/10/2025 | HI20R2-MWS<br>4/9/2025 | HI21-MWS<br>4/10/2025 | HI22-MWS<br>4/14/2025 | TM04-PZM006<br>4/15/2025 | TM08R-PZM007<br>4/9/2025 |
|------------------------------|------|-------|-------|-----------------------|----------------------|-----------------------|------------------------|-----------------------|-----------------------|--------------------------|--------------------------|
| SVOCs                        |      |       |       |                       |                      |                       |                        |                       |                       |                          |                          |
| 1,1-Biphenyl                 | SVOC | µg/L  | 0.83  | 2 U                   | 2 U                  | 0.76 J                | 2 U                    | 2 U                   | 2 U                   | 2 U                      | 2 U                      |
| 2,4-Dimethylphenol           | SVOC | µg/L  | 360   | 5 U                   | 5 U                  | 150                   | 8.1                    | 5 U                   | 5 U                   | 5 U                      | 5 U                      |
| 2-Methylnaphthalene          | SVOC | µg/L  | 36    | 0.08 J                | 0.05 J               | 3.7                   | 0.12                   | 0.1 U                 | 0.03 J                | 0.1 U                    | 0.1 U                    |
| 2-Methylphenol               | SVOC | µg/L  | 930   | 5 U                   | 5 U                  | 3.2 J                 | 5 U                    | 5 U                   | 5 U                   | 5 U                      | 5 U                      |
| 3&4-Methylphenol(m&p Cresol) | SVOC | µg/L  | 930   | 5 U                   | 3.3 J                | 57                    | 4.8 J                  | 5 U                   | 5 U                   | 5 U                      | 5 U                      |
| Acenaphthene                 | SVOC | µg/L  | 530   | 0.06 J                | 0.07 J               | 0.90                  | 0.17                   | 0.1 U                 | 0.1 U                 | 0.1 U                    | 0.1 U                    |
| Acenaphthylene               | SVOC | µg/L  | 530   | 0.08 J                | 0.06 J               | 0.56                  | 0.04 J                 | 0.1 U                 | 0.04 J                | 0.07 J                   | 0.1 U                    |
| Anthracene                   | SVOC | µg/L  | 1800  | 0.35                  | 0.26                 | 0.30                  | 0.05 J                 | 0.1 U                 | 0.08 J                | 0.08 J                   | 0.07 J                   |
| Benzo[a]anthracene           | SVOC | µg/L  | 0.03  | 0.06                  | 0.05 U               | 0.1 U                 | 0.05 U                 | 0.05 U                | 0.05 U                | 0.05 U                   | 0.05 U                   |
| Benzo[a]pyrene               | SVOC | µg/L  | 0.20  | 0.04 J                | 0.1 U                | 0.2 U                 | 0.1 U                  | 0.1 U                 | 0.1 U                 | 0.1 U                    | 0.1 U                    |
| Benzo[b]fluoranthene         | SVOC | µg/L  | 0.25  | 0.07                  | 0.05 U               | 0.1 U                 | 0.05 U                 | 0.05 U                | 0.07                  | 0.05 U                   | 0.05 U                   |
| Benzo[g,h,i]perylene         | SVOC | µg/L  | -     | 0.03 J                | 0.1 U                | 0.2 U                 | 0.1 U                  | 0.1 U                 | 0.1 U                 | 0.1 U                    | 0.1 U                    |
| bis(2-Ethylhexyl)phthalate   | SVOC | µg/L  | 6.0   | 3 U                   | 3 U                  | 3 U                   | 3 U                    | 3 U                   | 3 U                   | 3 U                      | 3 U                      |
| Carbazole                    | SVOC | µg/L  | -     | 2 U                   | 0.81 J               | 1.7 J                 | 2 U                    | 2 U                   | 2 U                   | 2 U                      | 2 U                      |
| Chrysene                     | SVOC | µg/L  | 25    | 0.04 J                | 0.1 U                | 0.2 U                 | 0.1 U                  | 0.1 U                 | 0.1 U                 | 0.1 U                    | 0.1 U                    |
| Fluoranthene                 | SVOC | µg/L  | 800   | 0.24                  | 0.05 J               | 0.37                  | 0.13                   | 0.1 U                 | 0.13                  | 0.05 J                   | 0.03 J                   |
| Fluorene                     | SVOC | µg/L  | 290   | 0.1 U                 | 0.04 J               | 1.0                   | 0.19                   | 0.1 U                 | 0.1 U                 | 0.1 U                    | 0.03 J                   |
| Indeno[1,2,3-c,d]pyrene      | SVOC | µg/L  | 0.25  | 0.04 J                | 0.1 U                | 0.2 U                 | 0.1 U                  | 0.1 U                 | 0.05 J                | 0.1 U                    | 0.1 U                    |
| Naphthalene                  | SVOC | µg/L  | 0.12  | 1.6                   | 0.27                 | 52                    | 0.44                   | 0.05 J                | 0.16                  | 0.12                     | 0.08 J                   |
| Pentachlorophenol            | SVOC | µg/L  | 1.0   | 10 U                  | 6.3 J                | 10 U                  | 10 U                   | 10 U                  | 10 U                  | 10 U                     | 10 U                     |
| Phenanthrene                 | SVOC | µg/L  | -     | 0.05                  | 0.05 U               | 1.6                   | 0.29                   | 0.05 U                | 0.07                  | 0.05 U                   | 0.05 U                   |
| Phenol                       | SVOC | µg/L  | 5800  | 5 U                   | 5 U                  | 4.2 J                 | 5 U                    | 5 U                   | 5 U                   | 5 U                      | 5 U                      |
| Pyrene                       | SVOC | µg/L  | 120   | 0.23                  | 0.1 U                | 0.24                  | 0.13                   | 0.1 U                 | 0.1 U                 | 0.06 J                   | 0.1 U                    |
| VOCs                         |      |       |       |                       |                      |                       |                        |                       |                       |                          |                          |
| 1,1-Dichloroethane           | VOC  | µg/L  | 2.7   | 0.75 U                | 0.75 U               | 0.75 U                | 0.75 U                 | 0.75 U                | 0.75 U                | 0.22 J                   | 0.75 U                   |
| Acetone                      | VOC  | µg/L  | 14000 | 7.4                   | 5.3                  | 2.7 J                 | 2.8 J                  | 5 U                   | 5.1                   | 5.8                      | 3.3 J                    |
| Benzene                      | VOC  | µg/L  | 5.0   | 0.5 U                 | 0.5 U                | 12                    | 0.32 J                 | 0.5 U                 | 0.23 J                | 64                       | 0.5 U                    |
| Ethylbenzene                 | VOC  | µg/L  | 700   | 0.5 U                 | 0.5 U                | 0.19 J                | 0.5 U                  | 0.5 U                 | 0.5 U                 | 0.71                     | 0.5 U                    |
| m&p-Xylene                   | VOC  | µg/L  | -     | 1 U                   | 1 U                  | 2.0                   | 1 U                    | 1 U                   | 1 U                   | 0.65 J                   | 1 U                      |
| o-Xylene                     | VOC  | µg/L  | -     | 1 U                   | 1 U                  | 1.0                   | 1 U                    | 1 U                   | 1 U                   | 1 U                      | 1 U                      |
| Toluene                      | VOC  | µg/L  | 1000  | 0.75 U                | 0.75 U               | 4.1                   | 0.75 U                 | 0.75 U                | 0.75 U                | 0.36 J                   | 0.75 U                   |
| Vinyl chloride               | VOC  | µg/L  | 2.0   | 1 U                   | 1 U                  | 1 U                   | 1 U                    | 1 U                   | 1 U                   | 0.15 J                   | 1 U                      |
| Xylenes                      | VOC  | µg/L  | 10000 | 1 U                   | 1 U                  | 3.0                   | 1 U                    | 1 U                   | 1 U                   | 0.65 J                   | 1 U                      |

Detections in bold

Exceedances of EPA Region 4 Surface Water Screening Levels (freshwater, chronic) in red

U: This analyte was not detected in the sample. The numeric value represents the sample quantitation/detection limit.

J: The positive result reported for this analyte is a quantitative estimate.

JB: The positive result reported for this analyte is a quantitative estimate and was detected above the reporting limits in the associated method blank.

B: The analyte was detected above the reporting limits in the associated method blank.

µg/L: micrograms per liter

PAL: Project Action Limit



Table 2 - Parcel B14 Groundwater Sampling  
Historic Groundwater Sampling Results

| Well ID                    | PAL | Benzene (µg/L) |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |
|----------------------------|-----|----------------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|
|                            |     | Dec-01         | Jul-04 | Oct-17 | Oct-21 | Jan-22 | Feb-22 | Mar-22 | Apr-22 | May-22 | Jun-22 | Jul-22 | Aug-22 | Sep-22 | Oct-22 | Jan-23 | Apr-23 | Jul-23 | Oct-23 | Feb-24 | Jun-24 | Aug-24 | Apr-25 |
| HI02-PZM006 / HI02R-PZM006 | 5   | 1.2            |        | 0.88 J | NS     | 1.5    | 0.61   | 0.89   | 0.85   | 0.75   | 0.66   | 0.71   | 0.78   | 0.93   | 0.62   | 0.74   | 0.74   |        |        | 0.5 U  | 0.29 J | 0.31 J | 0.42 J |
| HI04-PZM006                | 5   | 1 U            |        | 1 U    | NS     | 0.5 U  | 0.5 U  | 0.50 U | 0.5 U  |        |        |        |        |        |        |        |        |        |        |        |        |        |        |
| HI07-PZM005 / HI07R-PZM005 | 5   | 25             | 16     | 16.2   | NS     | 13.0   | 13.0   | 16.0   | 4.4    | 11     | 11     | 12.0   | 12.0   | 14     | 12     | 10     | 11     | 8.7    |        | 1.5    | 2.9    | 1.2    | 2.3    |
| HI10-MWS                   | 5   |                |        |        |        | 0.8    | 6.6    | 2.3    | 0.48 J | 1.0 U  | 0.34 J | 0.57   | 0.30 J | 0.47 J | 0.43 J | 0.50 U | 0.69   | 0.39 J | 0.97   | 0.92   | 0.88   | 1.2    | 1.0    |
| HI11-MWS                   | 5   |                |        |        |        | 0.18 J | 0.89   | 0.71   | 0.5 U  | 0.18 J | 0.50 U | 10 U   | 0.50 U | 0.21 J | 0.50 U | 0.50 U | 2.5 U  | 2.5 U  | 0.5 U  | 0.5 U  | 2 U    | 1.0    | 0.5 U  |
| HI12-MWS                   | 5   |                |        |        |        | 0.96   | 0.5 U  | 0.50 U | 0.5 U  | 0.16 J | 0.21 J | 0.36 J | 1.4    | 1.4    | 0.22 J | 0.50 U | 0.5 U  | 0.54   | 0.5 U  | 0.5 U  | 0.43 J | 1.0    | 0.5 U  |
| HI13-MWS                   | 5   |                |        |        |        | 0.23 J | 0.86   | 0.76   | 0.28 J | 0.76   | 1.5    | 5.1    | 5.8    | 7.7    | 0.50 U | 0.16 J | 0.8    | 1.2    | 3.6    | 0.5 U  | 0.46 J | 1.1    | 0.2 J  |
| HI14-MWS                   | 5   |                |        |        |        | 0.5 U  | 0.5 U  | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  |
| HI15-MWS                   | 5   |                |        |        |        | 0.43 J | 0.5 U  | 0.22 J | 0.5 U  | 0.17 J | 0.5 U  | 0.5 U  | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  |
| HI16-MWS                   | 5   |                |        |        |        | 0.5 U  | 0.5 U  | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  |
| HI17-MWS                   | 5   |                |        |        |        | 0.5 U  | 0.5 U  | 0.50 U | NS (b) | 0.5 U  | 0.5 U  | NS (b) | NS (b) | NS (b) | 0.50 U | NS (b) | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | NS*    |
| HI18-MWS                   | 5   |                |        |        |        | 0.5 U  | 0.5 U  | 0.50 U | 0.5 U  | 0.5 U  | 2.5 U  | 0.5 U  | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.23 J | 0.16 J | 0.16 J | 0.5 U  |
| HI19-MWS                   | 5   |                |        |        |        | 9.7    | 8.8    | 8.9    | 12     | 5.4    | 7.7    | 8.2    | 9.7    | 10     | 8.3    | 8.9    | 9.8    | 8.5    | 11     | 9.1    | 7.4    | 11     | 12     |
| HI20-MWS / HI20R2-MWS      | 5   |                |        |        |        | 0.94   | 1.1    | 1.2    | 0.50   | 0.72   | 1.10   | 1.3    | 1.3    | NS (c) | 1.4    | 1.4    | 1.4    |        |        |        |        | 0.35 J | 0.32 J |
| HI20R-MWS                  | 5   |                |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        |        | 0.5 U  | NS     | NS     |
| HI21-MWS                   | 5   |                |        |        |        |        |        |        |        | 0.50 U | 0.5 U  | 0.5 U  | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  |
| HI22-MWS                   | 5   |                |        |        |        |        | 2.5 U  | 0.50 U | 5 U    | 0.50 U | 2.5 U  | 5 U    | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 2.5 U  | NS     | 0.3 J  | 0.44 J | 0.46 J | 0.2 J  | 0.23 J |
| TM04-PZM006                | 5   | 1400           | 610    | 653    | 355    | 390    | 390    | 380    | 180    | 500    | 500    | 460    | 400    | 500    | 540    | 500    | 500    | 340    | 490    | 140    | 400    | 380    | 64     |
| TM08R-PZM007               | 5   | 1 U            |        | 1 U    | NS     | 0.5 U  | 0.5 U  | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 0.5 U  | 0.50 U | 0.50 U | 0.50 U | 0.50 U | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  | 0.5 U  |

Detections in bold  
Values in red indicate an exceedance of the Project Action Limit (PAL)

U: This analyte was not detected in the sample. The numeric value represents the sample quantitation / detection limit.  
J: The positive result reported for this analyte is a quantitative estimate.  
B: The analyte was detected above the reporting limits in the associated method blank.  
NS: Not Sampled  
µg/L: micrograms per liter  
PAL: Project Action Limit  
(a) Analysis was re-run due to B qualifier.  
(b) HI17-MWS was not sampled in January 2023 due to trace NAPL.  
(c) HI20-MWS was within an active demolition area and could not be safely accessed.  
\* Well was dry.

|  |                        |
|--|------------------------|
|  | Well not installed     |
|  | Well damaged/abandoned |



| Table 2 - Parcel B14 Groundwater Sampling<br>Historic Groundwater Sampling Results |      |                    |        |         |        |        |        |        |        |        |         |        |                     |        |        |        |        |         |        |         |        |        |        |
|------------------------------------------------------------------------------------|------|--------------------|--------|---------|--------|--------|--------|--------|--------|--------|---------|--------|---------------------|--------|--------|--------|--------|---------|--------|---------|--------|--------|--------|
| Well ID                                                                            | PAL  | Naphthalene (µg/L) |        |         |        |        |        |        |        |        |         |        |                     |        |        |        |        |         |        |         |        |        |        |
|                                                                                    |      | Dec-01             | Jul-04 | Oct-17  | Oct-21 | Jan-22 | Feb-22 | Mar-22 | Apr-22 | May-22 | Jun-22  | Jul-22 | Aug-22              | Sep-22 | Oct-22 | Jan-23 | Apr-23 | Jul-23  | Oct-23 | Feb-24  | Jun-24 | Aug-24 | Apr-25 |
| HI02-PZM006 / HI02R-PZM006                                                         | 0.12 | 7.3 J              |        | 5.1     | NS     | 5.2    | 2.5    | 7.3    | 8.6    | 7.3    | 6.2     | 5.3    | 8.6                 | 5.3    | 9.9    | 8.6    | 8.6    |         |        | 1.5     | 3.7    | 6.9    | 8.0    |
| HI04-PZM006                                                                        | 0.12 |                    |        | 0.15    | NS     | 0.1 U  | 0.1 U  | 1.0 U  | 0.06 J |        |         |        |                     |        |        |        |        |         |        |         |        |        |        |
| HI07-PZM005 / HI07R-PZM005                                                         | 0.12 | 40                 | 16     | 26.1    | NS     | 15     | 14.0   | 50     | 1.4    | 0.05 J | 8.1     | 14.0   | 16                  | 12     | 16     | 10     | 9.8    | 12      |        | 0.12    | 6.1    | 2.9    | 5.4    |
| HI10-MWS                                                                           | 0.12 |                    |        |         |        | 4.9    | 2.5    | 9.0    | 0.57   | 1.4    | 0.23    | 0.87   | 0.4                 | 0.63   | 0.81   | 0.16   | 2.8    | 2.2 B   | 33     | 31      | 30     | 32     | 33     |
| HI11-MWS                                                                           | 0.12 |                    |        |         |        | 0.07 J | 0.07 J | 1.0 U  | 0.10   | 0.11   | 0.17    | 0.20   | 0.10 U              | 0.10   | 0.12   | 0.11   | 0.12   | 0.08 JB | 0.38   | 0.18    | 0.17   | 0.91   | 0.92   |
| HI12-MWS                                                                           | 0.12 |                    |        |         |        | 0.12   | 0.1 U  | 0.82 J | 0.92   | 0.35   | 0.79    | 9.0    | 0.98                | 0.63   | 4.3    | 2.4    | 4.2    | 0.06 J  | 10     | 0.09 J  | 0.30   | 0.09 J | 0.20   |
| HI13-MWS                                                                           | 0.12 |                    |        |         |        | 0.07 J | 0.11   | 0.40 J | 0.09 J | 0.11   | 0.17    | 0.28   | 0.44                | 0.39   | 0.1 U  | 0.10 U | 0.15   | 0.12 B  | 0.72 B | 0.08 J  | 0.12   | 0.28   | 0.21   |
| HI14-MWS                                                                           | 0.12 |                    |        |         |        | 0.32   | 0.53   | 0.64 J | 0.23   | 0.24   | 0.34    | 0.54   | 0.36                | 0.55   | 0.29   | 0.28   | 0.28   | 0.48    | 0.16   | 0.16    | 0.32   | 0.6    | 0.81   |
| HI15-MWS                                                                           | 0.12 |                    |        |         |        | 0.42   | 0.06 J | 0.38 J | 1.4    | 0.22   | 0.10 U  | 0.07 J | 0.12                | 0.08 J | 0.07 J | 0.10 U | 0.1 U  | 0.1 U   | 0.09 J | 8.2     | 0.05 J | 0.08 J | 0.64   |
| HI16-MWS                                                                           | 0.12 |                    |        |         |        | 0.16   | 0.1 U  | 0.36 J | 0.08 J | 0.10 U | 0.09 J  | 0.14   | 0.10 J              | 0.07 J | 0.1 U  | 0.07 J | 0.1 U  | 0.1 U   | 0.1 U  | 0.06 JB | 0.19   | 0.09 J | 1.6    |
| HI17-MWS                                                                           | 0.12 |                    |        |         |        | 0.68   | 0.22   | 0.67 J | NS (b) | 0.32   | 0.13    | NS (b) | NS (b)              | NS (b) | 0.23   | NS (b) | 0.27   | 0.52    | 0.55   | 0.44 B  | 0.39   | 0.40   | NS     |
| HI18-MWS                                                                           | 0.12 |                    |        |         |        | 0.1 U  | 0.1 U  | 0.42 J | 0.08 J | 0.10 U | 0.05 JB | 0.1 U  | 0.26                | 0.10 U | 0.05 J | 0.07 J | 0.07 J | 0.1 U   | 0.18   | 0.44 B  | 1.1    | 1.6    | 0.27   |
| HI19-MWS                                                                           | 0.12 |                    |        |         |        | 33     | 18.0   | 62     | 62     | 21     | 16      | 18.0   | 25                  | 28     | 24     | 24     | 25     | 26      | 34     | 24      | 3.3    | 28     | 52     |
| HI20-MWS / HI20R2-MWS                                                              | 0.12 |                    |        |         |        | 1.8    | 4.5    | 6.2    | 1.8    | 2.9    | 4.2     | 4.6    | 6.4                 | NS (c) | 7.4    | 5.5    | 7.9    |         |        |         |        | 0.52   | 0.44   |
| HI20R-MWS                                                                          | 0.12 |                    |        |         |        |        |        |        |        |        |         |        |                     |        |        |        |        |         |        |         | 0.16   | NS     | NS     |
| HI21-MWS                                                                           | 0.12 |                    |        |         |        |        |        |        |        | 0.07 J | 0.06 J  | 0.06 J | 0.26 B / 0.07 J (a) | 0.08 J | 0.12   | 0.10 U | 0.1 U  | 0.1 U   | 0.1 U  | 0.1 U   | 0.1 U  | 0.38   | 0.05 J |
| HI22-MWS                                                                           | 0.12 |                    |        |         |        |        | 0.11   | 1.0 U  | 0.32   | 0.07 J | 0.32    | 0.11   | 0.06 J              | 0.10 U | 0.10 U | 0.10 U | 0.1 U  | NS      | 0.16 J | 0.27    | 0.08 J | 0.40   | 0.16   |
| TM04-PZM006                                                                        | 0.12 | 200                | 51     | 405     | 358    | 150 E  | 69     | 240    | 54     | 230    | 160     | 180    | 140                 | 210    | 160    | 210    | 68     | 0.1 U   | 170    | 0.12    | 13.0   | 130    | 0.12   |
| TM08R-PZM007                                                                       | 0.12 | 10 U               |        | 0.096 J | NS     | 0.1 U  | 0.10   | 1.0 U  | 0.1 U  | 0.10 U | 0.10 U  | 0.1 U  | 0.10 U              | 0.10 U | 0.10 U | 0.10 U | 0.1 U  | 0.1 U   | 0.1 U  | 0.05 JB | 0.1 U  | 0.13   | 0.08 J |

Detections in bold  
Values in red indicate an exceedance of the Project Action Limit (PAL)

U: This analyte was not detected in the sample. The numeric value represents the sample quantitation / detection limit.  
J: The positive result reported for this analyte is a quantitative estimate.  
B: The analyte was detected above the reporting limits in the associated method blank.  
E: Analyte concentration is greater than the reporting limit; however, the reported concentration should be considered estimated.  
NS: Not Sampled  
µg/L: micrograms per liter  
PAL: Project Action Limit  
(a) Analysis was re-run due to B qualifier.  
(b) HI17-MWS was not sampled due to trace NAPL.  
(c) HI20-MWS was within an active demoltion area and could not be safely accessed.

|  |                        |
|--|------------------------|
|  | Well not installed     |
|  | Well damaged/abandoned |

**Table 3 - Parcel B24 Groundwater Sampling (1st Half 2025)**  
**Summary of Organics Detected in Groundwater**

| Parameter               | Units | PAL    | B24-001-MWS<br>4/8/2025 | B24-002-MWS<br>4/15/2025 | TS03-PDP002<br>4/8/2025 |
|-------------------------|-------|--------|-------------------------|--------------------------|-------------------------|
| <b>SVOCs</b>            |       |        |                         |                          |                         |
| 2-Methylnaphthalene     | µg/L  | 36     | <b>0.6</b>              | 5 U                      | 5 U                     |
| Acenaphthene            | µg/L  | 530    | <b>0.46</b>             | <b>0.03 J</b>            | 0.1 U                   |
| Acenaphthylene          | µg/L  | 530    | <b>0.04 J</b>           | 0.1 U                    | 0.1 U                   |
| Anthracene              | µg/L  | 1,800  | <b>0.04 J</b>           | 0.1 U                    | 0.1 U                   |
| Fluoranthene            | µg/L  | 800    | <b>0.16</b>             | <b>0.04 J</b>            | 0.1 U                   |
| Fluorene                | µg/L  | 290    | <b>0.41</b>             | 2 U                      | 2 U                     |
| Naphthalene             | µg/L  | 0.12   | <b>3.6</b>              | <b>0.08 J</b>            | 2 U                     |
| Phenanthrene            | µg/L  | --     | <b>0.43</b>             | 0.05 U                   | 0.05 U                  |
| Pyrene                  | µg/L  | 120    | <b>0.12</b>             | <b>0.05 J</b>            | 0.1 U                   |
| <b>VOCs</b>             |       |        |                         |                          |                         |
| Acetone                 | µg/L  | 14,000 | <b>4.1 J</b>            | <b>8.2</b>               | <b>2.1 J</b>            |
| Benzene                 | µg/L  | 5.0    | <b>0.23 J</b>           | 0.5 U                    | 0.5 U                   |
| Bromodichloromethane    | µg/L  | 0.13   | 0.5 U                   | <b>4.5</b>               | 0.5 U                   |
| Bromoform               | µg/L  | 3.3    | 2 U                     | <b>7.9</b>               | 2 U                     |
| Chloroform              | µg/L  | 0.22   | 0.75 U                  | <b>2.9</b>               | 0.75 U                  |
| Dibromochloromethane    | µg/L  | 0.17   | 0.50 U                  | <b>6.9</b>               | 0.50 U                  |
| Methyl tert butyl ether | µg/L  | 14     | 1 U                     | <b>0.56 J</b>            | 1 U                     |

**Detections in bold**

**Values in red indicate an exceedance of the groundwater Project Action Limit (PAL)**

*U: This analyte was not detected in the sample. The numeric value represents the sample quantitation/detection limit.*

*J: The positive result reported for this analyte is a quantitative estimate.*

µg/L: micrograms per liter

PAL: Project Action Limit



**Table 4 - Parcel B14 Surface Water Sampling (1st Half 2025)**  
**Summary of Organics Detected in Surface Water**

| Parameter                                         | Units | EPA Region 4<br>Surface Water<br>(Freshwater,<br>Chronic) | TMC-BEND<br>4/16/2025 | TMC-<br>OUTLET<br>4/16/2025 | TMC-RAIL<br>BRIDGE<br>4/16/2025 | TMC-TM04<br>4/16/2025 |
|---------------------------------------------------|-------|-----------------------------------------------------------|-----------------------|-----------------------------|---------------------------------|-----------------------|
| <b>SVOCs by EPA 8270D / PAHs by EPA 8270D SIM</b> |       |                                                           |                       |                             |                                 |                       |
| Acenaphthene                                      | µg/L  | 15                                                        | <b>0.03 J</b>         | <i>1 U</i>                  | <b>0.09 J</b>                   | <b>0.04 J</b>         |
| Benzo[a]anthracene                                | µg/L  | 4.7                                                       | <i>1 U</i>            | <b>0.06</b>                 | <b>0.03 J</b>                   | <b>0.03 J</b>         |
| Benzo[b]fluoranthene                              | µg/L  | 0.25                                                      | <i>0.05 U</i>         | <b>0.05 J</b>               | <i>0.05 U</i>                   | <i>0.05 U</i>         |
| Benzo(g,h,i)perylene                              | µg/L  | 0.012                                                     | <i>1 U</i>            | <b>0.03 J</b>               | <i>1 U</i>                      | <i>1 U</i>            |
| Chrysene                                          | µg/L  | 25                                                        | <i>0.5 U</i>          | <b>0.04 J</b>               | <i>0.5 U</i>                    | <i>0.5 U</i>          |
| Fluoranthene                                      | µg/L  | 0.8                                                       | <b>0.03 J</b>         | <b>0.16</b>                 | <b>0.06 J</b>                   | <b>0.03 J</b>         |
| Fluorene                                          | µg/L  | 19                                                        | <b>0.04 J</b>         | <b>0.06 J</b>               | <b>0.10</b>                     | <b>0.04 J</b>         |
| Indeno(1,2,3-cd)pyrene                            | µg/L  | 0.012                                                     | <i>1 U</i>            | <b>0.03 J</b>               | <i>1 U</i>                      | <i>1 U</i>            |
| Phenanthrene                                      | µg/L  | 2.3                                                       | <b>0.04 J</b>         | <b>0.08</b>                 | <b>0.10</b>                     | <b>0.05 J</b>         |
| Pyrene                                            | µg/L  | 4.6                                                       | <i>1 U</i>            | <b>0.15</b>                 | <i>1 U</i>                      | <i>1 U</i>            |
| <b>VOCs by EPA 8260C</b>                          |       |                                                           |                       |                             |                                 |                       |
| Acetone                                           | µg/L  | 1700                                                      | <b>4.1 J</b>          | <b>5.2</b>                  | <b>3.1 J</b>                    | <b>4.2 J</b>          |
| Benzene                                           | µg/L  | 160                                                       | <b>0.17 J</b>         | <b>0.42 J</b>               | <b>0.39 J</b>                   | <b>0.30 J</b>         |
| Carbon disulfide                                  | µg/L  | 15                                                        | <i>1 U</i>            | <b>3.7 J</b>                | <i>1 U</i>                      | <i>1 U</i>            |
| Chloroform                                        | µg/L  | 140                                                       | <i>1 U</i>            | <b>0.42 J</b>               | <i>1 U</i>                      | <i>1 U</i>            |
| Toluene                                           | µg/L  | 62                                                        | <i>1 U</i>            | <i>1 U</i>                  | <b>0.29 J</b>                   | <b>0.20 J</b>         |

**Detections in bold**

**Exceedances of EPA Region 4 Surface Water Screening Levels (freshwater, chronic) in red**

*U: This analyte was not detected in the sample. The numeric value represents the sample quantitation/detection limit*

*B: The analyte was detected above the reporting limit in the associated method blank.*

*J: The positive result reported for this analyte is a quantitative estimate.*

µg/L: micrograms per liter

PAL: Project Action Limit

**Table 5 - Parcel B14 (1st Half 2025)  
Historic Surface Water Sampling Results**

| Benzene (µg/L)  |        |                       |        |        |         |         |         |         |         |         |         |          |         |        |         |         |        |         |         |
|-----------------|--------|-----------------------|--------|--------|---------|---------|---------|---------|---------|---------|---------|----------|---------|--------|---------|---------|--------|---------|---------|
| Location        | GW PAL | EPA<br>Region 4<br>SW | 2/4/22 | 2/9/22 | 3/21/22 | 4/19/22 | 5/18/22 | 6/16/22 | 7/29/22 | 8/23/22 | 9/19/22 | 10/26/22 | 1/12/23 | 4/6/23 | 7/18/23 | 10/4/23 | 2/8/24 | 6/11/24 | 4/16/25 |
| TMC-Bend        | 5      | 160                   | NS     | NS     | 0.50 U  | 0.5 U   | 0.50 U  | 0.50 U  | 0.5 U   | 0.50 U  | 0.21 J  | 0.50 U   | 0.50 U  | 0.5 U  | 0.25 J  | 0.5 U   | 0.38 J | 0.19 J  | 0.17 J  |
| TMC-Outlet      | 5      | 160                   | 1 U    | 1.84   | 0.50 U  | 0.54    | 0.82    | 0.50 U  | 0.17 J  | 2.1     | 0.05 U  | 1.0      | 0.20 J  | 0.62   | 0.62    | 0.5 U   | 0.50   | 0.27 J  | 0.42 J  |
| TMC-Rail Bridge | 5      | 160                   | 5 U    | 1 U    | 0.50 U  | 0.5 U   | 0.50 U  | 0.50 U  | 0.5 U   | 0.16 J  | 0.22 J  | 0.50 U   | 0.50 J  | 0.5 U  | 0.22 J  | 0.5 U   | 0.38 J | 0.21 J  | 0.39 J  |
| TMC-TM04        | 5      | 160                   | 5 U    | 3.49   | 0.16 J  | 2.0     | 1.1     | 0.50 U  | 0.5 U   | 1.2     | 0.29 J  | 0.50 U   | 0.20 J  | 0.51   | 0.51    | 0.5 U   | 0.52   | 0.25 J  | 0.30 J  |

| Naphthalene (µg/L) |        |                       |        |        |         |         |         |         |         |         |         |          |         |        |         |         |        |         |         |
|--------------------|--------|-----------------------|--------|--------|---------|---------|---------|---------|---------|---------|---------|----------|---------|--------|---------|---------|--------|---------|---------|
| Location           | GW PAL | EPA<br>Region 4<br>SW | 2/4/22 | 2/9/22 | 3/21/22 | 4/19/22 | 5/18/22 | 6/16/22 | 7/29/22 | 8/23/22 | 9/19/22 | 10/26/22 | 1/12/23 | 4/6/23 | 7/18/23 | 10/4/23 | 2/8/24 | 6/11/24 | 4/16/25 |
| TMC-Bend           | 0.12   | 21                    | NS     | NS     | 0.03 J  | 0.62    | 0.06 J  | 0.06 J  | 0.17    | 0.10 U  | 0.11    | 0.10 U   | 0.10 U  | 0.15   | 0.15    | 0.13    | 2 U    | 0.1 U   | 0.1 U   |
| TMC-Outlet         | 0.12   | 21                    | 1 U    | 3.33   | 0.03 J  | 1.2     | 0.10 U  | 0.10 U  | 0.22    | 0.10 U  | 0.06 J  | 0.10 U   | 0.10 U  | 0.1 U  | 0.1 U   | 0.1 U   | 0.88 J | 0.1 U   | 0.1 U   |
| TMC-Rail Bridge    | 0.12   | 21                    | 1 U    | 1 U    | 0.10 U  | 0.58    | 0.05 J  | 0.10 U  | 1.1     | 0.10 U  | 0.11    | 0.06 J   | 0.10 U  | 0.1 U  | 0.27    | 0.08 J  | 1.4 J  | 0.07 J  | 0.1 U   |
| TMC-TM04           | 0.12   | 21                    | 1 U    | 1.3    | 0.03 J  | 0.79    | 0.10 U  | 0.12    | 0.12    | 0.10 U  | 0.14    | 0.10 U   | 0.10 U  | 0.34   | 0.34    | 0.1 U   | 0.63 J | 0.1 U   | 0.1 U   |

**Detections in bold**

**Exceedances of EPA Region 4 Surface Water Screening Levels (freshwater, chronic) in red**

*U: This analyte was not detected in the sample. The numeric value represents the sample quantitation / detection limit.*

*J: The positive result reported for this analyte is a quantitative estimate.*

*B: The analyte was detected above the reporting limits in the associated method blank.*

µg/L: micrograms per liter

PAL: Project Action Limit

Starting in March 2022, naphthalene was analyzed by SIM PAH.



**Table 6 - Parcel B14  
Groundwater Elevations**

| Well Designation | Installation Date | Top of Casing<br>(ft, AMSL) | April 2025             |                                        |
|------------------|-------------------|-----------------------------|------------------------|----------------------------------------|
|                  |                   |                             | Depth to<br>Water (ft) | Groundwater<br>Elevation<br>(ft, AMSL) |
| B24-001-MWS      | 4/27/2022         | 11.30                       | 10.80                  | 0.50                                   |
| B24-002-MWS      | 4/28/2022         | 12.43                       | 11.31                  | 1.12                                   |
| HI02R-PZM006 (b) | 1/26/2024         | 11.61                       | 10.00                  | 1.61                                   |
| HI07R-PZM005 (b) | 1/25/2024         | 11.65                       | 10.08                  | 1.57                                   |
| HI10-MWS         | 12/21/2021        | 12.50                       | 11.24                  | 1.26                                   |
| HI11-MWS         | 12/30/2021        | 13.08                       | 15.02                  | -1.94                                  |
| HI12-MWS         | 12/30/2021        | 11.54                       | 9.11                   | 2.43                                   |
| HI13-MWS         | 12/22/2021        | 11.29                       | 10.26                  | 1.03                                   |
| HI14-MWS         | 12/22/2021        | 14.65                       | 13.61                  | 1.04                                   |
| HI15-MWS         | 1/3/2022          | 13.43                       | 12.20                  | 1.23                                   |
| HI16-MWS         | 12/22/2021        | 10.53                       | 9.24                   | 1.29                                   |
| HI17-MWS (a)     | 12/23/2021        | 10.11                       | DRY                    | DRY                                    |
| HI18-MWS         | 12/23/2021        | 9.35                        | 8.12                   | 1.23                                   |
| HI19-MWS         | 12/29/2021        | 10.18                       | 8.82                   | 1.36                                   |
| HI20R-MWS (a, b) | 1/25/2024         | 11.49                       | DRY                    | DRY                                    |
| HI20R2-MWS       | 3/20/2024         | (c)                         | 17.91                  | (c)                                    |
| HI21-MWS         | 4/27/2022         | 12.56                       | 11.96                  | 0.60                                   |
| HI22-MWS         | 12/30/2021        | 11.49                       | 12.95                  | -1.46                                  |
| TM04-PZM006      | 3/25/1983         | 12.00                       | 12.21                  | -0.21                                  |
| TM08R-PZM007     | 11/19/2021        | 9.92                        | 8.64                   | 1.28                                   |
| TS03-PDP002      | 8/11/2000         | 13.53                       | 12.97                  | 0.56                                   |

Notes:

- (a) Absorbent socks utilized to address minor NAPL impacts.
- (b) Replacement well installed in January 2024 and surveyed in March 2024.
- (c) Replacement well was installed in March 2024, has not been surveyed.

**Table 7 - Parcel B14**  
**NAPL Gauging Activities**

| Sample ID  | Installation Date | Well Total Depth (Feet bgs) | Screen Interval (Feet bgs) | Riser Stick-Up (Feet) | 1/6/2025                 |                           |                       | 2/18/2025                |                           |                       |
|------------|-------------------|-----------------------------|----------------------------|-----------------------|--------------------------|---------------------------|-----------------------|--------------------------|---------------------------|-----------------------|
|            |                   |                             |                            |                       | Depth to NAPL (Feet TOC) | Depth to Water (Feet TOC) | NAPL Thickness (Feet) | Depth to NAPL (Feet TOC) | Depth to Water (Feet TOC) | NAPL Thickness (Feet) |
| HI17-MWS   | 12/23/2021        | 20                          | 5-20                       | 2.13                  | NM                       | NM                        | NM                    | NM                       | NM                        | NM                    |
| HI20R-MWS  | 1/25/2024         | 15                          | 5-15                       | 3                     | <b>10.38</b>             | <b>10.4</b>               | <b>0.02</b>           | NM                       | NM                        | NM                    |
| HI20R2-MWS | 3/20/2024         | 27.00                       | 7- 22                      | 3.00                  | NA                       | <b>18.47</b>              | NA                    | NA                       | <b>17.94</b>              | NA                    |

NM = Not Measured

**SHADED** = NAPL Detection

bgs = below ground surface

TOC = Top of Casing

HI17-MWS: Well not accessible due to deep snow drifts in the area.

HI20R-MWS: Absorbent sock deployed.

HI20R2-MWS: No removal action taken.

HI17-MWS: Absorbent sock removed, trace product observed.

HI20R-MWS: Absorbent sock removed, 1/4 saturated with product.

HI20R2-MWS: No removal action taken.



**Table 7 - Parcel B14  
NAPL Gauging Activities**

| Sample ID  | Installation Date | Well Total Depth (Feet bgs) | Screen Interval (Feet bgs) | Riser Stick-Up (Feet) | 3/19/2025                |                           |                       | 4/14/2025                |                           |                       |
|------------|-------------------|-----------------------------|----------------------------|-----------------------|--------------------------|---------------------------|-----------------------|--------------------------|---------------------------|-----------------------|
|            |                   |                             |                            |                       | Depth to NAPL (Feet TOC) | Depth to Water (Feet TOC) | NAPL Thickness (Feet) | Depth to NAPL (Feet TOC) | Depth to Water (Feet TOC) | NAPL Thickness (Feet) |
| HI17-MWS   | 12/23/2021        | 20                          | 5-20                       | 2.13                  | Trace                    | 10.13                     | NM                    | NA                       | DRY                       | NM                    |
| HI20R-MWS  | 1/25/2024         | 15                          | 5-15                       | 3                     | Trace                    | 10.38                     | NM                    | NA                       | DRY                       | NM                    |
| HI20R2-MWS | 3/20/2024         | 27.00                       | 7- 22                      | 3.00                  | NA                       | 18.18                     | NA                    | NA                       | 17.91                     | NA                    |

NM = Not Measured

**SHADED** = NAPL Detection

bgs = below ground surface

TOC = Top of Casing

HI17-MWS: Absorbent sock deployed.

HI20R-MWS: Absorbent sock deployed.

HI20R2-MWS: No removal action taken.

HI17-MWS: Absorbent sock removed, trace product observed.

HI20R-MWS: Absorbent sock removed, trace product observed.

HI20R2-MWS: No removal action taken.

**Table 7 - Parcel B14  
NAPL Gauging Activities**

| Sample ID  | Installation Date | Well Total Depth (Feet bgs) | Screen Interval (Feet bgs) | Riser Stick-Up (Feet) | 5/9/2025                 |                           |                       | 6/18/2025                |                           |                       |
|------------|-------------------|-----------------------------|----------------------------|-----------------------|--------------------------|---------------------------|-----------------------|--------------------------|---------------------------|-----------------------|
|            |                   |                             |                            |                       | Depth to NAPL (Feet TOC) | Depth to Water (Feet TOC) | NAPL Thickness (Feet) | Depth to NAPL (Feet TOC) | Depth to Water (Feet TOC) | NAPL Thickness (Feet) |
| HI17-MWS   | 12/23/2021        | 20                          | 5-20                       | 2.13                  | Trace                    | 10.12                     | NM                    | NA                       | DRY                       | NM                    |
| HI20R-MWS  | 1/25/2024         | 15                          | 5-15                       | 3                     | Trace                    | 10.35                     | NM                    | NA                       | DRY                       | NM                    |
| HI20R2-MWS | 3/20/2024         | 27.00                       | 7- 22                      | 3.00                  | NA                       | 18.16                     | NA                    | NA                       | 17.83                     | NA                    |

NM = Not Measured

**SHADED** = NAPL Detection

bgs = below ground surface

TOC = Top of Casing

HI17-MWS: Absorbent sock deployed.

HI20R-MWS: Absorbent sock deployed.

HI20R2-MWS: No removal action taken.

HI17-MWS: Absorbent sock removed, 1/3 saturated with product.

HI20R-MWS: Absorbent sock removed, 1/4 saturated with product.

HI20R2-MWS: No removal action taken.



---

---

## APPENDIX A

---

---



## ANALYTICAL REPORT

|                 |                                                                             |
|-----------------|-----------------------------------------------------------------------------|
| Lab Number:     | L2521696                                                                    |
| Client:         | Tradepoint Atlantic<br>1600 Sparrows Point Boulevard<br>Baltimore, MD 21219 |
| ATTN:           | Matthew Newman                                                              |
| Phone:          | (443) 649-5073                                                              |
| Project Name:   | B14 1ST HALF 2025 GW                                                        |
| Project Number: | 21010214                                                                    |
| Report Date:    | 04/18/25                                                                    |

The original project report/data package is held by Pace Analytical Services. This report/data package is paginated and should be reproduced only in its entirety. Pace Analytical Services holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0826), IL (200077), IN (C-MA-03), KY (KY98045), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), OR (MA-1316), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #525-23-122-91930A1).

---

Eight Walkup Drive, Westborough, MA 01581-1019  
508-898-9220 (Fax) 508-898-9193 800-624-9220 - [www.alphalab.com](http://www.alphalab.com)





**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

| Lab<br>Sample ID | Client ID    | Matrix | Sample<br>Location | Collection<br>Date/Time | Receive Date |
|------------------|--------------|--------|--------------------|-------------------------|--------------|
| L2521696-01      | TS03-PDP002  | WATER  | B14                | 04/08/25 13:20          | 04/09/25     |
| L2521696-02      | B24-001-MWS  | WATER  | B14                | 04/08/25 13:55          | 04/09/25     |
| L2521696-03      | HI02R-PZM006 | WATER  | B14                | 04/08/25 14:45          | 04/09/25     |
| L2521696-04      | TB-001       | WATER  | B14                | 04/08/25 00:00          | 04/09/25     |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

### Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Pace Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments and solids are reported on a dry weight basis unless otherwise noted. Tissues are reported "as received" or on a wet weight basis, unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

**HOLD POLICY** - For samples submitted on hold, Pace's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Pace Project Manager and made arrangements for Pace to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.

---



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

### Case Narrative (continued)

#### Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

#### Semivolatile Organics

L2521696-01, -02, and -03: The volume of sample received for the analysis deviates from the laboratory's current volume requirements. An aliquot was taken from the original sample container, extracted, analyzed, and reported accordingly.

#### Semivolatile Organics by SIM

L2521696-01, -02, and -03: The volume of sample received for the analysis deviates from the laboratory's current volume requirements. An aliquot was taken from the original sample container, extracted, analyzed, and reported accordingly.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

*Tiffani Morrissey* - Tiffani Morrissey

Title: Technical Director/Representative

Date: 04/18/25

# ORGANICS

# VOLATILES



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-01  
**Client ID:** TS03-PDP002  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/15/25 13:13  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 2.1    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS****Lab ID:** L2521696-01**Date Collected:** 04/08/25 13:20**Client ID:** TS03-PDP002**Date Received:** 04/09/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 118        |           | 70-130              |
| Toluene-d8            | 97         |           | 70-130              |
| 4-Bromofluorobenzene  | 98         |           | 70-130              |
| Dibromofluoromethane  | 108        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-01  
**Client ID:** TS03-PDP002  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/15/25 13:13  
**Analyst:** MCM

| Parameter                                        | Result     | Qualifier | Units     | RL                  | MDL   | Dilution Factor |
|--------------------------------------------------|------------|-----------|-----------|---------------------|-------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |            |           |           |                     |       |                 |
| 1,1,2,2-Tetrachloroethane                        | ND         |           | ug/l      | 0.050               | 0.006 | 1               |
| Surrogate                                        | % Recovery |           | Qualifier | Acceptance Criteria |       |                 |
| 1,2-Dichloroethane-d4                            | 116        |           |           | 70-130              |       |                 |
| 4-Bromofluorobenzene                             | 95         |           |           | 70-130              |       |                 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-02  
**Client ID:** B24-001-MWS  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/15/25 13:39  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 4.1    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.23   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS****Lab ID:** L2521696-02**Date Collected:** 04/08/25 13:55**Client ID:** B24-001-MWS**Date Received:** 04/09/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 120        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 109        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-02  
**Client ID:** B24-001-MWS  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/15/25 13:39  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 116        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-03  
**Client ID:** HI02R-PZM006  
**Sample Location:** B14

**Date Collected:** 04/08/25 14:45  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/15/25 14:05  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 3.1    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.42   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521696-03  
 Client ID: HI02R-PZM006  
 Sample Location: B14

Date Collected: 04/08/25 14:45  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 120        |           | 70-130              |
| Toluene-d8            | 101        |           | 70-130              |
| 4-Bromofluorobenzene  | 95         |           | 70-130              |
| Dibromofluoromethane  | 110        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

Lab ID: L2521696-03  
 Client ID: HI02R-PZM006  
 Sample Location: B14

Date Collected: 04/08/25 14:45  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/15/25 14:05  
 Analyst: MCM

| Parameter                                        | Result     | Qualifier | Units     | RL                  | MDL   | Dilution Factor |
|--------------------------------------------------|------------|-----------|-----------|---------------------|-------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |            |           |           |                     |       |                 |
| 1,1,2,2-Tetrachloroethane                        | ND         |           | ug/l      | 0.050               | 0.006 | 1               |
| Surrogate                                        | % Recovery |           | Qualifier | Acceptance Criteria |       |                 |
| 1,2-Dichloroethane-d4                            | 115        |           |           | 70-130              |       |                 |
| 4-Bromofluorobenzene                             | 93         |           |           | 70-130              |       |                 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-04  
**Client ID:** TB-001  
**Sample Location:** B14

**Date Collected:** 04/08/25 00:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/15/25 12:20  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | ND     |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS****Lab ID:** L2521696-04**Date Collected:** 04/08/25 00:00**Client ID:** TB-001**Date Received:** 04/09/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 117        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 99         |           | 70-130              |
| Dibromofluoromethane  | 108        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-04  
**Client ID:** TB-001  
**Sample Location:** B14

**Date Collected:** 04/08/25 00:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/15/25 12:20  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 114        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 95         |           | 70-130              |                 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D-SIM(M)  
Analytical Date: 04/15/25 08:48  
Analyst: MCM

| Parameter                                                                                | Result | Qualifier | Units | RL    | MDL   |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-------|-------|
| Volatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-04 Batch: WG2054023-5 |        |           |       |       |       |
| 1,1,1,2-Tetrachloroethane                                                                | ND     |           | ug/l  | 0.050 | 0.006 |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 115       |           | 70-130                 |
| 4-Bromofluorobenzene  | 95        |           | 70-130                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 04/15/25 07:29  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2054026-5 |        |           |       |      |      |
| Dichlorodifluoromethane                                                              | ND     |           | ug/l  | 5.0  | 0.24 |
| Chloromethane                                                                        | ND     |           | ug/l  | 2.5  | 0.20 |
| Vinyl chloride                                                                       | ND     |           | ug/l  | 1.0  | 0.07 |
| Bromomethane                                                                         | ND     |           | ug/l  | 1.0  | 0.26 |
| Chloroethane                                                                         | ND     |           | ug/l  | 1.0  | 0.13 |
| Trichlorofluoromethane                                                               | ND     |           | ug/l  | 2.5  | 0.16 |
| 1,1-Dichloroethene                                                                   | ND     |           | ug/l  | 0.50 | 0.17 |
| Carbon disulfide                                                                     | ND     |           | ug/l  | 5.0  | 0.30 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                | ND     |           | ug/l  | 2.5  | 0.15 |
| Methylene chloride                                                                   | ND     |           | ug/l  | 2.5  | 0.68 |
| Acetone                                                                              | ND     |           | ug/l  | 5.0  | 1.5  |
| trans-1,2-Dichloroethene                                                             | ND     |           | ug/l  | 0.75 | 0.16 |
| Methyl Acetate                                                                       | ND     |           | ug/l  | 2.0  | 0.23 |
| Methyl tert butyl ether                                                              | ND     |           | ug/l  | 1.0  | 0.17 |
| 1,1-Dichloroethane                                                                   | ND     |           | ug/l  | 0.75 | 0.21 |
| cis-1,2-Dichloroethene                                                               | ND     |           | ug/l  | 0.50 | 0.19 |
| Cyclohexane                                                                          | ND     |           | ug/l  | 10   | 0.27 |
| Chloroform                                                                           | ND     |           | ug/l  | 0.75 | 0.22 |
| Carbon tetrachloride                                                                 | ND     |           | ug/l  | 0.50 | 0.13 |
| 1,1,1-Trichloroethane                                                                | ND     |           | ug/l  | 0.50 | 0.16 |
| 2-Butanone                                                                           | ND     |           | ug/l  | 5.0  | 1.9  |
| Benzene                                                                              | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,2-Dichloroethane                                                                   | ND     |           | ug/l  | 0.50 | 0.13 |
| Trichloroethene                                                                      | ND     |           | ug/l  | 0.50 | 0.18 |
| 1,2-Dichloropropane                                                                  | ND     |           | ug/l  | 1.0  | 0.14 |
| Bromodichloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.19 |
| cis-1,3-Dichloropropene                                                              | ND     |           | ug/l  | 0.50 | 0.14 |
| Toluene                                                                              | ND     |           | ug/l  | 0.75 | 0.20 |
| Tetrachloroethene                                                                    | ND     |           | ug/l  | 0.50 | 0.18 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D  
 Analytical Date: 04/15/25 07:29  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2054026-5 |        |           |       |      |      |
| 4-Methyl-2-pentanone                                                                 | ND     |           | ug/l  | 5.0  | 0.42 |
| trans-1,3-Dichloropropene                                                            | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,3-Dichloropropene, Total                                                           | ND     |           | ug/l  | 0.50 | 0.14 |
| 1,1,2-Trichloroethane                                                                | ND     |           | ug/l  | 0.75 | 0.14 |
| Dibromochloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.15 |
| 1,2-Dibromoethane                                                                    | ND     |           | ug/l  | 2.0  | 0.19 |
| 2-Hexanone                                                                           | ND     |           | ug/l  | 5.0  | 0.52 |
| Chlorobenzene                                                                        | ND     |           | ug/l  | 0.50 | 0.18 |
| Ethylbenzene                                                                         | ND     |           | ug/l  | 0.50 | 0.17 |
| p/m-Xylene                                                                           | ND     |           | ug/l  | 1.0  | 0.33 |
| o-Xylene                                                                             | ND     |           | ug/l  | 1.0  | 0.39 |
| Xylenes, Total                                                                       | ND     |           | ug/l  | 1.0  | 0.33 |
| Styrene                                                                              | ND     |           | ug/l  | 1.0  | 0.36 |
| Bromoform                                                                            | ND     |           | ug/l  | 2.0  | 0.25 |
| Isopropylbenzene                                                                     | ND     |           | ug/l  | 0.50 | 0.19 |
| 1,1,2,2-Tetrachloroethane                                                            | ND     |           | ug/l  | 0.50 | 0.17 |
| 1,3-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,4-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,2-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.18 |
| 1,2-Dibromo-3-chloropropane                                                          | ND     |           | ug/l  | 2.5  | 0.35 |
| 1,2,4-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.22 |
| 1,2,3-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.23 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
Analytical Date: 04/15/25 07:29  
Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL | MDL |
|--------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2054026-5 |        |           |       |    |     |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 118       |           | 70-130                 |
| Toluene-d8            | 100       |           | 70-130                 |
| 4-Bromofluorobenzene  | 98        |           | 70-130                 |
| Dibromofluoromethane  | 108       |           | 70-130                 |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521696

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2054023-3 WG2054023-4 |                          |             |                           |             |                             |            |             |                       |
| 1,1,2,2-Tetrachloroethane                                                                                   | 103                      |             | 98                        |             | 70-130                      | 5          |             | 25                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 114                      |             | 114                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 95                       |             | 95                        |             | 70-130                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521696

**Report Date:** 04/18/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2054026-3 WG2054026-4 |                  |      |                   |      |                     |     |      |               |
| Dichlorodifluoromethane                                                                                 | 81               |      | 80                |      | 36-147              | 1   |      | 20            |
| Chloromethane                                                                                           | 100              |      | 100               |      | 64-130              | 0   |      | 20            |
| Vinyl chloride                                                                                          | 100              |      | 99                |      | 55-140              | 1   |      | 20            |
| Bromomethane                                                                                            | 47               |      | 51                |      | 39-139              | 8   |      | 20            |
| Chloroethane                                                                                            | 120              |      | 110               |      | 55-138              | 9   |      | 20            |
| Trichlorofluoromethane                                                                                  | 100              |      | 100               |      | 62-150              | 0   |      | 20            |
| 1,1-Dichloroethene                                                                                      | 97               |      | 96                |      | 61-145              | 1   |      | 20            |
| Carbon disulfide                                                                                        | 100              |      | 100               |      | 51-130              | 0   |      | 20            |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                                   | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Methylene chloride                                                                                      | 99               |      | 96                |      | 70-130              | 3   |      | 20            |
| Acetone                                                                                                 | 120              |      | 120               |      | 58-148              | 0   |      | 20            |
| trans-1,2-Dichloroethene                                                                                | 100              |      | 96                |      | 70-130              | 4   |      | 20            |
| Methyl Acetate                                                                                          | 130              |      | 130               |      | 70-130              | 0   |      | 20            |
| Methyl tert butyl ether                                                                                 | 96               |      | 94                |      | 63-130              | 2   |      | 20            |
| 1,1-Dichloroethane                                                                                      | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| cis-1,2-Dichloroethene                                                                                  | 97               |      | 94                |      | 70-130              | 3   |      | 20            |
| Cyclohexane                                                                                             | 120              |      | 110               |      | 70-130              | 9   |      | 20            |
| Chloroform                                                                                              | 100              |      | 98                |      | 70-130              | 2   |      | 20            |
| Carbon tetrachloride                                                                                    | 100              |      | 100               |      | 63-132              | 0   |      | 20            |
| 1,1,1-Trichloroethane                                                                                   | 100              |      | 97                |      | 67-130              | 3   |      | 20            |
| 2-Butanone                                                                                              | 110              |      | 100               |      | 63-138              | 10  |      | 20            |
| Benzene                                                                                                 | 98               |      | 95                |      | 70-130              | 3   |      | 20            |
| 1,2-Dichloroethane                                                                                      | 110              |      | 100               |      | 70-130              | 10  |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521696

**Report Date:** 04/18/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2054026-3 WG2054026-4 |                  |      |                   |      |                     |     |      |               |
| Trichloroethene                                                                                         | 97               |      | 94                |      | 70-130              | 3   |      | 20            |
| 1,2-Dichloropropane                                                                                     | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Bromodichloromethane                                                                                    | 96               |      | 92                |      | 67-130              | 4   |      | 20            |
| cis-1,3-Dichloropropene                                                                                 | 91               |      | 90                |      | 70-130              | 1   |      | 20            |
| Toluene                                                                                                 | 96               |      | 93                |      | 70-130              | 3   |      | 20            |
| Tetrachloroethene                                                                                       | 95               |      | 94                |      | 70-130              | 1   |      | 20            |
| 4-Methyl-2-pentanone                                                                                    | 87               |      | 83                |      | 59-130              | 5   |      | 20            |
| trans-1,3-Dichloropropene                                                                               | 96               |      | 94                |      | 70-130              | 2   |      | 20            |
| 1,1,2-Trichloroethane                                                                                   | 98               |      | 95                |      | 70-130              | 3   |      | 20            |
| Dibromochloromethane                                                                                    | 88               |      | 88                |      | 63-130              | 0   |      | 20            |
| 1,2-Dibromoethane                                                                                       | 92               |      | 91                |      | 70-130              | 1   |      | 20            |
| 2-Hexanone                                                                                              | 96               |      | 94                |      | 57-130              | 2   |      | 20            |
| Chlorobenzene                                                                                           | 94               |      | 92                |      | 75-130              | 2   |      | 20            |
| Ethylbenzene                                                                                            | 97               |      | 96                |      | 70-130              | 1   |      | 20            |
| p/m-Xylene                                                                                              | 95               |      | 95                |      | 70-130              | 0   |      | 20            |
| o-Xylene                                                                                                | 95               |      | 95                |      | 70-130              | 0   |      | 20            |
| Styrene                                                                                                 | 90               |      | 90                |      | 70-130              | 0   |      | 20            |
| Bromoform                                                                                               | 83               |      | 82                |      | 54-136              | 1   |      | 20            |
| Isopropylbenzene                                                                                        | 95               |      | 93                |      | 70-130              | 2   |      | 20            |
| 1,1,2,2-Tetrachloroethane                                                                               | 92               |      | 91                |      | 67-130              | 1   |      | 20            |
| 1,3-Dichlorobenzene                                                                                     | 94               |      | 93                |      | 70-130              | 1   |      | 20            |
| 1,4-Dichlorobenzene                                                                                     | 93               |      | 91                |      | 70-130              | 2   |      | 20            |
| 1,2-Dichlorobenzene                                                                                     | 92               |      | 88                |      | 70-130              | 4   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521696

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b>                                                                                        | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|---------------------------------------------------------------------------------------------------------|--------------------------|-------------|--------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2054026-3 WG2054026-4 |                          |             |                          |             |                             |            |             |                       |
| 1,2-Dibromo-3-chloropropane                                                                             | 76                       |             | 77                       |             | 41-144                      | 1          |             | 20                    |
| 1,2,4-Trichlorobenzene                                                                                  | 87                       |             | 85                       |             | 70-130                      | 2          |             | 20                    |
| 1,2,3-Trichlorobenzene                                                                                  | 88                       |             | 89                       |             | 70-130                      | 1          |             | 20                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|--------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 112                      |             | 112                      |             | 70-130                         |
| Toluene-d8            | 101                      |             | 101                      |             | 70-130                         |
| 4-Bromofluorobenzene  | 97                       |             | 99                       |             | 70-130                         |
| Dibromofluoromethane  | 102                      |             | 102                      |             | 70-130                         |



# SEMIVOLATILES

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-01  
**Client ID:** TS03-PDP002  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 23:12  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-01  
**Client ID:** TS03-PDP002  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 53         |           | 21-120              |
| Phenol-d6            | 41         |           | 10-120              |
| Nitrobenzene-d5      | 68         |           | 23-120              |
| 2-Fluorobiphenyl     | 75         |           | 15-120              |
| 2,4,6-Tribromophenol | 76         |           | 10-120              |
| 4-Terphenyl-d14      | 82         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-01  
**Client ID:** TS03-PDP002  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 10:36  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 58         |           | 21-120              |
| Phenol-d6            | 48         |           | 10-120              |
| Nitrobenzene-d5      | 99         |           | 23-120              |
| 2-Fluorobiphenyl     | 85         |           | 15-120              |
| 2,4,6-Tribromophenol | 120        |           | 10-120              |
| 4-Terphenyl-d14      | 83         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-02  
**Client ID:** B24-001-MWS  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 23:35  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 3.2    |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | 0.40   | J         | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521696-02

Date Collected: 04/08/25 13:55

Client ID: B24-001-MWS

Date Received: 04/09/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | 0.40   | J         | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | 0.42   | J         | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 47         |           | 21-120              |
| Phenol-d6            | 36         |           | 10-120              |
| Nitrobenzene-d5      | 62         |           | 23-120              |
| 2-Fluorobiphenyl     | 71         |           | 15-120              |
| 2,4,6-Tribromophenol | 71         |           | 10-120              |
| 4-Terphenyl-d14      | 76         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-02  
**Client ID:** B24-001-MWS  
**Sample Location:** B14

**Date Collected:** 04/08/25 13:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 10:52  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 3.6    |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.60   |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.46   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.41   |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.43   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.16   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.12   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 58         |           | 21-120              |
| Phenol-d6            | 48         |           | 10-120              |
| Nitrobenzene-d5      | 97         |           | 23-120              |
| 2-Fluorobiphenyl     | 83         |           | 15-120              |
| 2,4,6-Tribromophenol | 121        | Q         | 10-120              |
| 4-Terphenyl-d14      | 81         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-03  
**Client ID:** HI02R-PZM006  
**Sample Location:** B14

**Date Collected:** 04/08/25 14:45  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 23:58  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | 5.5    |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 9.3    |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | 1.0    | J         | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | 0.24   | J         | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521696-03  
 Client ID: HI02R-PZM006  
 Sample Location: B14

Date Collected: 04/08/25 14:45  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | 1.3    | J         | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | 0.87   | J         | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | 0.83   | J         | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | 0.42   | J         | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 57         |           | 21-120              |
| Phenol-d6            | 44         |           | 10-120              |
| Nitrobenzene-d5      | 77         |           | 23-120              |
| 2-Fluorobiphenyl     | 80         |           | 15-120              |
| 2,4,6-Tribromophenol | 91         |           | 10-120              |
| 4-Terphenyl-d14      | 88         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521696-03  
**Client ID:** HI02R-PZM006  
**Sample Location:** B14

**Date Collected:** 04/08/25 14:45  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 11:09  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 8.0    |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.90   |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.05   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 1.4    |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 1.0    |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.96   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.97   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.15   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.10   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 63         |           | 21-120              |
| Phenol-d6            | 53         |           | 10-120              |
| Nitrobenzene-d5      | 110        |           | 23-120              |
| 2-Fluorobiphenyl     | 74         |           | 15-120              |
| 2,4,6-Tribromophenol | 132        | Q         | 10-120              |
| 4-Terphenyl-d14      | 79         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW**Project Number:** 21010214**Lab Number:** L2521696**Report Date:** 04/18/25

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270E  
 Analytical Date: 04/16/25 20:08  
 Analyst: SLR

Extraction Method: EPA 3510C  
 Extraction Date: 04/13/25 14:21

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-03 Batch: WG2053156-1 |        |           |       |     |      |
| Benzaldehyde                                                                             | ND     |           | ug/l  | 5.0 | 1.1  |
| Phenol                                                                                   | ND     |           | ug/l  | 5.0 | 0.35 |
| Bis(2-chloroethyl)ether                                                                  | ND     |           | ug/l  | 2.0 | 0.39 |
| 2-Chlorophenol                                                                           | ND     |           | ug/l  | 2.0 | 0.65 |
| 2-Methylphenol                                                                           | ND     |           | ug/l  | 5.0 | 2.3  |
| Bis(2-chloroisopropyl)ether                                                              | ND     |           | ug/l  | 2.0 | 0.40 |
| Acetophenone                                                                             | ND     |           | ug/l  | 5.0 | 0.92 |
| n-Nitrosodi-n-propylamine                                                                | ND     |           | ug/l  | 5.0 | 0.91 |
| 3-Methylphenol/4-Methylphenol                                                            | ND     |           | ug/l  | 5.0 | 1.4  |
| Hexachloroethane                                                                         | ND     |           | ug/l  | 2.0 | 0.20 |
| Nitrobenzene                                                                             | ND     |           | ug/l  | 1.4 | 0.20 |
| Isophorone                                                                               | ND     |           | ug/l  | 5.0 | 0.86 |
| 2,4-Dimethylphenol                                                                       | ND     |           | ug/l  | 5.0 | 2.0  |
| Bis(2-chloroethoxy)methane                                                               | ND     |           | ug/l  | 5.0 | 0.84 |
| 2,4-Dichlorophenol                                                                       | ND     |           | ug/l  | 5.0 | 1.7  |
| Naphthalene                                                                              | ND     |           | ug/l  | 2.0 | 0.54 |
| 4-Chloroaniline                                                                          | ND     |           | ug/l  | 3.7 | 0.47 |
| Hexachlorobutadiene                                                                      | ND     |           | ug/l  | 2.0 | 0.36 |
| Caprolactam                                                                              | ND     |           | ug/l  | 10  | 1.2  |
| 2-Methylnaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.37 |
| Hexachlorocyclopentadiene                                                                | ND     |           | ug/l  | 20  | 1.2  |
| 1,2,4,5-Tetrachlorobenzene                                                               | ND     |           | ug/l  | 1.7 | 0.24 |
| 2,4,6-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| 2,4,5-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| Biphenyl                                                                                 | ND     |           | ug/l  | 2.0 | 0.20 |
| 2-Chloronaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.35 |
| 2-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.0  |
| 2,6-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.84 |
| Acenaphthylene                                                                           | ND     |           | ug/l  | 2.0 | 0.32 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 20:08  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-03 Batch: WG2053156-1 |        |           |       |     |      |
| Acenaphthene                                                                             | ND     |           | ug/l  | 2.0 | 0.40 |
| 2,4-Dinitrophenol                                                                        | ND     |           | ug/l  | 20  | 5.4  |
| 2,4-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.54 |
| 2,3,4,6-Tetrachlorophenol                                                                | ND     |           | ug/l  | 5.0 | 2.2  |
| Diethyl phthalate                                                                        | ND     |           | ug/l  | 5.0 | 0.76 |
| Fluorene                                                                                 | ND     |           | ug/l  | 2.0 | 0.44 |
| 4-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.4  |
| NDPA/DPA                                                                                 | ND     |           | ug/l  | 2.0 | 0.92 |
| Hexachlorobenzene                                                                        | ND     |           | ug/l  | 2.0 | 0.45 |
| Pentachlorophenol                                                                        | ND     |           | ug/l  | 10  | 2.5  |
| Phenanthrene                                                                             | ND     |           | ug/l  | 2.0 | 0.42 |
| Anthracene                                                                               | ND     |           | ug/l  | 2.0 | 0.47 |
| Carbazole                                                                                | ND     |           | ug/l  | 2.0 | 0.31 |
| Di-n-butylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 0.96 |
| Fluoranthene                                                                             | ND     |           | ug/l  | 2.0 | 0.41 |
| Pyrene                                                                                   | ND     |           | ug/l  | 2.0 | 0.41 |
| 3,3'-Dichlorobenzidine                                                                   | ND     |           | ug/l  | 5.0 | 1.8  |
| Benzo(a)anthracene                                                                       | ND     |           | ug/l  | 2.0 | 0.32 |
| Chrysene                                                                                 | ND     |           | ug/l  | 1.5 | 0.22 |
| Bis(2-ethylhexyl)phthalate                                                               | ND     |           | ug/l  | 3.0 | 1.4  |
| Di-n-octylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 2.3  |
| Benzo(b)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.53 |
| Benzo(k)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.62 |
| Benzo(a)pyrene                                                                           | ND     |           | ug/l  | 2.0 | 0.37 |
| Indeno(1,2,3-cd)pyrene                                                                   | ND     |           | ug/l  | 2.0 | 0.48 |
| Dibenzo(a,h)anthracene                                                                   | ND     |           | ug/l  | 2.0 | 0.29 |
| Benzo(ghi)perylene                                                                       | ND     |           | ug/l  | 2.0 | 0.37 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8270E  
 Analytical Date: 04/16/25 20:08  
 Analyst: SLR

Extraction Method: EPA 3510C  
 Extraction Date: 04/13/25 14:21

| Parameter                                                                                | Result | Qualifier | Units | RL | MDL |
|------------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-03 Batch: WG2053156-1 |        |           |       |    |     |

| Surrogate            | %Recovery | Qualifier | Acceptance<br>Criteria |
|----------------------|-----------|-----------|------------------------|
| 2-Fluorophenol       | 53        |           | 21-120                 |
| Phenol-d6            | 37        |           | 10-120                 |
| Nitrobenzene-d5      | 68        |           | 23-120                 |
| 2-Fluorobiphenyl     | 72        |           | 15-120                 |
| 2,4,6-Tribromophenol | 67        |           | 10-120                 |
| 4-Terphenyl-d14      | 77        |           | 41-149                 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 06:54  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                                                                    | Result | Qualifier | Units | RL   | MDL  |
|----------------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-03 Batch: WG2053157-1 |        |           |       |      |      |
| Naphthalene                                                                                  | ND     |           | ug/l  | 0.10 | 0.02 |
| 2-Methylnaphthalene                                                                          | ND     |           | ug/l  | 0.10 | 0.03 |
| Acenaphthylene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Acenaphthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluorene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Phenanthrene                                                                                 | ND     |           | ug/l  | 0.05 | 0.04 |
| Anthracene                                                                                   | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluoranthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.03 |
| Pyrene                                                                                       | ND     |           | ug/l  | 0.10 | 0.04 |
| Benzo(a)anthracene                                                                           | ND     |           | ug/l  | 0.05 | 0.03 |
| Chrysene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(b)fluoranthene                                                                         | ND     |           | ug/l  | 0.05 | 0.03 |
| Benzo(k)fluoranthene                                                                         | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(a)pyrene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Indeno(1,2,3-cd)pyrene                                                                       | ND     |           | ug/l  | 0.10 | 0.02 |
| Dibenzo(a,h)anthracene                                                                       | ND     |           | ug/l  | 0.05 | 0.02 |
| Benzo(ghi)perylene                                                                           | ND     |           | ug/l  | 0.10 | 0.02 |

| Surrogate            | %Recovery  | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 76         |           | 21-120              |
| Phenol-d6            | 60         |           | 10-120              |
| Nitrobenzene-d5      | <b>121</b> | Q         | 23-120              |
| 2-Fluorobiphenyl     | 106        |           | 15-120              |
| 2,4,6-Tribromophenol | <b>140</b> | Q         | 10-120              |
| 4-Terphenyl-d14      | 109        |           | 41-149              |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521696

**Report Date:** 04/18/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-03 Batch: WG2053156-2 WG2053156-3 |                  |      |                   |      |                     |     |      |               |
| Benzaldehyde                                                                                                | 78               |      | 77                |      | 40-140              | 1   |      | 30            |
| Phenol                                                                                                      | 41               |      | 40                |      | 12-110              | 2   |      | 30            |
| Bis(2-chloroethyl)ether                                                                                     | 71               |      | 74                |      | 40-140              | 4   |      | 30            |
| 2-Chlorophenol                                                                                              | 70               |      | 69                |      | 27-123              | 1   |      | 30            |
| 2-Methylphenol                                                                                              | 70               |      | 70                |      | 30-130              | 0   |      | 30            |
| Bis(2-chloroisopropyl)ether                                                                                 | 77               |      | 77                |      | 40-140              | 0   |      | 30            |
| Acetophenone                                                                                                | 80               |      | 79                |      | 39-129              | 1   |      | 30            |
| n-Nitrosodi-n-propylamine                                                                                   | 71               |      | 69                |      | 29-132              | 3   |      | 30            |
| 3-Methylphenol/4-Methylphenol                                                                               | 67               |      | 72                |      | 30-130              | 7   |      | 30            |
| Hexachloroethane                                                                                            | 63               |      | 65                |      | 40-140              | 3   |      | 30            |
| Nitrobenzene                                                                                                | 68               |      | 69                |      | 40-140              | 1   |      | 30            |
| Isophorone                                                                                                  | 68               |      | 69                |      | 40-140              | 1   |      | 30            |
| 2,4-Dimethylphenol                                                                                          | 64               |      | 59                |      | 30-130              | 8   |      | 30            |
| Bis(2-chloroethoxy)methane                                                                                  | 72               |      | 73                |      | 40-140              | 1   |      | 30            |
| 2,4-Dichlorophenol                                                                                          | 81               |      | 80                |      | 30-130              | 1   |      | 30            |
| Naphthalene                                                                                                 | 71               |      | 72                |      | 40-140              | 1   |      | 30            |
| 4-Chloroaniline                                                                                             | 48               |      | 57                |      | 40-140              | 17  |      | 30            |
| Hexachlorobutadiene                                                                                         | 64               |      | 68                |      | 40-140              | 6   |      | 30            |
| Caprolactam                                                                                                 | 40               |      | 44                |      | 10-130              | 10  |      | 30            |
| 2-Methylnaphthalene                                                                                         | 74               |      | 75                |      | 40-140              | 1   |      | 30            |
| Hexachlorocyclopentadiene                                                                                   | 67               |      | 70                |      | 40-140              | 4   |      | 30            |
| 1,2,4,5-Tetrachlorobenzene                                                                                  | 73               |      | 78                |      | 2-134               | 7   |      | 30            |
| 2,4,6-Trichlorophenol                                                                                       | 80               |      | 81                |      | 30-130              | 1   |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521696

**Report Date:** 04/18/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-03 Batch: WG2053156-2 WG2053156-3 |                  |      |                   |      |                     |     |      |               |
| 2,4,5-Trichlorophenol                                                                                       | 88               |      | 92                |      | 30-130              | 4   |      | 30            |
| Biphenyl                                                                                                    | 81               |      | 82                |      | 40-140              | 1   |      | 30            |
| 2-Chloronaphthalene                                                                                         | 75               |      | 76                |      | 40-140              | 1   |      | 30            |
| 2-Nitroaniline                                                                                              | 86               |      | 89                |      | 52-143              | 3   |      | 30            |
| 2,6-Dinitrotoluene                                                                                          | 85               |      | 89                |      | 40-140              | 5   |      | 30            |
| Acenaphthylene                                                                                              | 75               |      | 78                |      | 45-123              | 4   |      | 30            |
| Acenaphthene                                                                                                | 76               |      | 77                |      | 37-111              | 1   |      | 30            |
| 2,4-Dinitrophenol                                                                                           | 81               |      | 80                |      | 20-130              | 1   |      | 30            |
| 2,4-Dinitrotoluene                                                                                          | 82               |      | 84                |      | 48-143              | 2   |      | 30            |
| 2,3,4,6-Tetrachlorophenol                                                                                   | 88               |      | 87                |      | 54-145              | 1   |      | 30            |
| Diethyl phthalate                                                                                           | 78               |      | 78                |      | 40-140              | 0   |      | 30            |
| Fluorene                                                                                                    | 77               |      | 78                |      | 40-140              | 1   |      | 30            |
| 4-Nitroaniline                                                                                              | 76               |      | 78                |      | 51-143              | 3   |      | 30            |
| NDPA/DPA                                                                                                    | 81               |      | 82                |      | 40-140              | 1   |      | 30            |
| Hexachlorobenzene                                                                                           | 81               |      | 82                |      | 40-140              | 1   |      | 30            |
| Pentachlorophenol                                                                                           | 90               |      | 85                |      | 9-103               | 6   |      | 30            |
| Phenanthrene                                                                                                | 79               |      | 82                |      | 40-140              | 4   |      | 30            |
| Anthracene                                                                                                  | 82               |      | 85                |      | 40-140              | 4   |      | 30            |
| Carbazole                                                                                                   | 81               |      | 83                |      | 55-144              | 2   |      | 30            |
| Di-n-butylphthalate                                                                                         | 83               |      | 85                |      | 40-140              | 2   |      | 30            |
| Fluoranthene                                                                                                | 84               |      | 86                |      | 40-140              | 2   |      | 30            |
| Pyrene                                                                                                      | 86               |      | 88                |      | 26-127              | 2   |      | 30            |
| 3,3'-Dichlorobenzidine                                                                                      | 64               |      | 72                |      | 40-140              | 12  |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521696

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-03 Batch: WG2053156-2 WG2053156-3 |                          |             |                           |             |                             |            |             |                       |
| Benzo(a)anthracene                                                                                          | 80                       |             | 86                        |             | 40-140                      | 7          |             | 30                    |
| Chrysene                                                                                                    | 82                       |             | 86                        |             | 40-140                      | 5          |             | 30                    |
| Bis(2-ethylhexyl)phthalate                                                                                  | 88                       |             | 90                        |             | 40-140                      | 2          |             | 30                    |
| Di-n-octylphthalate                                                                                         | 85                       |             | 89                        |             | 40-140                      | 5          |             | 30                    |
| Benzo(b)fluoranthene                                                                                        | 84                       |             | 90                        |             | 40-140                      | 7          |             | 30                    |
| Benzo(k)fluoranthene                                                                                        | 87                       |             | 92                        |             | 40-140                      | 6          |             | 30                    |
| Benzo(a)pyrene                                                                                              | 87                       |             | 91                        |             | 40-140                      | 4          |             | 30                    |
| Indeno(1,2,3-cd)pyrene                                                                                      | 84                       |             | 89                        |             | 40-140                      | 6          |             | 30                    |
| Dibenzo(a,h)anthracene                                                                                      | 82                       |             | 87                        |             | 40-140                      | 6          |             | 30                    |
| Benzo(ghi)perylene                                                                                          | 80                       |             | 82                        |             | 40-140                      | 2          |             | 30                    |

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 54                       |             | 55                        |             | 21-120                         |
| Phenol-d6            | 42                       |             | 42                        |             | 10-120                         |
| Nitrobenzene-d5      | 67                       |             | 68                        |             | 23-120                         |
| 2-Fluorobiphenyl     | 74                       |             | 73                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 74                       |             | 77                        |             | 10-120                         |
| 4-Terphenyl-d14      | 76                       |             | 80                        |             | 41-149                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521696

**Report Date:** 04/18/25

| Parameter                                                                                                       | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-----------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-03 Batch: WG2053157-2 WG2053157-3 |                  |      |                   |      |                     |     |      |               |
| Naphthalene                                                                                                     | 69               |      | 75                |      | 40-140              | 8   |      | 40            |
| 2-Methylnaphthalene                                                                                             | 72               |      | 78                |      | 40-140              | 8   |      | 40            |
| Acenaphthylene                                                                                                  | 86               |      | 93                |      | 40-140              | 8   |      | 40            |
| Acenaphthene                                                                                                    | 75               |      | 82                |      | 37-111              | 9   |      | 40            |
| Fluorene                                                                                                        | 82               |      | 88                |      | 40-140              | 7   |      | 40            |
| Phenanthrene                                                                                                    | 78               |      | 85                |      | 40-140              | 9   |      | 40            |
| Anthracene                                                                                                      | 83               |      | 90                |      | 40-140              | 8   |      | 40            |
| Fluoranthene                                                                                                    | 82               |      | 88                |      | 40-140              | 7   |      | 40            |
| Pyrene                                                                                                          | 80               |      | 87                |      | 26-127              | 8   |      | 40            |
| Benzo(a)anthracene                                                                                              | 80               |      | 87                |      | 40-140              | 8   |      | 40            |
| Chrysene                                                                                                        | 83               |      | 90                |      | 40-140              | 8   |      | 40            |
| Benzo(b)fluoranthene                                                                                            | 87               |      | 94                |      | 40-140              | 8   |      | 40            |
| Benzo(k)fluoranthene                                                                                            | 90               |      | 99                |      | 40-140              | 10  |      | 40            |
| Benzo(a)pyrene                                                                                                  | 94               |      | 103               |      | 40-140              | 9   |      | 40            |
| Indeno(1,2,3-cd)pyrene                                                                                          | 104              |      | 114               |      | 40-140              | 9   |      | 40            |
| Dibenzo(a,h)anthracene                                                                                          | 103              |      | 112               |      | 40-140              | 8   |      | 40            |
| Benzo(ghi)perylene                                                                                              | 97               |      | 105               |      | 40-140              | 8   |      | 40            |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521696

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b>                                                                                                | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-----------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-03 Batch: WG2053157-2 WG2053157-3 |                          |             |                           |             |                             |            |             |                       |

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 58                       |             | 61                        |             | 21-120                         |
| Phenol-d6            | 48                       |             | 50                        |             | 10-120                         |
| Nitrobenzene-d5      | 87                       |             | 93                        |             | 23-120                         |
| 2-Fluorobiphenyl     | 76                       |             | 81                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 106                      |             | 115                       |             | 10-120                         |
| 4-Terphenyl-d14      | 71                       |             | 78                        |             | 41-149                         |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

**Cooler Information**

|               |                     |
|---------------|---------------------|
| <b>Cooler</b> | <b>Custody Seal</b> |
| A             | Absent              |

**Container Information**

| <b>Container ID</b> | <b>Container Type</b>   | <b>Cooler</b> | <b>Initial pH</b> | <b>Final pH</b> | <b>Temp deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen Date/Time</b> | <b>Analysis(*)</b>               |
|---------------------|-------------------------|---------------|-------------------|-----------------|-------------------|-------------|-------------|-------------------------|----------------------------------|
| L2521696-01A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-01B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-01C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-01D        | Amber 250ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2521696-01E        | Amber 250ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2521696-02A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-02B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-02C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-02D        | Amber 250ml unpreserved | A             | 10                | 10              | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2521696-02E        | Amber 250ml unpreserved | A             | 10                | 10              | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2521696-03A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-03B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-03C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-03D        | Amber 250ml unpreserved | A             | 10                | 10              | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521696-03E        | Amber 250ml unpreserved | A             | 10                | 10              | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521696-04A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-04B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-04C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521696-04D        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |

**Container Comments**

L2521696-01D      amber-a.25 received

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

Serial\_No:04182514:16  
**Lab Number:** L2521696  
**Report Date:** 04/18/25

**Container Information**

| Container ID | Container Type | Cooler | Initial<br>pH | Final<br>pH | Temp<br>deg C | Pres | Seal | Frozen<br>Date/Time | Analysis(*) |
|--------------|----------------|--------|---------------|-------------|---------------|------|------|---------------------|-------------|
|--------------|----------------|--------|---------------|-------------|---------------|------|------|---------------------|-------------|

**Container Comments**

|              |                     |
|--------------|---------------------|
| L2521696-01E | amber-a.25 received |
| L2521696-02D | amber-a.25 received |
| L2521696-02E | amber-a.25 received |
| L2521696-03D | amber-a.25 received |
| L2521696-03E | amber-a.25 received |

\*Values in parentheses indicate holding time in days



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

## GLOSSARY

### Acronyms

|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| DL       | - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                |
| EDL      | - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).                                                                                         |
| EMPC     | - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.                                                                                                                                                                                                                               |
| EPA      | - Environmental Protection Agency.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| LCS      | - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                          |
| LCSD     | - Laboratory Control Sample Duplicate: Refer to LCS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |
| LFB      | - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                         |
| LOD      | - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                                                               |
| LOQ      | - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)<br><br>Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| MDL      | - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                          |
| MS       | - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.                                                                                                                                                                                                    |
| MSD      | - Matrix Spike Sample Duplicate: Refer to MS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NA       | - Not Applicable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| NC       | - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.                                                                                                                                                                                                                                                                                                                                                                           |
| NDPA/DPA | - N-Nitrosodiphenylamine/Diphenylamine.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| NI       | - Not Ignitable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| NP       | - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NR       | - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.                                                                                                                                                                                                                                                                                                                                                                    |
| RL       | - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                                                                                                                                   |
| RPD      | - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.                                                                  |
| SRM      | - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.                                                                                                                                                                                                                                                                                                                                                                     |
| STLP     | - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| TEF      | - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.                                                                                                                                                                                                                                                                                                                                                                                             |
| TEQ      | - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.                                                                                                                                                                                                                                                                                                                                                        |
| TIC      | - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.                                                                                                                                                                                                                                                                      |

Report Format: DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521696  
**Report Date:** 04/18/25

### Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

### Terms

**Analytical Method:** Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

**Chlordane:** The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

**Difference:** With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

**Final pH:** As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

**Frozen Date/Time:** With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**Initial pH:** As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

**PAH Total:** With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

**PFAS Total:** With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**Total:** With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

### Data Qualifiers

- A** - Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F** - The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The lower value for the two columns has been reported due to obvious interference.
- J** - Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

**Report Format:** DU Report with 'J' Qualifiers





**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25**Data Qualifiers**

Identified Compounds (TICs). For calculated parameters, this represents that one or more values used in the calculation were estimated.

- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND** - Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ** - Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- R** - Analytical results are from sample re-analysis.
- RE** - Analytical results are from sample re-extraction.
- S** - Analytical results are from modified screening analysis.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521696**Project Number:** 21010214**Report Date:** 04/18/25

## REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

## LIMITATION OF LIABILITIES

Pace Analytical Services performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Pace Analytical Services shall be to re-perform the work at it's own expense. In no event shall Pace Analytical Services be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Pace Analytical Services.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



**Pace Analytical Services LLC**Facility: **Northeast**Department: **Quality Assurance**Title: **Certificate/Approval Program Summary**ID No.: **17873**

Revision 27

Published Date: 01/24/2025

Page 1 of 2

**Certification Information****The following analytes are not included in our Primary NELAP Scope of Accreditation:****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****EPA 624.1:** m/p-xylene, o-xylene, Naphthalene**EPA 625.1:** alpha-Terpineol**EPA 8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****SM 2540D:** TSS.**EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**MADEP-APH.****Nonpotable Water:** EPA RSK-175 Dissolved Gases**Biological Tissue Matrix:** EPA 3050B**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048****EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**Nonpotable Water:** EPA RSK-175 Dissolved Gases**The following test method is not included in our New Jersey Secondary NELAP Scope of Accreditation:****Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Determination of Selected Perfluorinated Alkyl Substances by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry Isotope Dilution (via Alpha SOP 23528)****The following analytes are included in our Massachusetts DEP Scope of Accreditation****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****Drinking Water****EPA 300.0:** Chloride, Nitrate-N, Fluoride, Sulfate; **EPA 353.2:** Nitrate-N, Nitrite-N; **SM4500NO3-F:** Nitrate-N, Nitrite-N; **SM4500F-C, SM4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B****EPA 524.2:** THMs and VOCs; **EPA 504.1:** EDB, DBCP.**Microbiology:** SM9215B; SM9223-P/A, SM9223B-Colilert-QT, SM9222D.**Non-Potable Water****SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH:** Ammonia-N and Kjeldahl-N, **EPA 350.1:** Ammonia-N, **LCHAT 10-107-06-1-B:** Ammonia-N, **EPA 351.1, SM4500NO3-F, EPA 353.2:** Nitrate-N, **SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300:** Chloride, Sulfate, Nitrate.**EPA 624.1:** Volatile Halocarbons & Aromatics,**EPA 608.3:** Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs**EPA 625.1:** SVOC (Acid/Base/Neutral Extractables).**Microbiology:** SM9223B-Colilert-QT; Enterolert-QT, EPA 1600, EPA 1603, SM9222D.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Drinking Water****EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1 Hg. EPA 522, EPA 537.1.****Non-Potable Water****EPA 200.7:** Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.**EPA 200.8:** Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.**EPA 245.1 Hg.****SM2340B**

**Pace Analytical Services LLC**ID No.:**17873**Facility: **Northeast**

Revision 27

Department: **Quality Assurance**

Published Date: 01/24/2025

Title: **Certificate/Approval Program Summary**

Page 2 of 2

**Certification IDs:****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581**

CT PH-0826, IL 200077, IN C-MA-03, KY JY98045, ME MA00086, MD 348, MA M-MA086, NH 2064, NJ MA935, NY 11148, NC (DW) 25700, NC (NPW/SCM) 666, OR MA-1316, PA 68-03671, RI LAO00065, TX T104704476, VT VT-0935, VA 460195

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048**

CT PH-0825, ANAB/DoD L2474, IL 200081, IN C-MA-04, KY KY98046, LA 3090, ME MA00030, MI 9110, MN 025-999-495, NH 2062, NJ MA015, NY 11627, NC (NPW/SCM) 685, OR MA-0262, PA 68-02089, RI LAO00299, TX T-104704419, VT VT-0015, VA 460194, WA C954

**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048**

ANAB/DoD L2474, ME MA01156, MN 025-999-498, NH 2249, NJ MA025, NY 12191, OR 4203, TX T104704583, VA 460311, WA C1104.

---

For a complete listing of analytes and methods, please contact your Project Manager.







## ANALYTICAL REPORT

|                 |                                                                             |
|-----------------|-----------------------------------------------------------------------------|
| Lab Number:     | L2521646                                                                    |
| Client:         | Tradepoint Atlantic<br>1600 Sparrows Point Boulevard<br>Baltimore, MD 21219 |
| ATTN:           | Matthew Newman                                                              |
| Phone:          | (443) 649-5073                                                              |
| Project Name:   | B14 1ST HALF 2025 GW                                                        |
| Project Number: | 21010214                                                                    |
| Report Date:    | 04/18/25                                                                    |

The original project report/data package is held by Pace Analytical Services. This report/data package is paginated and should be reproduced only in its entirety. Pace Analytical Services holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0826), IL (200077), IN (C-MA-03), KY (KY98045), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), OR (MA-1316), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #525-23-122-91930A1).

---

Eight Walkup Drive, Westborough, MA 01581-1019  
508-898-9220 (Fax) 508-898-9193 800-624-9220 - [www.alphalab.com](http://www.alphalab.com)



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

| Lab<br>Sample ID | Client ID    | Matrix | Sample<br>Location | Collection<br>Date/Time | Receive Date |
|------------------|--------------|--------|--------------------|-------------------------|--------------|
| L2521646-01      | HI20R2-MWS   | WATER  | B14                | 04/09/25 11:00          | 04/09/25     |
| L2521646-02      | HI07R-PZM005 | WATER  | B14                | 04/09/25 11:55          | 04/09/25     |
| L2521646-03      | HI18-MWS     | WATER  | B14                | 04/09/25 13:30          | 04/09/25     |
| L2521646-04      | TM08R-PZM007 | WATER  | B14                | 04/09/25 14:20          | 04/09/25     |
| L2521646-05      | TB-001       | WATER  | B14                | 04/09/25 00:00          | 04/09/25     |
| L2521646-06      | TB-002       | WATER  | B14                | 04/09/25 00:00          | 04/09/25     |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

### Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Pace Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments and solids are reported on a dry weight basis unless otherwise noted. Tissues are reported "as received" or on a wet weight basis, unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

**HOLD POLICY** - For samples submitted on hold, Pace's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Pace Project Manager and made arrangements for Pace to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.

---

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

### Case Narrative (continued)

#### Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

#### Sample Receipt

L2521646-01 through -06: The collection date was specified by the client.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:  Tiffani Morrissey

Title: Technical Director/Representative

Date: 04/18/25

# ORGANICS



# VOLATILES

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-01  
**Client ID:** HI20R2-MWS  
**Sample Location:** B14

**Date Collected:** 04/09/25 11:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 09:12  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 2.8    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.32   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS****Lab ID:** L2521646-01**Date Collected:** 04/09/25 11:00**Client ID:** HI20R2-MWS**Date Received:** 04/09/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 115        |           | 70-130              |
| Toluene-d8            | 98         |           | 70-130              |
| 4-Bromofluorobenzene  | 98         |           | 70-130              |
| Dibromofluoromethane  | 105        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

Lab ID: L2521646-01  
 Client ID: HI20R2-MWS  
 Sample Location: B14

Date Collected: 04/09/25 11:00  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/17/25 09:12  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 114        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 94         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-02  
**Client ID:** HI07R-PZM005  
**Sample Location:** B14

**Date Collected:** 04/09/25 11:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 09:38  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 4.6    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | 0.26   | J         | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 2.3    |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | 1.5    |           | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521646-02  
 Client ID: HI07R-PZM005  
 Sample Location: B14

Date Collected: 04/09/25 11:55  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | 0.88   | J         | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | 0.88   | J         | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 116        |           | 70-130              |
| Toluene-d8            | 100        |           | 70-130              |
| 4-Bromofluorobenzene  | 100        |           | 70-130              |
| Dibromofluoromethane  | 106        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-02  
**Client ID:** HI07R-PZM005  
**Sample Location:** B14

**Date Collected:** 04/09/25 11:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/17/25 09:38  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 114        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-03  
**Client ID:** HI18-MWS  
**Sample Location:** B14

**Date Collected:** 04/09/25 13:30  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 10:05  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 5.3    |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS****Lab ID:** L2521646-03**Date Collected:** 04/09/25 13:30**Client ID:** HI18-MWS**Date Received:** 04/09/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 115        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 98         |           | 70-130              |
| Dibromofluoromethane  | 106        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

Lab ID: L2521646-03  
 Client ID: HI18-MWS  
 Sample Location: B14

Date Collected: 04/09/25 13:30  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/17/25 10:05  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 114        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 94         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-04  
**Client ID:** TM08R-PZM007  
**Sample Location:** B14

**Date Collected:** 04/09/25 14:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 10:31  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 3.3    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

**Lab ID:** L2521646-04  
**Client ID:** TM08R-PZM007  
**Sample Location:** B14

**Date Collected:** 04/09/25 14:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 117        |           | 70-130              |
| Toluene-d8            | 100        |           | 70-130              |
| 4-Bromofluorobenzene  | 99         |           | 70-130              |
| Dibromofluoromethane  | 106        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-04  
**Client ID:** TM08R-PZM007  
**Sample Location:** B14

**Date Collected:** 04/09/25 14:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/17/25 10:31  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 114        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 95         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-05  
**Client ID:** TB-001  
**Sample Location:** B14

**Date Collected:** 04/09/25 00:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 07:26  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | ND     |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS****Lab ID:** L2521646-05**Date Collected:** 04/09/25 00:00**Client ID:** TB-001**Date Received:** 04/09/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 118        |           | 70-130              |
| Toluene-d8            | 100        |           | 70-130              |
| 4-Bromofluorobenzene  | 99         |           | 70-130              |
| Dibromofluoromethane  | 106        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

Lab ID: L2521646-05  
 Client ID: TB-001  
 Sample Location: B14

Date Collected: 04/09/25 00:00  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/17/25 07:26  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 115        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 95         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-06  
**Client ID:** TB-002  
**Sample Location:** B14

**Date Collected:** 04/09/25 00:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 06:33  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | ND     |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



Project Name: B14 1ST HALF 2025 GW

Lab Number: L2521646

Project Number: 21010214

Report Date: 04/18/25

## SAMPLE RESULTS

Lab ID: L2521646-06

Date Collected: 04/09/25 00:00

Client ID: TB-002

Date Received: 04/09/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 116        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 100        |           | 70-130              |
| Dibromofluoromethane  | 107        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-06  
**Client ID:** TB-002  
**Sample Location:** B14

**Date Collected:** 04/09/25 00:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/17/25 06:33  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 114        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 96         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 04/17/25 06:07  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-06 Batch: WG2055118-5 |        |           |       |      |      |
| Dichlorodifluoromethane                                                              | ND     |           | ug/l  | 5.0  | 0.24 |
| Chloromethane                                                                        | ND     |           | ug/l  | 2.5  | 0.20 |
| Vinyl chloride                                                                       | ND     |           | ug/l  | 1.0  | 0.07 |
| Bromomethane                                                                         | ND     |           | ug/l  | 1.0  | 0.26 |
| Chloroethane                                                                         | ND     |           | ug/l  | 1.0  | 0.13 |
| Trichlorofluoromethane                                                               | ND     |           | ug/l  | 2.5  | 0.16 |
| 1,1-Dichloroethene                                                                   | ND     |           | ug/l  | 0.50 | 0.17 |
| Carbon disulfide                                                                     | ND     |           | ug/l  | 5.0  | 0.30 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                | ND     |           | ug/l  | 2.5  | 0.15 |
| Methylene chloride                                                                   | ND     |           | ug/l  | 2.5  | 0.68 |
| Acetone                                                                              | ND     |           | ug/l  | 5.0  | 1.5  |
| trans-1,2-Dichloroethene                                                             | ND     |           | ug/l  | 0.75 | 0.16 |
| Methyl Acetate                                                                       | ND     |           | ug/l  | 2.0  | 0.23 |
| Methyl tert butyl ether                                                              | ND     |           | ug/l  | 1.0  | 0.17 |
| 1,1-Dichloroethane                                                                   | ND     |           | ug/l  | 0.75 | 0.21 |
| cis-1,2-Dichloroethene                                                               | ND     |           | ug/l  | 0.50 | 0.19 |
| Cyclohexane                                                                          | ND     |           | ug/l  | 10   | 0.27 |
| Chloroform                                                                           | ND     |           | ug/l  | 0.75 | 0.22 |
| Carbon tetrachloride                                                                 | ND     |           | ug/l  | 0.50 | 0.13 |
| 1,1,1-Trichloroethane                                                                | ND     |           | ug/l  | 0.50 | 0.16 |
| 2-Butanone                                                                           | ND     |           | ug/l  | 5.0  | 1.9  |
| Benzene                                                                              | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,2-Dichloroethane                                                                   | ND     |           | ug/l  | 0.50 | 0.13 |
| Trichloroethene                                                                      | ND     |           | ug/l  | 0.50 | 0.18 |
| 1,2-Dichloropropane                                                                  | ND     |           | ug/l  | 1.0  | 0.14 |
| Bromodichloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.19 |
| cis-1,3-Dichloropropene                                                              | ND     |           | ug/l  | 0.50 | 0.14 |
| Toluene                                                                              | ND     |           | ug/l  | 0.75 | 0.20 |
| Tetrachloroethene                                                                    | ND     |           | ug/l  | 0.50 | 0.18 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 06:07  
**Analyst:** MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-06 Batch: WG2055118-5 |        |           |       |      |      |
| 4-Methyl-2-pentanone                                                                 | ND     |           | ug/l  | 5.0  | 0.42 |
| trans-1,3-Dichloropropene                                                            | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,3-Dichloropropene, Total                                                           | ND     |           | ug/l  | 0.50 | 0.14 |
| 1,1,2-Trichloroethane                                                                | ND     |           | ug/l  | 0.75 | 0.14 |
| Dibromochloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.15 |
| 1,2-Dibromoethane                                                                    | ND     |           | ug/l  | 2.0  | 0.19 |
| 2-Hexanone                                                                           | ND     |           | ug/l  | 5.0  | 0.52 |
| Chlorobenzene                                                                        | ND     |           | ug/l  | 0.50 | 0.18 |
| Ethylbenzene                                                                         | ND     |           | ug/l  | 0.50 | 0.17 |
| p/m-Xylene                                                                           | ND     |           | ug/l  | 1.0  | 0.33 |
| o-Xylene                                                                             | ND     |           | ug/l  | 1.0  | 0.39 |
| Xylenes, Total                                                                       | ND     |           | ug/l  | 1.0  | 0.33 |
| Styrene                                                                              | ND     |           | ug/l  | 1.0  | 0.36 |
| Bromoform                                                                            | ND     |           | ug/l  | 2.0  | 0.25 |
| Isopropylbenzene                                                                     | ND     |           | ug/l  | 0.50 | 0.19 |
| 1,1,2,2-Tetrachloroethane                                                            | ND     |           | ug/l  | 0.50 | 0.17 |
| 1,3-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,4-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,2-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.18 |
| 1,2-Dibromo-3-chloropropane                                                          | ND     |           | ug/l  | 2.5  | 0.35 |
| 1,2,4-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.22 |
| 1,2,3-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.23 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
Analytical Date: 04/17/25 06:07  
Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL | MDL |
|--------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-06 Batch: WG2055118-5 |        |           |       |    |     |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 119       |           | 70-130                 |
| Toluene-d8            | 99        |           | 70-130                 |
| 4-Bromofluorobenzene  | 98        |           | 70-130                 |
| Dibromofluoromethane  | 106       |           | 70-130                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D-SIM(M)  
Analytical Date: 04/17/25 06:07  
Analyst: MCM

| Parameter                                                                                | Result | Qualifier | Units | RL    | MDL   |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-------|-------|
| Volatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-06 Batch: WG2055119-5 |        |           |       |       |       |
| 1,1,1,2-Tetrachloroethane                                                                | ND     |           | ug/l  | 0.050 | 0.006 |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 115       |           | 70-130                 |
| 4-Bromofluorobenzene  | 96        |           | 70-130                 |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521646

**Report Date:** 04/18/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-06 Batch: WG2055118-3 WG2055118-4 |                  |      |                   |      |                     |     |      |               |
| Dichlorodifluoromethane                                                                                 | 110              |      | 100               |      | 36-147              | 10  |      | 20            |
| Chloromethane                                                                                           | 97               |      | 95                |      | 64-130              | 2   |      | 20            |
| Vinyl chloride                                                                                          | 110              |      | 100               |      | 55-140              | 10  |      | 20            |
| Bromomethane                                                                                            | 64               |      | 66                |      | 39-139              | 3   |      | 20            |
| Chloroethane                                                                                            | 120              |      | 120               |      | 55-138              | 0   |      | 20            |
| Trichlorofluoromethane                                                                                  | 110              |      | 110               |      | 62-150              | 0   |      | 20            |
| 1,1-Dichloroethene                                                                                      | 100              |      | 98                |      | 61-145              | 2   |      | 20            |
| Carbon disulfide                                                                                        | 110              |      | 100               |      | 51-130              | 10  |      | 20            |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                                   | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Methylene chloride                                                                                      | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| Acetone                                                                                                 | 120              |      | 110               |      | 58-148              | 9   |      | 20            |
| trans-1,2-Dichloroethene                                                                                | 100              |      | 97                |      | 70-130              | 3   |      | 20            |
| Methyl Acetate                                                                                          | 120              |      | 120               |      | 70-130              | 0   |      | 20            |
| Methyl tert butyl ether                                                                                 | 100              |      | 99                |      | 63-130              | 1   |      | 20            |
| 1,1-Dichloroethane                                                                                      | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| cis-1,2-Dichloroethene                                                                                  | 98               |      | 96                |      | 70-130              | 2   |      | 20            |
| Cyclohexane                                                                                             | 120              |      | 110               |      | 70-130              | 9   |      | 20            |
| Chloroform                                                                                              | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Carbon tetrachloride                                                                                    | 100              |      | 100               |      | 63-132              | 0   |      | 20            |
| 1,1,1-Trichloroethane                                                                                   | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| 2-Butanone                                                                                              | 96               |      | 93                |      | 63-138              | 3   |      | 20            |
| Benzene                                                                                                 | 100              |      | 96                |      | 70-130              | 4   |      | 20            |
| 1,2-Dichloroethane                                                                                      | 110              |      | 110               |      | 70-130              | 0   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521646

**Report Date:** 04/18/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-06 Batch: WG2055118-3 WG2055118-4 |                  |      |                   |      |                     |     |      |               |
| Trichloroethene                                                                                         | 98               |      | 94                |      | 70-130              | 4   |      | 20            |
| 1,2-Dichloropropane                                                                                     | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Bromodichloromethane                                                                                    | 99               |      | 95                |      | 67-130              | 4   |      | 20            |
| cis-1,3-Dichloropropene                                                                                 | 94               |      | 92                |      | 70-130              | 2   |      | 20            |
| Toluene                                                                                                 | 98               |      | 96                |      | 70-130              | 2   |      | 20            |
| Tetrachloroethene                                                                                       | 99               |      | 95                |      | 70-130              | 4   |      | 20            |
| 4-Methyl-2-pentanone                                                                                    | 89               |      | 88                |      | 59-130              | 1   |      | 20            |
| trans-1,3-Dichloropropene                                                                               | 99               |      | 99                |      | 70-130              | 0   |      | 20            |
| 1,1,2-Trichloroethane                                                                                   | 100              |      | 98                |      | 70-130              | 2   |      | 20            |
| Dibromochloromethane                                                                                    | 92               |      | 92                |      | 63-130              | 0   |      | 20            |
| 1,2-Dibromoethane                                                                                       | 96               |      | 94                |      | 70-130              | 2   |      | 20            |
| 2-Hexanone                                                                                              | 95               |      | 91                |      | 57-130              | 4   |      | 20            |
| Chlorobenzene                                                                                           | 96               |      | 94                |      | 75-130              | 2   |      | 20            |
| Ethylbenzene                                                                                            | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| p/m-Xylene                                                                                              | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| o-Xylene                                                                                                | 100              |      | 95                |      | 70-130              | 5   |      | 20            |
| Styrene                                                                                                 | 95               |      | 95                |      | 70-130              | 0   |      | 20            |
| Bromoform                                                                                               | 89               |      | 88                |      | 54-136              | 1   |      | 20            |
| Isopropylbenzene                                                                                        | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| 1,1,2,2-Tetrachloroethane                                                                               | 100              |      | 96                |      | 67-130              | 4   |      | 20            |
| 1,3-Dichlorobenzene                                                                                     | 100              |      | 97                |      | 70-130              | 3   |      | 20            |
| 1,4-Dichlorobenzene                                                                                     | 99               |      | 95                |      | 70-130              | 4   |      | 20            |
| 1,2-Dichlorobenzene                                                                                     | 98               |      | 96                |      | 70-130              | 2   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521646

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b>                                                                                        | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|---------------------------------------------------------------------------------------------------------|--------------------------|-------------|--------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-06 Batch: WG2055118-3 WG2055118-4 |                          |             |                          |             |                             |            |             |                       |
| 1,2-Dibromo-3-chloropropane                                                                             | 82                       |             | 80                       |             | 41-144                      | 2          |             | 20                    |
| 1,2,4-Trichlorobenzene                                                                                  | 94                       |             | 89                       |             | 70-130                      | 5          |             | 20                    |
| 1,2,3-Trichlorobenzene                                                                                  | 92                       |             | 91                       |             | 70-130                      | 1          |             | 20                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|--------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 115                      |             | 112                      |             | 70-130                         |
| Toluene-d8            | 101                      |             | 100                      |             | 70-130                         |
| 4-Bromofluorobenzene  | 102                      |             | 100                      |             | 70-130                         |
| Dibromofluoromethane  | 103                      |             | 103                      |             | 70-130                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521646

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-06 Batch: WG2055119-3 WG2055119-4 |                          |             |                           |             |                             |            |             |                       |
| 1,1,1,2-Tetrachloroethane                                                                                   | 100                      |             | 96                        |             | 70-130                      | 4          |             | 25                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 114                      |             | 115                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 98                       |             | 97                        |             | 70-130                         |

# SEMIVOLATILES

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-01  
**Client ID:** HI20R2-MWS  
**Sample Location:** B14

**Date Collected:** 04/09/25 11:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 00:21  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | 4.8    | J         | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | 8.1    |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521646-01

Date Collected: 04/09/25 11:00

Client ID: HI20R2-MWS

Date Received: 04/09/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 49         |           | 21-120              |
| Phenol-d6            | 38         |           | 10-120              |
| Nitrobenzene-d5      | 68         |           | 23-120              |
| 2-Fluorobiphenyl     | 74         |           | 15-120              |
| 2,4,6-Tribromophenol | 76         |           | 10-120              |
| 4-Terphenyl-d14      | 78         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-01  
**Client ID:** HI20R2-MWS  
**Sample Location:** B14

**Date Collected:** 04/09/25 11:00  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 08:54  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.44   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.12   |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.17   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.19   |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.29   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.05   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.13   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.13   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 53         |           | 21-120              |
| Phenol-d6            | 43         |           | 10-120              |
| Nitrobenzene-d5      | 95         |           | 23-120              |
| 2-Fluorobiphenyl     | 82         |           | 15-120              |
| 2,4,6-Tribromophenol | 117        |           | 10-120              |
| 4-Terphenyl-d14      | 74         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-02  
**Client ID:** HI07R-PZM005  
**Sample Location:** B14

**Date Collected:** 04/09/25 11:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 00:44  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | 3.6    | J         | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 5.2    |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | 0.73   | J         | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | 0.26   | J         | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521646-02  
 Client ID: HI07R-PZM005  
 Sample Location: B14

Date Collected: 04/09/25 11:55  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | 0.51   | J         | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | 0.91   | J         | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | 8.0    | J         | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | 1.5    | J         | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | 0.92   | J         | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | 1.0    | J         | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | 0.59   | J         | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 49         |           | 21-120              |
| Phenol-d6            | 37         |           | 10-120              |
| Nitrobenzene-d5      | 69         |           | 23-120              |
| 2-Fluorobiphenyl     | 75         |           | 15-120              |
| 2,4,6-Tribromophenol | 78         |           | 10-120              |
| 4-Terphenyl-d14      | 89         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-02  
**Client ID:** HI07R-PZM005  
**Sample Location:** B14

**Date Collected:** 04/09/25 11:55  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 09:45  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 5.4    |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.85   |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.36   |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.55   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.76   |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 1.5    |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.28   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.89   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.56   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | 0.05   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 54         |           | 21-120              |
| Phenol-d6            | 43         |           | 10-120              |
| Nitrobenzene-d5      | 98         |           | 23-120              |
| 2-Fluorobiphenyl     | 83         |           | 15-120              |
| 2,4,6-Tribromophenol | 122        | Q         | 10-120              |
| 4-Terphenyl-d14      | 75         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-03  
**Client ID:** HI18-MWS  
**Sample Location:** B14

**Date Collected:** 04/09/25 13:30  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 01:08  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | 3.3    | J         | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 4.4    |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | 0.64   | J         | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521646-03

Date Collected: 04/09/25 13:30

Client ID: HI18-MWS

Date Received: 04/09/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | 0.47   | J         | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | 0.81   | J         | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | 6.3    | J         | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | 1.3    | J         | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | 0.81   | J         | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | 0.80   | J         | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | 0.49   | J         | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 46         |           | 21-120              |
| Phenol-d6            | 33         |           | 10-120              |
| Nitrobenzene-d5      | 63         |           | 23-120              |
| 2-Fluorobiphenyl     | 67         |           | 15-120              |
| 2,4,6-Tribromophenol | 73         |           | 10-120              |
| 4-Terphenyl-d14      | 73         |           | 41-149              |





**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-03  
**Client ID:** HI18-MWS  
**Sample Location:** B14

**Date Collected:** 04/09/25 13:30  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 12:18  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.27   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.05   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.06   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.07   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.26   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.05   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 50         |           | 21-120              |
| Phenol-d6            | 41         |           | 10-120              |
| Nitrobenzene-d5      | 88         |           | 23-120              |
| 2-Fluorobiphenyl     | 75         |           | 15-120              |
| 2,4,6-Tribromophenol | 105        |           | 10-120              |
| 4-Terphenyl-d14      | 61         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-04  
**Client ID:** TM08R-PZM007  
**Sample Location:** B14

**Date Collected:** 04/09/25 14:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 01:31  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**SAMPLE RESULTS**

Lab ID: L2521646-04  
 Client ID: TM08R-PZM007  
 Sample Location: B14

Date Collected: 04/09/25 14:20  
 Date Received: 04/09/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 48         |           | 21-120              |
| Phenol-d6            | 38         |           | 10-120              |
| Nitrobenzene-d5      | 65         |           | 23-120              |
| 2-Fluorobiphenyl     | 68         |           | 15-120              |
| 2,4,6-Tribromophenol | 73         |           | 10-120              |
| 4-Terphenyl-d14      | 73         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**SAMPLE RESULTS**

**Lab ID:** L2521646-04  
**Client ID:** TM08R-PZM007  
**Sample Location:** B14

**Date Collected:** 04/09/25 14:20  
**Date Received:** 04/09/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 10:19  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.08   | J         | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.03   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.07   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.03   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 54         |           | 21-120              |
| Phenol-d6            | 44         |           | 10-120              |
| Nitrobenzene-d5      | 92         |           | 23-120              |
| 2-Fluorobiphenyl     | 79         |           | 15-120              |
| 2,4,6-Tribromophenol | 114        |           | 10-120              |
| 4-Terphenyl-d14      | 70         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 20:08  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2053156-1 |        |           |       |     |      |
| Benzaldehyde                                                                             | ND     |           | ug/l  | 5.0 | 1.1  |
| Phenol                                                                                   | ND     |           | ug/l  | 5.0 | 0.35 |
| Bis(2-chloroethyl)ether                                                                  | ND     |           | ug/l  | 2.0 | 0.39 |
| 2-Chlorophenol                                                                           | ND     |           | ug/l  | 2.0 | 0.65 |
| 2-Methylphenol                                                                           | ND     |           | ug/l  | 5.0 | 2.3  |
| Bis(2-chloroisopropyl)ether                                                              | ND     |           | ug/l  | 2.0 | 0.40 |
| Acetophenone                                                                             | ND     |           | ug/l  | 5.0 | 0.92 |
| n-Nitrosodi-n-propylamine                                                                | ND     |           | ug/l  | 5.0 | 0.91 |
| 3-Methylphenol/4-Methylphenol                                                            | ND     |           | ug/l  | 5.0 | 1.4  |
| Hexachloroethane                                                                         | ND     |           | ug/l  | 2.0 | 0.20 |
| Nitrobenzene                                                                             | ND     |           | ug/l  | 1.4 | 0.20 |
| Isophorone                                                                               | ND     |           | ug/l  | 5.0 | 0.86 |
| 2,4-Dimethylphenol                                                                       | ND     |           | ug/l  | 5.0 | 2.0  |
| Bis(2-chloroethoxy)methane                                                               | ND     |           | ug/l  | 5.0 | 0.84 |
| 2,4-Dichlorophenol                                                                       | ND     |           | ug/l  | 5.0 | 1.7  |
| Naphthalene                                                                              | ND     |           | ug/l  | 2.0 | 0.54 |
| 4-Chloroaniline                                                                          | ND     |           | ug/l  | 3.7 | 0.47 |
| Hexachlorobutadiene                                                                      | ND     |           | ug/l  | 2.0 | 0.36 |
| Caprolactam                                                                              | ND     |           | ug/l  | 10  | 1.2  |
| 2-Methylnaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.37 |
| Hexachlorocyclopentadiene                                                                | ND     |           | ug/l  | 20  | 1.2  |
| 1,2,4,5-Tetrachlorobenzene                                                               | ND     |           | ug/l  | 1.7 | 0.24 |
| 2,4,6-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| 2,4,5-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| Biphenyl                                                                                 | ND     |           | ug/l  | 2.0 | 0.20 |
| 2-Chloronaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.35 |
| 2-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.0  |
| 2,6-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.84 |
| Acenaphthylene                                                                           | ND     |           | ug/l  | 2.0 | 0.32 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 20:08  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2053156-1 |        |           |       |     |      |
| Acenaphthene                                                                             | ND     |           | ug/l  | 2.0 | 0.40 |
| 2,4-Dinitrophenol                                                                        | ND     |           | ug/l  | 20  | 5.4  |
| 2,4-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.54 |
| 2,3,4,6-Tetrachlorophenol                                                                | ND     |           | ug/l  | 5.0 | 2.2  |
| Diethyl phthalate                                                                        | ND     |           | ug/l  | 5.0 | 0.76 |
| Fluorene                                                                                 | ND     |           | ug/l  | 2.0 | 0.44 |
| 4-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.4  |
| NDPA/DPA                                                                                 | ND     |           | ug/l  | 2.0 | 0.92 |
| Hexachlorobenzene                                                                        | ND     |           | ug/l  | 2.0 | 0.45 |
| Pentachlorophenol                                                                        | ND     |           | ug/l  | 10  | 2.5  |
| Phenanthrene                                                                             | ND     |           | ug/l  | 2.0 | 0.42 |
| Anthracene                                                                               | ND     |           | ug/l  | 2.0 | 0.47 |
| Carbazole                                                                                | ND     |           | ug/l  | 2.0 | 0.31 |
| Di-n-butylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 0.96 |
| Fluoranthene                                                                             | ND     |           | ug/l  | 2.0 | 0.41 |
| Pyrene                                                                                   | ND     |           | ug/l  | 2.0 | 0.41 |
| 3,3'-Dichlorobenzidine                                                                   | ND     |           | ug/l  | 5.0 | 1.8  |
| Benzo(a)anthracene                                                                       | ND     |           | ug/l  | 2.0 | 0.32 |
| Chrysene                                                                                 | ND     |           | ug/l  | 1.5 | 0.22 |
| Bis(2-ethylhexyl)phthalate                                                               | ND     |           | ug/l  | 3.0 | 1.4  |
| Di-n-octylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 2.3  |
| Benzo(b)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.53 |
| Benzo(k)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.62 |
| Benzo(a)pyrene                                                                           | ND     |           | ug/l  | 2.0 | 0.37 |
| Indeno(1,2,3-cd)pyrene                                                                   | ND     |           | ug/l  | 2.0 | 0.48 |
| Dibenzo(a,h)anthracene                                                                   | ND     |           | ug/l  | 2.0 | 0.29 |
| Benzo(ghi)perylene                                                                       | ND     |           | ug/l  | 2.0 | 0.37 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8270E  
 Analytical Date: 04/16/25 20:08  
 Analyst: SLR

Extraction Method: EPA 3510C  
 Extraction Date: 04/13/25 14:21

| Parameter                                                                                | Result | Qualifier | Units | RL | MDL |
|------------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2053156-1 |        |           |       |    |     |

| Surrogate            | %Recovery | Qualifier | Acceptance<br>Criteria |
|----------------------|-----------|-----------|------------------------|
| 2-Fluorophenol       | 53        |           | 21-120                 |
| Phenol-d6            | 37        |           | 10-120                 |
| Nitrobenzene-d5      | 68        |           | 23-120                 |
| 2-Fluorobiphenyl     | 72        |           | 15-120                 |
| 2,4,6-Tribromophenol | 67        |           | 10-120                 |
| 4-Terphenyl-d14      | 77        |           | 41-149                 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/14/25 06:54  
**Analyst:** KMH

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/13/25 14:21

| Parameter                                                                                    | Result | Qualifier | Units | RL   | MDL  |
|----------------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-04 Batch: WG2053157-1 |        |           |       |      |      |
| Naphthalene                                                                                  | ND     |           | ug/l  | 0.10 | 0.02 |
| 2-Methylnaphthalene                                                                          | ND     |           | ug/l  | 0.10 | 0.03 |
| Acenaphthylene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Acenaphthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluorene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Phenanthrene                                                                                 | ND     |           | ug/l  | 0.05 | 0.04 |
| Anthracene                                                                                   | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluoranthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.03 |
| Pyrene                                                                                       | ND     |           | ug/l  | 0.10 | 0.04 |
| Benzo(a)anthracene                                                                           | ND     |           | ug/l  | 0.05 | 0.03 |
| Chrysene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(b)fluoranthene                                                                         | ND     |           | ug/l  | 0.05 | 0.03 |
| Benzo(k)fluoranthene                                                                         | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(a)pyrene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Indeno(1,2,3-cd)pyrene                                                                       | ND     |           | ug/l  | 0.10 | 0.02 |
| Dibenzo(a,h)anthracene                                                                       | ND     |           | ug/l  | 0.05 | 0.02 |
| Benzo(ghi)perylene                                                                           | ND     |           | ug/l  | 0.10 | 0.02 |

| Surrogate            | %Recovery  | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 76         |           | 21-120              |
| Phenol-d6            | 60         |           | 10-120              |
| Nitrobenzene-d5      | <b>121</b> | Q         | 23-120              |
| 2-Fluorobiphenyl     | 106        |           | 15-120              |
| 2,4,6-Tribromophenol | <b>140</b> | Q         | 10-120              |
| 4-Terphenyl-d14      | 109        |           | 41-149              |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521646

**Report Date:** 04/18/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2053156-2 WG2053156-3 |                  |      |                   |      |                     |     |      |               |
| Benzaldehyde                                                                                                | 78               |      | 77                |      | 40-140              | 1   |      | 30            |
| Phenol                                                                                                      | 41               |      | 40                |      | 12-110              | 2   |      | 30            |
| Bis(2-chloroethyl)ether                                                                                     | 71               |      | 74                |      | 40-140              | 4   |      | 30            |
| 2-Chlorophenol                                                                                              | 70               |      | 69                |      | 27-123              | 1   |      | 30            |
| 2-Methylphenol                                                                                              | 70               |      | 70                |      | 30-130              | 0   |      | 30            |
| Bis(2-chloroisopropyl)ether                                                                                 | 77               |      | 77                |      | 40-140              | 0   |      | 30            |
| Acetophenone                                                                                                | 80               |      | 79                |      | 39-129              | 1   |      | 30            |
| n-Nitrosodi-n-propylamine                                                                                   | 71               |      | 69                |      | 29-132              | 3   |      | 30            |
| 3-Methylphenol/4-Methylphenol                                                                               | 67               |      | 72                |      | 30-130              | 7   |      | 30            |
| Hexachloroethane                                                                                            | 63               |      | 65                |      | 40-140              | 3   |      | 30            |
| Nitrobenzene                                                                                                | 68               |      | 69                |      | 40-140              | 1   |      | 30            |
| Isophorone                                                                                                  | 68               |      | 69                |      | 40-140              | 1   |      | 30            |
| 2,4-Dimethylphenol                                                                                          | 64               |      | 59                |      | 30-130              | 8   |      | 30            |
| Bis(2-chloroethoxy)methane                                                                                  | 72               |      | 73                |      | 40-140              | 1   |      | 30            |
| 2,4-Dichlorophenol                                                                                          | 81               |      | 80                |      | 30-130              | 1   |      | 30            |
| Naphthalene                                                                                                 | 71               |      | 72                |      | 40-140              | 1   |      | 30            |
| 4-Chloroaniline                                                                                             | 48               |      | 57                |      | 40-140              | 17  |      | 30            |
| Hexachlorobutadiene                                                                                         | 64               |      | 68                |      | 40-140              | 6   |      | 30            |
| Caprolactam                                                                                                 | 40               |      | 44                |      | 10-130              | 10  |      | 30            |
| 2-Methylnaphthalene                                                                                         | 74               |      | 75                |      | 40-140              | 1   |      | 30            |
| Hexachlorocyclopentadiene                                                                                   | 67               |      | 70                |      | 40-140              | 4   |      | 30            |
| 1,2,4,5-Tetrachlorobenzene                                                                                  | 73               |      | 78                |      | 2-134               | 7   |      | 30            |
| 2,4,6-Trichlorophenol                                                                                       | 80               |      | 81                |      | 30-130              | 1   |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521646

**Project Number:** 21010214

**Report Date:** 04/18/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2053156-2 WG2053156-3 |                  |      |                   |      |                     |     |      |               |
| 2,4,5-Trichlorophenol                                                                                       | 88               |      | 92                |      | 30-130              | 4   |      | 30            |
| Biphenyl                                                                                                    | 81               |      | 82                |      | 40-140              | 1   |      | 30            |
| 2-Chloronaphthalene                                                                                         | 75               |      | 76                |      | 40-140              | 1   |      | 30            |
| 2-Nitroaniline                                                                                              | 86               |      | 89                |      | 52-143              | 3   |      | 30            |
| 2,6-Dinitrotoluene                                                                                          | 85               |      | 89                |      | 40-140              | 5   |      | 30            |
| Acenaphthylene                                                                                              | 75               |      | 78                |      | 45-123              | 4   |      | 30            |
| Acenaphthene                                                                                                | 76               |      | 77                |      | 37-111              | 1   |      | 30            |
| 2,4-Dinitrophenol                                                                                           | 81               |      | 80                |      | 20-130              | 1   |      | 30            |
| 2,4-Dinitrotoluene                                                                                          | 82               |      | 84                |      | 48-143              | 2   |      | 30            |
| 2,3,4,6-Tetrachlorophenol                                                                                   | 88               |      | 87                |      | 54-145              | 1   |      | 30            |
| Diethyl phthalate                                                                                           | 78               |      | 78                |      | 40-140              | 0   |      | 30            |
| Fluorene                                                                                                    | 77               |      | 78                |      | 40-140              | 1   |      | 30            |
| 4-Nitroaniline                                                                                              | 76               |      | 78                |      | 51-143              | 3   |      | 30            |
| NDPA/DPA                                                                                                    | 81               |      | 82                |      | 40-140              | 1   |      | 30            |
| Hexachlorobenzene                                                                                           | 81               |      | 82                |      | 40-140              | 1   |      | 30            |
| Pentachlorophenol                                                                                           | 90               |      | 85                |      | 9-103               | 6   |      | 30            |
| Phenanthrene                                                                                                | 79               |      | 82                |      | 40-140              | 4   |      | 30            |
| Anthracene                                                                                                  | 82               |      | 85                |      | 40-140              | 4   |      | 30            |
| Carbazole                                                                                                   | 81               |      | 83                |      | 55-144              | 2   |      | 30            |
| Di-n-butylphthalate                                                                                         | 83               |      | 85                |      | 40-140              | 2   |      | 30            |
| Fluoranthene                                                                                                | 84               |      | 86                |      | 40-140              | 2   |      | 30            |
| Pyrene                                                                                                      | 86               |      | 88                |      | 26-127              | 2   |      | 30            |
| 3,3'-Dichlorobenzidine                                                                                      | 64               |      | 72                |      | 40-140              | 12  |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521646

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2053156-2 WG2053156-3 |                          |             |                           |             |                             |            |             |                       |
| Benzo(a)anthracene                                                                                          | 80                       |             | 86                        |             | 40-140                      | 7          |             | 30                    |
| Chrysene                                                                                                    | 82                       |             | 86                        |             | 40-140                      | 5          |             | 30                    |
| Bis(2-ethylhexyl)phthalate                                                                                  | 88                       |             | 90                        |             | 40-140                      | 2          |             | 30                    |
| Di-n-octylphthalate                                                                                         | 85                       |             | 89                        |             | 40-140                      | 5          |             | 30                    |
| Benzo(b)fluoranthene                                                                                        | 84                       |             | 90                        |             | 40-140                      | 7          |             | 30                    |
| Benzo(k)fluoranthene                                                                                        | 87                       |             | 92                        |             | 40-140                      | 6          |             | 30                    |
| Benzo(a)pyrene                                                                                              | 87                       |             | 91                        |             | 40-140                      | 4          |             | 30                    |
| Indeno(1,2,3-cd)pyrene                                                                                      | 84                       |             | 89                        |             | 40-140                      | 6          |             | 30                    |
| Dibenzo(a,h)anthracene                                                                                      | 82                       |             | 87                        |             | 40-140                      | 6          |             | 30                    |
| Benzo(ghi)perylene                                                                                          | 80                       |             | 82                        |             | 40-140                      | 2          |             | 30                    |

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 54                       |             | 55                        |             | 21-120                         |
| Phenol-d6            | 42                       |             | 42                        |             | 10-120                         |
| Nitrobenzene-d5      | 67                       |             | 68                        |             | 23-120                         |
| 2-Fluorobiphenyl     | 74                       |             | 73                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 74                       |             | 77                        |             | 10-120                         |
| 4-Terphenyl-d14      | 76                       |             | 80                        |             | 41-149                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2521646

**Report Date:** 04/18/25

| Parameter                                                                                                       | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-----------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2053157-2 WG2053157-3 |                  |      |                   |      |                     |     |      |               |
| Naphthalene                                                                                                     | 69               |      | 75                |      | 40-140              | 8   |      | 40            |
| 2-Methylnaphthalene                                                                                             | 72               |      | 78                |      | 40-140              | 8   |      | 40            |
| Acenaphthylene                                                                                                  | 86               |      | 93                |      | 40-140              | 8   |      | 40            |
| Acenaphthene                                                                                                    | 75               |      | 82                |      | 37-111              | 9   |      | 40            |
| Fluorene                                                                                                        | 82               |      | 88                |      | 40-140              | 7   |      | 40            |
| Phenanthrene                                                                                                    | 78               |      | 85                |      | 40-140              | 9   |      | 40            |
| Anthracene                                                                                                      | 83               |      | 90                |      | 40-140              | 8   |      | 40            |
| Fluoranthene                                                                                                    | 82               |      | 88                |      | 40-140              | 7   |      | 40            |
| Pyrene                                                                                                          | 80               |      | 87                |      | 26-127              | 8   |      | 40            |
| Benzo(a)anthracene                                                                                              | 80               |      | 87                |      | 40-140              | 8   |      | 40            |
| Chrysene                                                                                                        | 83               |      | 90                |      | 40-140              | 8   |      | 40            |
| Benzo(b)fluoranthene                                                                                            | 87               |      | 94                |      | 40-140              | 8   |      | 40            |
| Benzo(k)fluoranthene                                                                                            | 90               |      | 99                |      | 40-140              | 10  |      | 40            |
| Benzo(a)pyrene                                                                                                  | 94               |      | 103               |      | 40-140              | 9   |      | 40            |
| Indeno(1,2,3-cd)pyrene                                                                                          | 104              |      | 114               |      | 40-140              | 9   |      | 40            |
| Dibenzo(a,h)anthracene                                                                                          | 103              |      | 112               |      | 40-140              | 8   |      | 40            |
| Benzo(ghi)perylene                                                                                              | 97               |      | 105               |      | 40-140              | 8   |      | 40            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2521646

**Project Number:** 21010214

**Report Date:** 04/18/25

| <b>Parameter</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
|------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2053157-2 WG2053157-3

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 58                       |             | 61                        |             | 21-120                         |
| Phenol-d6            | 48                       |             | 50                        |             | 10-120                         |
| Nitrobenzene-d5      | 87                       |             | 93                        |             | 23-120                         |
| 2-Fluorobiphenyl     | 76                       |             | 81                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 106                      |             | 115                       |             | 10-120                         |
| 4-Terphenyl-d14      | 71                       |             | 78                        |             | 41-149                         |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2521646**Project Number:** 21010214**Report Date:** 04/18/25**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

**Cooler Information**

| <b>Cooler</b> | <b>Custody Seal</b> |
|---------------|---------------------|
| A             | Absent              |
| B             | Absent              |

**Container Information**

| <b>Container ID</b> | <b>Container Type</b>   | <b>Cooler</b> | <b>Initial pH</b> | <b>Final pH</b> | <b>Temp deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen Date/Time</b> | <b>Analysis(*)</b>               |
|---------------------|-------------------------|---------------|-------------------|-----------------|-------------------|-------------|-------------|-------------------------|----------------------------------|
| L2521646-01A        | Vial HCl preserved      | B             | NA                |                 | 2.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-01B        | Vial HCl preserved      | B             | NA                |                 | 2.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-01C        | Vial HCl preserved      | B             | NA                |                 | 2.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-01D        | Amber 100ml unpreserved | B             | 7                 | 7               | 2.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521646-01E        | Amber 100ml unpreserved | B             | 7                 | 7               | 2.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521646-02A        | Vial HCl preserved      | B             | NA                |                 | 2.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-02B        | Vial HCl preserved      | B             | NA                |                 | 2.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-02C        | Vial HCl preserved      | B             | NA                |                 | 2.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-02D        | Amber 100ml unpreserved | B             | 11                | 11              | 2.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2521646-02E        | Amber 100ml unpreserved | B             | 11                | 11              | 2.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2521646-03A        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-03B        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-03C        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-03D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.7               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521646-03E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.7               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521646-04A        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-04B        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-04C        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-04D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.7               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521646-04E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.7               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2521646-05A        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2521646-05B        | Vial HCl preserved      | A             | NA                |                 | 3.7               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

Serial\_No:04182514:18  
**Lab Number:** L2521646  
**Report Date:** 04/18/25

**Container Information**

| <b>Container ID</b> | <b>Container Type</b> | <b>Cooler</b> | <b>Initial<br/>pH</b> | <b>Final<br/>pH</b> | <b>Temp<br/>deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen<br/>Date/Time</b> | <b>Analysis(*)</b>          |
|---------------------|-----------------------|---------------|-----------------------|---------------------|-----------------------|-------------|-------------|-----------------------------|-----------------------------|
| L2521646-05C        | Vial HCl preserved    | A             | NA                    |                     | 3.7                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |
| L2521646-05D        | Vial HCl preserved    | A             | NA                    |                     | 3.7                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |
| L2521646-06A        | Vial HCl preserved    | B             | NA                    |                     | 2.8                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |
| L2521646-06B        | Vial HCl preserved    | B             | NA                    |                     | 2.8                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |
| L2521646-06C        | Vial HCl preserved    | B             | NA                    |                     | 2.8                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |
| L2521646-06D        | Vial HCl preserved    | B             | NA                    |                     | 2.8                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

## GLOSSARY

### Acronyms

|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| DL       | - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                |
| EDL      | - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).                                                                                         |
| EMPC     | - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.                                                                                                                                                                                                                               |
| EPA      | - Environmental Protection Agency.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| LCS      | - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                          |
| LCSD     | - Laboratory Control Sample Duplicate: Refer to LCS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |
| LFB      | - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                         |
| LOD      | - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                                                               |
| LOQ      | - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)<br><br>Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| MDL      | - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                          |
| MS       | - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.                                                                                                                                                                                                    |
| MSD      | - Matrix Spike Sample Duplicate: Refer to MS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NA       | - Not Applicable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| NC       | - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.                                                                                                                                                                                                                                                                                                                                                                           |
| NDPA/DPA | - N-Nitrosodiphenylamine/Diphenylamine.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| NI       | - Not Ignitable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| NP       | - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NR       | - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.                                                                                                                                                                                                                                                                                                                                                                    |
| RL       | - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                                                                                                                                   |
| RPD      | - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.                                                                  |
| SRM      | - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.                                                                                                                                                                                                                                                                                                                                                                     |
| STLP     | - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| TEF      | - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.                                                                                                                                                                                                                                                                                                                                                                                             |
| TEQ      | - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.                                                                                                                                                                                                                                                                                                                                                        |
| TIC      | - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.                                                                                                                                                                                                                                                                      |

Report Format: DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

### Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

### Terms

**Analytical Method:** Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

**Chlordane:** The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

**Difference:** With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

**Final pH:** As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

**Frozen Date/Time:** With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**Initial pH:** As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

**PAH Total:** With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

**PFAS Total:** With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**Total:** With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

### Data Qualifiers

- A** - Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F** - The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The lower value for the two columns has been reported due to obvious interference.
- J** - Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

**Report Format:** DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

#### Data Qualifiers

Identified Compounds (TICs). For calculated parameters, this represents that one or more values used in the calculation were estimated.

- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND** - Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ** - Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- R** - Analytical results are from sample re-analysis.
- RE** - Analytical results are from sample re-extraction.
- S** - Analytical results are from modified screening analysis.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2521646  
**Report Date:** 04/18/25

## REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

## LIMITATION OF LIABILITIES

Pace Analytical Services performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Pace Analytical Services shall be to re-perform the work at it's own expense. In no event shall Pace Analytical Services be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Pace Analytical Services.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



**Pace Analytical Services LLC**Facility: **Northeast**Department: **Quality Assurance**Title: **Certificate/Approval Program Summary**ID No.: **17873**

Revision 27

Published Date: 01/24/2025

Page 1 of 2

**Certification Information**

The following analytes are not included in our Primary NELAP Scope of Accreditation:

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****EPA 624.1:** m/p-xylene, o-xylene, Naphthalene**EPA 625.1:** alpha-Terpineol**EPA 8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****SM 2540D:** TSS.**EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**MADEP-APH.****Nonpotable Water:** EPA RSK-175 Dissolved Gases**Biological Tissue Matrix:** EPA 3050B**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048****EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**Nonpotable Water:** EPA RSK-175 Dissolved Gases

The following test method is not included in our New Jersey Secondary NELAP Scope of Accreditation:

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Determination of Selected Perfluorinated Alkyl Substances by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry Isotope Dilution (via Alpha SOP 23528)**

The following analytes are included in our Massachusetts DEP Scope of Accreditation

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****Drinking Water****EPA 300.0:** Chloride, Nitrate-N, Fluoride, Sulfate; **EPA 353.2:** Nitrate-N, Nitrite-N; **SM4500NO3-F:** Nitrate-N, Nitrite-N; **SM4500F-C, SM4500CN-CE,****EPA 180.1, SM2130B, SM4500Cl-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B****EPA 524.2:** THMs and VOCs; **EPA 504.1:** EDB, DBCP.**Microbiology:** SM9215B; SM9223-P/A, SM9223B-Colilert-QT, SM9222D.**Non-Potable Water****SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH:** Ammonia-N and Kjeldahl-N, **EPA 350.1:**Ammonia-N, **LACHAT 10-107-06-1-B:** Ammonia-N, **EPA 351.1, SM4500NO3-F, EPA 353.2:** Nitrate-N, **SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300:** Chloride, Sulfate, Nitrate.**EPA 624.1:** Volatile Halocarbons & Aromatics,**EPA 608.3:** Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs**EPA 625.1:** SVOC (Acid/Base/Neutral Extractables).**Microbiology:** SM9223B-Colilert-QT; Enterolert-QT, EPA 1600, EPA 1603, SM9222D.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Drinking Water****EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1 Hg.****EPA 522, EPA 537.1.****Non-Potable Water****EPA 200.7:** Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.**EPA 200.8:** Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.**EPA 245.1 Hg.****SM2340B**

**Pace Analytical Services LLC**ID No.: **17873**Facility: **Northeast**

Revision 27

Department: **Quality Assurance**

Published Date: 01/24/2025

Title: **Certificate/Approval Program Summary**

Page 2 of 2

**Certification IDs:****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581**

CT PH-0826, IL 200077, IN C-MA-03, KY JY98045, ME MA00086, MD 348, MA M-MA086, NH 2064, NJ MA935, NY 11148, NC (DW) 25700, NC (NPW/SCM) 666, OR MA-1316, PA 68-03671, RI LAO00065, TX T104704476, VT VT-0935, VA 460195

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048**

CT PH-0825, ANAB/DoD L2474, IL 200081, IN C-MA-04, KY KY98046, LA 3090, ME MA00030, MI 9110, MN 025-999-495, NH 2062, NJ MA015, NY 11627, NC (NPW/SCM) 685, OR MA-0262, PA 68-02089, RI LAO00299, TX T-104704419, VT VT-0015, VA 460194, WA C954

**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048**

ANAB/DoD L2474, ME MA01156, MN 025-999-498, NH 2249, NJ MA025, NY 12191, OR 4203, TX T104704583, VA 460311, WA C1104.

---

For a complete listing of analytes and methods, please contact your Project Manager.







## ANALYTICAL REPORT

|                 |                                                                             |
|-----------------|-----------------------------------------------------------------------------|
| Lab Number:     | L2522789                                                                    |
| Client:         | Tradepoint Atlantic<br>1600 Sparrows Point Boulevard<br>Baltimore, MD 21219 |
| ATTN:           | Matthew Newman                                                              |
| Phone:          | (443) 649-5073                                                              |
| Project Name:   | B14 1ST HALF 2025 GW                                                        |
| Project Number: | 21010214                                                                    |
| Report Date:    | 04/22/25                                                                    |

The original project report/data package is held by Pace Analytical Services. This report/data package is paginated and should be reproduced only in its entirety. Pace Analytical Services holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0826), IL (200077), IN (C-MA-03), KY (KY98045), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), OR (MA-1316), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #525-23-122-91930A1).

---

Eight Walkup Drive, Westborough, MA 01581-1019  
508-898-9220 (Fax) 508-898-9193 800-624-9220 - [www.alphalab.com](http://www.alphalab.com)



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

| Lab<br>Sample ID | Client ID | Matrix | Sample<br>Location | Collection<br>Date/Time | Receive Date |
|------------------|-----------|--------|--------------------|-------------------------|--------------|
| L2522789-01      | HI22-MWS  | WATER  | B14                | 04/14/25 11:00          | 04/14/25     |
| L2522789-02      | HI12-MWS  | WATER  | B14                | 04/14/25 12:10          | 04/14/25     |
| L2522789-03      | HI10-MWS  | WATER  | B14                | 04/14/25 14:20          | 04/14/25     |
| L2522789-04      | HI11-MWS  | WATER  | B14                | 04/14/25 15:05          | 04/14/25     |
| L2522789-05      | TB-001    | WATER  | B14                | 04/14/25 00:00          | 04/14/25     |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

### Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Pace Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments and solids are reported on a dry weight basis unless otherwise noted. Tissues are reported "as received" or on a wet weight basis, unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

**HOLD POLICY** - For samples submitted on hold, Pace's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Pace Project Manager and made arrangements for Pace to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.

---

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Case Narrative (continued)**

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature: *Tiffani Morrissey* - Tiffani Morrissey

Title: Technical Director/Representative

Date: 04/22/25

# ORGANICS

# VOLATILES



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-01  
**Client ID:** HI22-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 11:00  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/21/25 07:59  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 5.1    |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.23   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25**SAMPLE RESULTS****Lab ID:** L2522789-01**Date Collected:** 04/14/25 11:00**Client ID:** HI22-MWS**Date Received:** 04/14/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 122        |           | 70-130              |
| Toluene-d8            | 100        |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 108        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

Lab ID: L2522789-01  
 Client ID: HI22-MWS  
 Sample Location: B14

Date Collected: 04/14/25 11:00  
 Date Received: 04/14/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/21/25 07:59  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 118        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-02  
**Client ID:** HI12-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 12:10  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/21/25 07:07  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 4.5    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25**SAMPLE RESULTS****Lab ID:** L2522789-02**Date Collected:** 04/14/25 12:10**Client ID:** HI12-MWS**Date Received:** 04/14/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 123        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 98         |           | 70-130              |
| Dibromofluoromethane  | 114        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-02  
**Client ID:** HI12-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 12:10  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/21/25 07:07  
**Analyst:** MCM

| Parameter                                        | Result     | Qualifier | Units     | RL                  | MDL   | Dilution Factor |
|--------------------------------------------------|------------|-----------|-----------|---------------------|-------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |            |           |           |                     |       |                 |
| 1,1,2,2-Tetrachloroethane                        | ND         |           | ug/l      | 0.050               | 0.006 | 1               |
| Surrogate                                        | % Recovery |           | Qualifier | Acceptance Criteria |       |                 |
| 1,2-Dichloroethane-d4                            | 122        |           |           | 70-130              |       |                 |
| 4-Bromofluorobenzene                             | 94         |           |           | 70-130              |       |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-03  
**Client ID:** HI10-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 14:20  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/21/25 08:26  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 2.9    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 1.0    |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | 0.46   | J         | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25**SAMPLE RESULTS****Lab ID:** L2522789-03**Date Collected:** 04/14/25 14:20**Client ID:** HI10-MWS**Date Received:** 04/14/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | 0.36   | J         | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | 0.36   | J         | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 123        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 113        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

Lab ID: L2522789-03  
 Client ID: HI10-MWS  
 Sample Location: B14

Date Collected: 04/14/25 14:20  
 Date Received: 04/14/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/21/25 08:26  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 119        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 91         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-04  
**Client ID:** HI11-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 15:05  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/21/25 07:33  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 5.2    |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

Project Name: B14 1ST HALF 2025 GW

Lab Number: L2522789

Project Number: 21010214

Report Date: 04/22/25

## SAMPLE RESULTS

Lab ID: L2522789-04

Date Collected: 04/14/25 15:05

Client ID: HI11-MWS

Date Received: 04/14/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 123        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 110        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-04  
**Client ID:** HI11-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 15:05  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/21/25 07:33  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 119        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 92         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-05  
**Client ID:** TB-001  
**Sample Location:** B14

**Date Collected:** 04/14/25 00:00  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/21/25 06:19  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | ND     |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



Project Name: B14 1ST HALF 2025 GW

Lab Number: L2522789

Project Number: 21010214

Report Date: 04/22/25

## SAMPLE RESULTS

Lab ID: L2522789-05

Date Collected: 04/14/25 00:00

Client ID: TB-001

Date Received: 04/14/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 123        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 98         |           | 70-130              |
| Dibromofluoromethane  | 110        |           | 70-130              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-05  
**Client ID:** TB-001  
**Sample Location:** B14

**Date Collected:** 04/14/25 00:00  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/21/25 06:19  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 120        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 04/21/25 05:53  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2056533-5 |        |           |       |      |      |
| Dichlorodifluoromethane                                                              | ND     |           | ug/l  | 5.0  | 0.24 |
| Chloromethane                                                                        | ND     |           | ug/l  | 2.5  | 0.20 |
| Vinyl chloride                                                                       | ND     |           | ug/l  | 1.0  | 0.07 |
| Bromomethane                                                                         | ND     |           | ug/l  | 1.0  | 0.26 |
| Chloroethane                                                                         | ND     |           | ug/l  | 1.0  | 0.13 |
| Trichlorofluoromethane                                                               | ND     |           | ug/l  | 2.5  | 0.16 |
| 1,1-Dichloroethene                                                                   | ND     |           | ug/l  | 0.50 | 0.17 |
| Carbon disulfide                                                                     | ND     |           | ug/l  | 5.0  | 0.30 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                | ND     |           | ug/l  | 2.5  | 0.15 |
| Methylene chloride                                                                   | ND     |           | ug/l  | 2.5  | 0.68 |
| Acetone                                                                              | ND     |           | ug/l  | 5.0  | 1.5  |
| trans-1,2-Dichloroethene                                                             | ND     |           | ug/l  | 0.75 | 0.16 |
| Methyl Acetate                                                                       | ND     |           | ug/l  | 2.0  | 0.23 |
| Methyl tert butyl ether                                                              | ND     |           | ug/l  | 1.0  | 0.17 |
| 1,1-Dichloroethane                                                                   | ND     |           | ug/l  | 0.75 | 0.21 |
| cis-1,2-Dichloroethene                                                               | ND     |           | ug/l  | 0.50 | 0.19 |
| Cyclohexane                                                                          | ND     |           | ug/l  | 10   | 0.27 |
| Chloroform                                                                           | ND     |           | ug/l  | 0.75 | 0.22 |
| Carbon tetrachloride                                                                 | ND     |           | ug/l  | 0.50 | 0.13 |
| 1,1,1-Trichloroethane                                                                | ND     |           | ug/l  | 0.50 | 0.16 |
| 2-Butanone                                                                           | ND     |           | ug/l  | 5.0  | 1.9  |
| Benzene                                                                              | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,2-Dichloroethane                                                                   | ND     |           | ug/l  | 0.50 | 0.13 |
| Trichloroethene                                                                      | ND     |           | ug/l  | 0.50 | 0.18 |
| 1,2-Dichloropropane                                                                  | ND     |           | ug/l  | 1.0  | 0.14 |
| Bromodichloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.19 |
| cis-1,3-Dichloropropene                                                              | ND     |           | ug/l  | 0.50 | 0.14 |
| Toluene                                                                              | ND     |           | ug/l  | 0.75 | 0.20 |
| Tetrachloroethene                                                                    | ND     |           | ug/l  | 0.50 | 0.18 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8260D  
**Analytical Date:** 04/21/25 05:53  
**Analyst:** MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2056533-5 |        |           |       |      |      |
| 4-Methyl-2-pentanone                                                                 | ND     |           | ug/l  | 5.0  | 0.42 |
| trans-1,3-Dichloropropene                                                            | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,3-Dichloropropene, Total                                                           | ND     |           | ug/l  | 0.50 | 0.14 |
| 1,1,2-Trichloroethane                                                                | ND     |           | ug/l  | 0.75 | 0.14 |
| Dibromochloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.15 |
| 1,2-Dibromoethane                                                                    | ND     |           | ug/l  | 2.0  | 0.19 |
| 2-Hexanone                                                                           | ND     |           | ug/l  | 5.0  | 0.52 |
| Chlorobenzene                                                                        | ND     |           | ug/l  | 0.50 | 0.18 |
| Ethylbenzene                                                                         | ND     |           | ug/l  | 0.50 | 0.17 |
| p/m-Xylene                                                                           | ND     |           | ug/l  | 1.0  | 0.33 |
| o-Xylene                                                                             | ND     |           | ug/l  | 1.0  | 0.39 |
| Xylenes, Total                                                                       | ND     |           | ug/l  | 1.0  | 0.33 |
| Styrene                                                                              | ND     |           | ug/l  | 1.0  | 0.36 |
| Bromoform                                                                            | ND     |           | ug/l  | 2.0  | 0.25 |
| Isopropylbenzene                                                                     | ND     |           | ug/l  | 0.50 | 0.19 |
| 1,1,2,2-Tetrachloroethane                                                            | ND     |           | ug/l  | 0.50 | 0.17 |
| 1,3-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,4-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,2-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.18 |
| 1,2-Dibromo-3-chloropropane                                                          | ND     |           | ug/l  | 2.5  | 0.35 |
| 1,2,4-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.22 |
| 1,2,3-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.23 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
Analytical Date: 04/21/25 05:53  
Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL | MDL |
|--------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2056533-5 |        |           |       |    |     |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 122       |           | 70-130                 |
| Toluene-d8            | 99        |           | 70-130                 |
| 4-Bromofluorobenzene  | 100       |           | 70-130                 |
| Dibromofluoromethane  | 110       |           | 70-130                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D-SIM(M)  
Analytical Date: 04/21/25 05:53  
Analyst: MCM

| Parameter                                                                                | Result | Qualifier | Units | RL    | MDL   |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-------|-------|
| Volatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-05 Batch: WG2056541-5 |        |           |       |       |       |
| 1,1,1,2-Tetrachloroethane                                                                | ND     |           | ug/l  | 0.050 | 0.006 |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 119       |           | 70-130                 |
| 4-Bromofluorobenzene  | 93        |           | 70-130                 |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522789

**Report Date:** 04/22/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2056533-3 WG2056533-4 |                  |      |                   |      |                     |     |      |               |
| Dichlorodifluoromethane                                                                                 | 100              |      | 110               |      | 36-147              | 10  |      | 20            |
| Chloromethane                                                                                           | 120              |      | 120               |      | 64-130              | 0   |      | 20            |
| Vinyl chloride                                                                                          | 110              |      | 110               |      | 55-140              | 0   |      | 20            |
| Bromomethane                                                                                            | 45               |      | 52                |      | 39-139              | 14  |      | 20            |
| Chloroethane                                                                                            | 130              |      | 130               |      | 55-138              | 0   |      | 20            |
| Trichlorofluoromethane                                                                                  | 110              |      | 110               |      | 62-150              | 0   |      | 20            |
| 1,1-Dichloroethene                                                                                      | 100              |      | 100               |      | 61-145              | 0   |      | 20            |
| Carbon disulfide                                                                                        | 120              |      | 120               |      | 51-130              | 0   |      | 20            |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                                   | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Methylene chloride                                                                                      | 100              |      | 110               |      | 70-130              | 10  |      | 20            |
| Acetone                                                                                                 | 120              |      | 130               |      | 58-148              | 8   |      | 20            |
| trans-1,2-Dichloroethene                                                                                | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Methyl Acetate                                                                                          | 140              | Q    | 140               | Q    | 70-130              | 0   |      | 20            |
| Methyl tert butyl ether                                                                                 | 98               |      | 100               |      | 63-130              | 2   |      | 20            |
| 1,1-Dichloroethane                                                                                      | 120              |      | 120               |      | 70-130              | 0   |      | 20            |
| cis-1,2-Dichloroethene                                                                                  | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Cyclohexane                                                                                             | 120              |      | 120               |      | 70-130              | 0   |      | 20            |
| Chloroform                                                                                              | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Carbon tetrachloride                                                                                    | 100              |      | 110               |      | 63-132              | 10  |      | 20            |
| 1,1,1-Trichloroethane                                                                                   | 100              |      | 110               |      | 67-130              | 10  |      | 20            |
| 2-Butanone                                                                                              | 110              |      | 110               |      | 63-138              | 0   |      | 20            |
| Benzene                                                                                                 | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| 1,2-Dichloroethane                                                                                      | 120              |      | 120               |      | 70-130              | 0   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522789

**Report Date:** 04/22/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2056533-3 WG2056533-4 |                  |      |                   |      |                     |     |      |               |
| Trichloroethene                                                                                         | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| 1,2-Dichloropropane                                                                                     | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Bromodichloromethane                                                                                    | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| cis-1,3-Dichloropropene                                                                                 | 93               |      | 96                |      | 70-130              | 3   |      | 20            |
| Toluene                                                                                                 | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Tetrachloroethene                                                                                       | 97               |      | 96                |      | 70-130              | 1   |      | 20            |
| 4-Methyl-2-pentanone                                                                                    | 90               |      | 89                |      | 59-130              | 1   |      | 20            |
| trans-1,3-Dichloropropene                                                                               | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| 1,1,2-Trichloroethane                                                                                   | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Dibromochloromethane                                                                                    | 93               |      | 95                |      | 63-130              | 2   |      | 20            |
| 1,2-Dibromoethane                                                                                       | 95               |      | 98                |      | 70-130              | 3   |      | 20            |
| 2-Hexanone                                                                                              | 98               |      | 98                |      | 57-130              | 0   |      | 20            |
| Chlorobenzene                                                                                           | 98               |      | 97                |      | 75-130              | 1   |      | 20            |
| Ethylbenzene                                                                                            | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| p/m-Xylene                                                                                              | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| o-Xylene                                                                                                | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Styrene                                                                                                 | 95               |      | 100               |      | 70-130              | 5   |      | 20            |
| Bromoform                                                                                               | 88               |      | 91                |      | 54-136              | 3   |      | 20            |
| Isopropylbenzene                                                                                        | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| 1,1,2,2-Tetrachloroethane                                                                               | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| 1,3-Dichlorobenzene                                                                                     | 99               |      | 100               |      | 70-130              | 1   |      | 20            |
| 1,4-Dichlorobenzene                                                                                     | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| 1,2-Dichlorobenzene                                                                                     | 98               |      | 100               |      | 70-130              | 2   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522789

**Project Number:** 21010214

**Report Date:** 04/22/25

| <b>Parameter</b>                                                                                        | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|---------------------------------------------------------------------------------------------------------|--------------------------|-------------|--------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2056533-3 WG2056533-4 |                          |             |                          |             |                             |            |             |                       |
| 1,2-Dibromo-3-chloropropane                                                                             | 80                       |             | 83                       |             | 41-144                      | 4          |             | 20                    |
| 1,2,4-Trichlorobenzene                                                                                  | 90                       |             | 94                       |             | 70-130                      | 4          |             | 20                    |
| 1,2,3-Trichlorobenzene                                                                                  | 91                       |             | 97                       |             | 70-130                      | 6          |             | 20                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|--------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 118                      |             | 117                      |             | 70-130                         |
| Toluene-d8            | 101                      |             | 100                      |             | 70-130                         |
| 4-Bromofluorobenzene  | 94                       |             | 97                       |             | 70-130                         |
| Dibromofluoromethane  | 106                      |             | 104                      |             | 70-130                         |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522789

**Project Number:** 21010214

**Report Date:** 04/22/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-05 Batch: WG2056541-3 WG2056541-4 |                          |             |                           |             |                             |            |             |                       |
| 1,1,1,2-Tetrachloroethane                                                                                   | 110                      |             | 106                       |             | 70-130                      | 4          |             | 25                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 119                      |             | 121                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 92                       |             | 91                        |             | 70-130                         |

# SEMIVOLATILES

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-01  
**Client ID:** HI22-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 11:00  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/19/25 14:30  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25**SAMPLE RESULTS**

Lab ID: L2522789-01

Date Collected: 04/14/25 11:00

Client ID: HI22-MWS

Date Received: 04/14/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 41         |           | 21-120              |
| Phenol-d6            | 34         |           | 10-120              |
| Nitrobenzene-d5      | 50         |           | 23-120              |
| 2-Fluorobiphenyl     | 49         |           | 15-120              |
| 2,4,6-Tribromophenol | 58         |           | 10-120              |
| 4-Terphenyl-d14      | 68         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-01  
**Client ID:** HI22-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 11:00  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/19/25 14:02  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.16   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.03   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.07   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.08   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.13   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | 0.07   |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | 0.05   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 71         |           | 23-120              |
| 2-Fluorobiphenyl | 61         |           | 15-120              |
| 4-Terphenyl-d14  | 78         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-02  
**Client ID:** HI12-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 12:10  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/19/25 14:52  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25**SAMPLE RESULTS****Lab ID:** L2522789-02**Date Collected:** 04/14/25 12:10**Client ID:** HI12-MWS**Date Received:** 04/14/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 50         |           | 21-120              |
| Phenol-d6            | 42         |           | 10-120              |
| Nitrobenzene-d5      | 72         |           | 23-120              |
| 2-Fluorobiphenyl     | 67         |           | 15-120              |
| 2,4,6-Tribromophenol | 91         |           | 10-120              |
| 4-Terphenyl-d14      | 96         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-02  
**Client ID:** HI12-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 12:10  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/19/25 14:36  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.20   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.07   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.09   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.10   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.08   | J         | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | 0.09   |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | 0.07   |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | 0.03   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 95         |           | 23-120              |
| 2-Fluorobiphenyl | 74         |           | 15-120              |
| 4-Terphenyl-d14  | 99         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-03  
**Client ID:** HI10-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 14:20  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/19/25 15:14  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 32     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | 1.7    | J         | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | 0.32   | J         | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25**SAMPLE RESULTS**

Lab ID: L2522789-03

Date Collected: 04/14/25 14:20

Client ID: HI10-MWS

Date Received: 04/14/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | 0.48   | J         | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | 1.8    | J         | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | 1.4    | J         | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | 2.3    |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | 4.2    |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | 0.56   | J         | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | 0.47   | J         | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 56         |           | 21-120              |
| Phenol-d6            | 39         |           | 10-120              |
| Nitrobenzene-d5      | 76         |           | 23-120              |
| 2-Fluorobiphenyl     | 67         |           | 15-120              |
| 2,4,6-Tribromophenol | 80         |           | 10-120              |
| 4-Terphenyl-d14      | 90         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-03  
**Client ID:** HI10-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 14:20  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/19/25 14:53  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 33     |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 1.9    |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.57   |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 2.2    |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 1.8    |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 2.8    |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.40   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.72   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.53   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | 0.07   |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | 0.03   | J         | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 100        |           | 23-120              |
| 2-Fluorobiphenyl | 76         |           | 15-120              |
| 4-Terphenyl-d14  | 103        |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-04  
**Client ID:** HI11-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 15:05  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/19/25 15:37  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 0.83   | J         | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25**SAMPLE RESULTS****Lab ID:** L2522789-04**Date Collected:** 04/14/25 15:05**Client ID:** HI11-MWS**Date Received:** 04/14/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 44         |           | 21-120              |
| Phenol-d6            | 34         |           | 10-120              |
| Nitrobenzene-d5      | 65         |           | 23-120              |
| 2-Fluorobiphenyl     | 56         |           | 15-120              |
| 2,4,6-Tribromophenol | 81         |           | 10-120              |
| 4-Terphenyl-d14      | 90         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**SAMPLE RESULTS**

**Lab ID:** L2522789-04  
**Client ID:** HI11-MWS  
**Sample Location:** B14

**Date Collected:** 04/14/25 15:05  
**Date Received:** 04/14/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/19/25 15:10  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.92   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.05   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.07   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.06   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.12   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.07   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.06   | J         | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | 0.04   | J         | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 85         |           | 23-120              |
| 2-Fluorobiphenyl | 70         |           | 15-120              |
| 4-Terphenyl-d14  | 90         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/19/25 13:01  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2055925-1 |        |           |       |     |      |
| Benzaldehyde                                                                             | ND     |           | ug/l  | 5.0 | 1.1  |
| Phenol                                                                                   | ND     |           | ug/l  | 5.0 | 0.35 |
| Bis(2-chloroethyl)ether                                                                  | ND     |           | ug/l  | 2.0 | 0.39 |
| 2-Chlorophenol                                                                           | ND     |           | ug/l  | 2.0 | 0.65 |
| 2-Methylphenol                                                                           | ND     |           | ug/l  | 5.0 | 2.3  |
| Bis(2-chloroisopropyl)ether                                                              | ND     |           | ug/l  | 2.0 | 0.40 |
| Acetophenone                                                                             | ND     |           | ug/l  | 5.0 | 0.92 |
| n-Nitrosodi-n-propylamine                                                                | ND     |           | ug/l  | 5.0 | 0.91 |
| 3-Methylphenol/4-Methylphenol                                                            | ND     |           | ug/l  | 5.0 | 1.4  |
| Hexachloroethane                                                                         | ND     |           | ug/l  | 2.0 | 0.20 |
| Nitrobenzene                                                                             | ND     |           | ug/l  | 1.4 | 0.20 |
| Isophorone                                                                               | ND     |           | ug/l  | 5.0 | 0.86 |
| 2,4-Dimethylphenol                                                                       | ND     |           | ug/l  | 5.0 | 2.0  |
| Bis(2-chloroethoxy)methane                                                               | ND     |           | ug/l  | 5.0 | 0.84 |
| 2,4-Dichlorophenol                                                                       | ND     |           | ug/l  | 5.0 | 1.7  |
| Naphthalene                                                                              | ND     |           | ug/l  | 2.0 | 0.54 |
| 4-Chloroaniline                                                                          | ND     |           | ug/l  | 3.7 | 0.47 |
| Hexachlorobutadiene                                                                      | ND     |           | ug/l  | 2.0 | 0.36 |
| Caprolactam                                                                              | ND     |           | ug/l  | 10  | 1.2  |
| 2-Methylnaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.37 |
| Hexachlorocyclopentadiene                                                                | ND     |           | ug/l  | 20  | 1.2  |
| 1,2,4,5-Tetrachlorobenzene                                                               | ND     |           | ug/l  | 1.7 | 0.24 |
| 2,4,6-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| 2,4,5-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| Biphenyl                                                                                 | ND     |           | ug/l  | 2.0 | 0.20 |
| 2-Chloronaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.35 |
| 2-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.0  |
| 2,6-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.84 |
| Acenaphthylene                                                                           | ND     |           | ug/l  | 2.0 | 0.32 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/19/25 13:01  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2055925-1 |        |           |       |     |      |
| Acenaphthene                                                                             | ND     |           | ug/l  | 2.0 | 0.40 |
| 2,4-Dinitrophenol                                                                        | ND     |           | ug/l  | 20  | 5.4  |
| 2,4-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.54 |
| 2,3,4,6-Tetrachlorophenol                                                                | ND     |           | ug/l  | 5.0 | 2.2  |
| Diethyl phthalate                                                                        | ND     |           | ug/l  | 5.0 | 0.76 |
| Fluorene                                                                                 | ND     |           | ug/l  | 2.0 | 0.44 |
| 4-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.4  |
| NDPA/DPA                                                                                 | ND     |           | ug/l  | 2.0 | 0.92 |
| Hexachlorobenzene                                                                        | ND     |           | ug/l  | 2.0 | 0.45 |
| Pentachlorophenol                                                                        | ND     |           | ug/l  | 10  | 2.5  |
| Phenanthrene                                                                             | ND     |           | ug/l  | 2.0 | 0.42 |
| Anthracene                                                                               | ND     |           | ug/l  | 2.0 | 0.47 |
| Carbazole                                                                                | ND     |           | ug/l  | 2.0 | 0.31 |
| Di-n-butylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 0.96 |
| Fluoranthene                                                                             | ND     |           | ug/l  | 2.0 | 0.41 |
| Pyrene                                                                                   | ND     |           | ug/l  | 2.0 | 0.41 |
| 3,3'-Dichlorobenzidine                                                                   | ND     |           | ug/l  | 5.0 | 1.8  |
| Benzo(a)anthracene                                                                       | ND     |           | ug/l  | 2.0 | 0.32 |
| Chrysene                                                                                 | ND     |           | ug/l  | 1.5 | 0.22 |
| Bis(2-ethylhexyl)phthalate                                                               | ND     |           | ug/l  | 3.0 | 1.4  |
| Di-n-octylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 2.3  |
| Benzo(b)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.53 |
| Benzo(k)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.62 |
| Benzo(a)pyrene                                                                           | ND     |           | ug/l  | 2.0 | 0.37 |
| Indeno(1,2,3-cd)pyrene                                                                   | ND     |           | ug/l  | 2.0 | 0.48 |
| Dibenzo(a,h)anthracene                                                                   | ND     |           | ug/l  | 2.0 | 0.29 |
| Benzo(ghi)perylene                                                                       | ND     |           | ug/l  | 2.0 | 0.37 |





**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8270E  
Analytical Date: 04/19/25 13:01  
Analyst: SLR

Extraction Method: EPA 3510C  
Extraction Date: 04/19/25 01:15

| Parameter                                                                                | Result | Qualifier | Units | RL | MDL |
|------------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2055925-1 |        |           |       |    |     |

| Surrogate            | %Recovery | Qualifier | Acceptance<br>Criteria |
|----------------------|-----------|-----------|------------------------|
| 2-Fluorophenol       | 41        |           | 21-120                 |
| Phenol-d6            | 33        |           | 10-120                 |
| Nitrobenzene-d5      | 56        |           | 23-120                 |
| 2-Fluorobiphenyl     | 47        |           | 15-120                 |
| 2,4,6-Tribromophenol | 65        |           | 10-120                 |
| 4-Terphenyl-d14      | 73        |           | 41-149                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/19/25 13:11  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/19/25 01:15

| Parameter                                                                                    | Result | Qualifier | Units | RL   | MDL  |
|----------------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-04 Batch: WG2055927-1 |        |           |       |      |      |
| Naphthalene                                                                                  | 0.04   | J         | ug/l  | 0.10 | 0.02 |
| 2-Methylnaphthalene                                                                          | ND     |           | ug/l  | 0.10 | 0.03 |
| Acenaphthylene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Acenaphthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluorene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Phenanthrene                                                                                 | ND     |           | ug/l  | 0.05 | 0.04 |
| Anthracene                                                                                   | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluoranthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.03 |
| Pyrene                                                                                       | ND     |           | ug/l  | 0.10 | 0.04 |
| Benzo(a)anthracene                                                                           | ND     |           | ug/l  | 0.05 | 0.03 |
| Chrysene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(b)fluoranthene                                                                         | ND     |           | ug/l  | 0.05 | 0.03 |
| Benzo(k)fluoranthene                                                                         | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(a)pyrene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Indeno(1,2,3-cd)pyrene                                                                       | ND     |           | ug/l  | 0.10 | 0.02 |
| Dibenzo(a,h)anthracene                                                                       | ND     |           | ug/l  | 0.05 | 0.02 |
| Benzo(ghi)perylene                                                                           | ND     |           | ug/l  | 0.10 | 0.02 |

| Surrogate        | %Recovery | Qualifier | Acceptance Criteria |
|------------------|-----------|-----------|---------------------|
| Nitrobenzene-d5  | 68        |           | 23-120              |
| 2-Fluorobiphenyl | 48        |           | 15-120              |
| 4-Terphenyl-d14  | 75        |           | 41-149              |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522789

**Report Date:** 04/22/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2055925-2 WG2055925-3 |                  |      |                   |      |                     |     |      |               |
| Benzaldehyde                                                                                                | 80               |      | 80                |      | 40-140              | 0   |      | 30            |
| Phenol                                                                                                      | 44               |      | 40                |      | 12-110              | 10  |      | 30            |
| Bis(2-chloroethyl)ether                                                                                     | 78               |      | 74                |      | 40-140              | 5   |      | 30            |
| 2-Chlorophenol                                                                                              | 64               |      | 64                |      | 27-123              | 0   |      | 30            |
| 2-Methylphenol                                                                                              | 68               |      | 70                |      | 30-130              | 3   |      | 30            |
| Bis(2-chloroisopropyl)ether                                                                                 | 74               |      | 74                |      | 40-140              | 0   |      | 30            |
| Acetophenone                                                                                                | 81               |      | 82                |      | 39-129              | 1   |      | 30            |
| n-Nitrosodi-n-propylamine                                                                                   | 75               |      | 80                |      | 29-132              | 6   |      | 30            |
| 3-Methylphenol/4-Methylphenol                                                                               | 72               |      | 72                |      | 30-130              | 0   |      | 30            |
| Hexachloroethane                                                                                            | 36               | Q    | 50                |      | 40-140              | 33  | Q    | 30            |
| Nitrobenzene                                                                                                | 68               |      | 71                |      | 40-140              | 4   |      | 30            |
| Isophorone                                                                                                  | 73               |      | 72                |      | 40-140              | 1   |      | 30            |
| 2,4-Dimethylphenol                                                                                          | 68               |      | 65                |      | 30-130              | 5   |      | 30            |
| Bis(2-chloroethoxy)methane                                                                                  | 73               |      | 74                |      | 40-140              | 1   |      | 30            |
| 2,4-Dichlorophenol                                                                                          | 72               |      | 74                |      | 30-130              | 3   |      | 30            |
| Naphthalene                                                                                                 | 56               |      | 62                |      | 40-140              | 10  |      | 30            |
| 4-Chloroaniline                                                                                             | 58               |      | 52                |      | 40-140              | 11  |      | 30            |
| Hexachlorobutadiene                                                                                         | 33               | Q    | 47                |      | 40-140              | 35  | Q    | 30            |
| Caprolactam                                                                                                 | 36               |      | 34                |      | 10-130              | 6   |      | 30            |
| 2-Methylnaphthalene                                                                                         | 60               |      | 64                |      | 40-140              | 6   |      | 30            |
| Hexachlorocyclopentadiene                                                                                   | 34               | Q    | 44                |      | 40-140              | 26  |      | 30            |
| 1,2,4,5-Tetrachlorobenzene                                                                                  | 60               |      | 65                |      | 2-134               | 8   |      | 30            |
| 2,4,6-Trichlorophenol                                                                                       | 69               |      | 71                |      | 30-130              | 3   |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522789

**Report Date:** 04/22/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2055925-2 WG2055925-3 |                  |      |                   |      |                     |     |      |               |
| 2,4,5-Trichlorophenol                                                                                       | 75               |      | 76                |      | 30-130              | 1   |      | 30            |
| Biphenyl                                                                                                    | 66               |      | 77                |      | 40-140              | 15  |      | 30            |
| 2-Chloronaphthalene                                                                                         | 62               |      | 70                |      | 40-140              | 12  |      | 30            |
| 2-Nitroaniline                                                                                              | 74               |      | 80                |      | 52-143              | 8   |      | 30            |
| 2,6-Dinitrotoluene                                                                                          | 77               |      | 79                |      | 40-140              | 3   |      | 30            |
| Acenaphthylene                                                                                              | 68               |      | 72                |      | 45-123              | 6   |      | 30            |
| Acenaphthene                                                                                                | 69               |      | 79                |      | 37-111              | 14  |      | 30            |
| 2,4-Dinitrophenol                                                                                           | 41               |      | 47                |      | 20-130              | 14  |      | 30            |
| 2,4-Dinitrotoluene                                                                                          | 76               |      | 85                |      | 48-143              | 11  |      | 30            |
| 2,3,4,6-Tetrachlorophenol                                                                                   | 84               |      | 87                |      | 54-145              | 4   |      | 30            |
| Diethyl phthalate                                                                                           | 80               |      | 87                |      | 40-140              | 8   |      | 30            |
| Fluorene                                                                                                    | 74               |      | 81                |      | 40-140              | 9   |      | 30            |
| 4-Nitroaniline                                                                                              | 89               |      | 97                |      | 51-143              | 9   |      | 30            |
| NDPA/DPA                                                                                                    | 78               |      | 84                |      | 40-140              | 7   |      | 30            |
| Hexachlorobenzene                                                                                           | 73               |      | 80                |      | 40-140              | 9   |      | 30            |
| Pentachlorophenol                                                                                           | 80               |      | 86                |      | 9-103               | 7   |      | 30            |
| Phenanthrene                                                                                                | 82               |      | 87                |      | 40-140              | 6   |      | 30            |
| Anthracene                                                                                                  | 84               |      | 88                |      | 40-140              | 5   |      | 30            |
| Carbazole                                                                                                   | 88               |      | 92                |      | 55-144              | 4   |      | 30            |
| Di-n-butylphthalate                                                                                         | 85               |      | 90                |      | 40-140              | 6   |      | 30            |
| Fluoranthene                                                                                                | 87               |      | 90                |      | 40-140              | 3   |      | 30            |
| Pyrene                                                                                                      | 84               |      | 88                |      | 26-127              | 5   |      | 30            |
| 3,3'-Dichlorobenzidine                                                                                      | 71               |      | 71                |      | 40-140              | 0   |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522789

**Project Number:** 21010214

**Report Date:** 04/22/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2055925-2 WG2055925-3 |                          |             |                           |             |                             |            |             |                       |
| Benzo(a)anthracene                                                                                          | 88                       |             | 91                        |             | 40-140                      | 3          |             | 30                    |
| Chrysene                                                                                                    | 87                       |             | 92                        |             | 40-140                      | 6          |             | 30                    |
| Bis(2-ethylhexyl)phthalate                                                                                  | 90                       |             | 96                        |             | 40-140                      | 6          |             | 30                    |
| Di-n-octylphthalate                                                                                         | 91                       |             | 99                        |             | 40-140                      | 8          |             | 30                    |
| Benzo(b)fluoranthene                                                                                        | 89                       |             | 96                        |             | 40-140                      | 8          |             | 30                    |
| Benzo(k)fluoranthene                                                                                        | 86                       |             | 95                        |             | 40-140                      | 10         |             | 30                    |
| Benzo(a)pyrene                                                                                              | 91                       |             | 96                        |             | 40-140                      | 5          |             | 30                    |
| Indeno(1,2,3-cd)pyrene                                                                                      | 84                       |             | 88                        |             | 40-140                      | 5          |             | 30                    |
| Dibenzo(a,h)anthracene                                                                                      | 88                       |             | 93                        |             | 40-140                      | 6          |             | 30                    |
| Benzo(ghi)perylene                                                                                          | 89                       |             | 92                        |             | 40-140                      | 3          |             | 30                    |

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 44                       |             | 52                        |             | 21-120                         |
| Phenol-d6            | 37                       |             | 41                        |             | 10-120                         |
| Nitrobenzene-d5      | 60                       |             | 74                        |             | 23-120                         |
| 2-Fluorobiphenyl     | 45                       |             | 57                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 71                       |             | 82                        |             | 10-120                         |
| 4-Terphenyl-d14      | 75                       |             | 81                        |             | 41-149                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522789

**Report Date:** 04/22/25

| Parameter                                                                                                       | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-----------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2055927-2 WG2055927-3 |                  |      |                   |      |                     |     |      |               |
| Naphthalene                                                                                                     | 48               |      | 55                |      | 40-140              | 14  |      | 40            |
| 2-Methylnaphthalene                                                                                             | 51               |      | 59                |      | 40-140              | 15  |      | 40            |
| Acenaphthylene                                                                                                  | 70               |      | 76                |      | 40-140              | 8   |      | 40            |
| Acenaphthene                                                                                                    | 60               |      | 66                |      | 37-111              | 10  |      | 40            |
| Fluorene                                                                                                        | 70               |      | 75                |      | 40-140              | 7   |      | 40            |
| Phenanthrene                                                                                                    | 67               |      | 72                |      | 40-140              | 7   |      | 40            |
| Anthracene                                                                                                      | 71               |      | 76                |      | 40-140              | 7   |      | 40            |
| Fluoranthene                                                                                                    | 77               |      | 85                |      | 40-140              | 10  |      | 40            |
| Pyrene                                                                                                          | 77               |      | 86                |      | 26-127              | 11  |      | 40            |
| Benzo(a)anthracene                                                                                              | 70               |      | 75                |      | 40-140              | 7   |      | 40            |
| Chrysene                                                                                                        | 72               |      | 77                |      | 40-140              | 7   |      | 40            |
| Benzo(b)fluoranthene                                                                                            | 76               |      | 83                |      | 40-140              | 9   |      | 40            |
| Benzo(k)fluoranthene                                                                                            | 73               |      | 79                |      | 40-140              | 8   |      | 40            |
| Benzo(a)pyrene                                                                                                  | 81               |      | 88                |      | 40-140              | 8   |      | 40            |
| Indeno(1,2,3-cd)pyrene                                                                                          | 95               |      | 103               |      | 40-140              | 8   |      | 40            |
| Dibenzo(a,h)anthracene                                                                                          | 90               |      | 98                |      | 40-140              | 9   |      | 40            |
| Benzo(ghi)perylene                                                                                              | 83               |      | 91                |      | 40-140              | 9   |      | 40            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522789

**Project Number:** 21010214

**Report Date:** 04/22/25

| <b>Parameter</b>                                                                                                | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-----------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2055927-2 WG2055927-3 |                          |             |                           |             |                             |            |             |                       |

| <b>Surrogate</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| Nitrobenzene-d5  | 65                       |             | 74                        |             | 23-120                         |
| 2-Fluorobiphenyl | 45                       |             | 57                        |             | 15-120                         |
| 4-Terphenyl-d14  | 69                       |             | 77                        |             | 41-149                         |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

Serial\_No:04222513:37  
**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

**Cooler Information**

|               |                     |
|---------------|---------------------|
| <b>Cooler</b> | <b>Custody Seal</b> |
| A             | Absent              |

**Container Information**

| Container ID | Container Type          | Cooler | Initial pH | Final pH | Temp deg C | Pres | Seal   | Frozen Date/Time | Analysis(*)                      |
|--------------|-------------------------|--------|------------|----------|------------|------|--------|------------------|----------------------------------|
| L2522789-01A | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-01B | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-01C | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-01D | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522789-01E | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522789-02A | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-02B | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-02C | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-02D | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522789-02E | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522789-03A | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-03B | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-03C | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-03D | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522789-03E | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522789-04A | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-04B | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-04C | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-04D | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522789-04E | Amber 100ml unpreserved | A      | 7          | 7        | 3.3        | Y    | Absent |                  | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522789-05A | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-05B | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |
| L2522789-05C | Vial HCl preserved      | A      | NA         |          | 3.3        | Y    | Absent |                  | PA-8260-SIM(14),PA-8260(14)      |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

Serial\_No:04222513:37  
**Lab Number:** L2522789  
**Report Date:** 04/22/25

**Container Information**

| <b>Container ID</b> | <b>Container Type</b> | <b>Cooler</b> | <b>Initial<br/>pH</b> | <b>Final<br/>pH</b> | <b>Temp<br/>deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen<br/>Date/Time</b> | <b>Analysis(*)</b>          |
|---------------------|-----------------------|---------------|-----------------------|---------------------|-----------------------|-------------|-------------|-----------------------------|-----------------------------|
| L2522789-05D        | Vial HCl preserved    | A             | NA                    |                     | 3.3                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

## GLOSSARY

### Acronyms

|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| DL       | - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                |
| EDL      | - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).                                                                                         |
| EMPC     | - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.                                                                                                                                                                                                                               |
| EPA      | - Environmental Protection Agency.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| LCS      | - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                          |
| LCSD     | - Laboratory Control Sample Duplicate: Refer to LCS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |
| LFB      | - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                         |
| LOD      | - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                                                               |
| LOQ      | - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)<br><br>Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| MDL      | - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                          |
| MS       | - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.                                                                                                                                                                                                    |
| MSD      | - Matrix Spike Sample Duplicate: Refer to MS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NA       | - Not Applicable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| NC       | - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.                                                                                                                                                                                                                                                                                                                                                                           |
| NDPA/DPA | - N-Nitrosodiphenylamine/Diphenylamine.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| NI       | - Not Ignitable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| NP       | - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NR       | - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.                                                                                                                                                                                                                                                                                                                                                                    |
| RL       | - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                                                                                                                                   |
| RPD      | - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.                                                                  |
| SRM      | - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.                                                                                                                                                                                                                                                                                                                                                                     |
| STLP     | - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| TEF      | - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.                                                                                                                                                                                                                                                                                                                                                                                             |
| TEQ      | - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.                                                                                                                                                                                                                                                                                                                                                        |
| TIC      | - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.                                                                                                                                                                                                                                                                      |

Report Format: DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

### Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

### Terms

**Analytical Method:** Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

**Chlordane:** The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

**Difference:** With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

**Final pH:** As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

**Frozen Date/Time:** With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**Initial pH:** As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

**PAH Total:** With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

**PFAS Total:** With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**Total:** With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

### Data Qualifiers

- A** - Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F** - The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The lower value for the two columns has been reported due to obvious interference.
- J** - Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

**Report Format:** DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522789  
**Report Date:** 04/22/25

#### Data Qualifiers

Identified Compounds (TICs). For calculated parameters, this represents that one or more values used in the calculation were estimated.

- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND** - Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ** - Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- R** - Analytical results are from sample re-analysis.
- RE** - Analytical results are from sample re-extraction.
- S** - Analytical results are from modified screening analysis.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522789**Project Number:** 21010214**Report Date:** 04/22/25

## REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

## LIMITATION OF LIABILITIES

Pace Analytical Services performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Pace Analytical Services shall be to re-perform the work at it's own expense. In no event shall Pace Analytical Services be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Pace Analytical Services.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



**Pace Analytical Services LLC**Facility: **Northeast**Department: **Quality Assurance**Title: **Certificate/Approval Program Summary**ID No.: **17873**

Revision 27

Published Date: 01/24/2025

Page 1 of 2

**Certification Information**

The following analytes are not included in our Primary NELAP Scope of Accreditation:

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****EPA 624.1:** m/p-xylene, o-xylene, Naphthalene**EPA 625.1:** alpha-Terpineol**EPA 8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****SM 2540D:** TSS.**EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**MADEP-APH.****Nonpotable Water:** EPA RSK-175 Dissolved Gases**Biological Tissue Matrix:** EPA 3050B**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048****EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**Nonpotable Water:** EPA RSK-175 Dissolved Gases

The following test method is not included in our New Jersey Secondary NELAP Scope of Accreditation:

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Determination of Selected Perfluorinated Alkyl Substances by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry Isotope Dilution (via Alpha SOP 23528)**

The following analytes are included in our Massachusetts DEP Scope of Accreditation

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****Drinking Water****EPA 300.0:** Chloride, Nitrate-N, Fluoride, Sulfate; **EPA 353.2:** Nitrate-N, Nitrite-N; **SM4500NO3-F:** Nitrate-N, Nitrite-N; **SM4500F-C, SM4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B****EPA 524.2:** THMs and VOCs; **EPA 504.1:** EDB, DBCP.**Microbiology:** SM9215B; SM9223-P/A, SM9223B-Colilert-QT, SM9222D.**Non-Potable Water****SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH:** Ammonia-N and Kjeldahl-N, **EPA 350.1:** Ammonia-N, **LACHAT 10-107-06-1-B:** Ammonia-N, **EPA 351.1, SM4500NO3-F, EPA 353.2:** Nitrate-N, **SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300:** Chloride, Sulfate, Nitrate.**EPA 624.1:** Volatile Halocarbons & Aromatics,**EPA 608.3:** Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs**EPA 625.1:** SVOC (Acid/Base/Neutral Extractables).**Microbiology:** SM9223B-Colilert-QT; Enterolert-QT, EPA 1600, EPA 1603, SM9222D.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Drinking Water****EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1 Hg. EPA 522, EPA 537.1.****Non-Potable Water****EPA 200.7:** Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.**EPA 200.8:** Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.**EPA 245.1 Hg.****SM2340B**

**Pace Analytical Services LLC**ID No.: **17873**Facility: **Northeast**

Revision 27

Department: **Quality Assurance**

Published Date: 01/24/2025

Title: **Certificate/Approval Program Summary**

Page 2 of 2

**Certification IDs:****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581**

CT PH-0826, IL 200077, IN C-MA-03, KY JY98045, ME MA00086, MD 348, MA M-MA086, NH 2064, NJ MA935, NY 11148, NC (DW) 25700, NC (NPW/SCM) 666, OR MA-1316, PA 68-03671, RI LAO00065, TX T104704476, VT VT-0935, VA 460195

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048**

CT PH-0825, ANAB/DoD L2474, IL 200081, IN C-MA-04, KY KY98046, LA 3090, ME MA00030, MI 9110, MN 025-999-495, NH 2062, NJ MA015, NY 11627, NC (NPW/SCM) 685, OR MA-0262, PA 68-02089, RI LAO00299, TX T-104704419, VT VT-0015, VA 460194, WA C954

**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048**

ANAB/DoD L2474, ME MA01156, MN 025-999-498, NH 2249, NJ MA025, NY 12191, OR 4203, TX T104704583, VA 460311, WA C1104.

---

For a complete listing of analytes and methods, please contact your Project Manager.





# CHAIN OF CUSTODY

PAGE 1 OF 1

WESTBORO, MA  
TEL: 508-898-8220  
FAX: 508-898-9193

MANSFIELD, MA  
TEL: 508-822-9300  
FAX: 508-822-3288

## Client Information

Client: Tridepoint Atlantic  
Address: 6995 Bethlehem Blvd  
Sparrows Point, MD  
Phone:  
Fax:

Email: J.Barnes@armgroup.net  
K.Guile@armgroup.net  
☐ These samples have been previously analyzed by Pace

Other Project Specific Requirements/Comments/Detection Limits:

## Project Information

Project Name: B14 1st Half 2025 GW  
Project Location: B14  
Project #: 21010214  
Project Manager: Kaye Guile  
Pace Quote #:

## Turn-Around Time

☒ Standard ☐ RUSH (only confirmed if pre-approved!)

Date Due: (5-day) Time:

Date Rec'd in Lab: 4/15/25Pace Job #: L2522789

## Report Information - Data Deliverables

☐ FAX ☐ EMAIL  
☐ ADEx ☐ Add'l Deliverables

## Billing Information

☐ Same as Client info PO #:

## Regulatory Requirements/Report Limits

State /Fed Program Criteria

| ANALYSIS                    |   |  |  |  |  |  |  |  |  | SAMPLE HANDLING                                                                                                                                                                                                |   | TOTAL # BOTTLES |
|-----------------------------|---|--|--|--|--|--|--|--|--|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|---|-----------------|
| VOCs 8260<br>SVOCs 8270 Sim |   |  |  |  |  |  |  |  |  | Filtration _____<br><input type="checkbox"/> Done<br><input type="checkbox"/> Not needed<br><input type="checkbox"/> Lab to do<br>Preservation<br><input type="checkbox"/> Lab to do<br>(Please specify below) |   |                 |
| Sample Specific Comments    |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
| X                           | X |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                | 5 |                 |
| X                           | X |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                | 5 |                 |
| X                           | X |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                | 5 |                 |
| X                           | X |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                | 5 |                 |
| X                           |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                | 4 |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |
|                             |   |  |  |  |  |  |  |  |  |                                                                                                                                                                                                                |   |                 |





## ANALYTICAL REPORT

|                 |                                                                             |
|-----------------|-----------------------------------------------------------------------------|
| Lab Number:     | L2522156                                                                    |
| Client:         | Tradepoint Atlantic<br>1600 Sparrows Point Boulevard<br>Baltimore, MD 21219 |
| ATTN:           | Matthew Newman                                                              |
| Phone:          | (443) 649-5073                                                              |
| Project Name:   | B14 1ST HALF 2025 GW                                                        |
| Project Number: | 21010214                                                                    |
| Report Date:    | 04/23/25                                                                    |

The original project report/data package is held by Pace Analytical Services. This report/data package is paginated and should be reproduced only in its entirety. Pace Analytical Services holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0826), IL (200077), IN (C-MA-03), KY (KY98045), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), OR (MA-1316), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #525-23-122-91930A1).

---

Eight Walkup Drive, Westborough, MA 01581-1019  
508-898-9220 (Fax) 508-898-9193 800-624-9220 - [www.alphalab.com](http://www.alphalab.com)



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

| Lab<br>Sample ID | Client ID | Matrix | Sample<br>Location | Collection<br>Date/Time | Receive Date |
|------------------|-----------|--------|--------------------|-------------------------|--------------|
| L2522156-01      | HI21-MWS  | WATER  | B14                | 04/10/25 10:30          | 04/10/25     |
| L2522156-02      | HI19-MWS  | WATER  | B14                | 04/10/25 11:25          | 04/10/25     |
| L2522156-03      | HI16-MWS  | WATER  | B14                | 04/10/25 13:05          | 04/10/25     |
| L2522156-04      | HI15-MWS  | WATER  | B14                | 04/10/25 13:50          | 04/10/25     |
| L2522156-05      | HI14-MWS  | WATER  | B14                | 04/10/25 14:25          | 04/10/25     |
| L2522156-06      | TB-001    | WATER  | B14                | 04/10/25 00:00          | 04/10/25     |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

### Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Pace Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments and solids are reported on a dry weight basis unless otherwise noted. Tissues are reported "as received" or on a wet weight basis, unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

**HOLD POLICY** - For samples submitted on hold, Pace's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Pace Project Manager and made arrangements for Pace to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.

---

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

### Case Narrative (continued)

#### Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

#### Semivolatile Organics

L2522156-05: The volume of sample received for the analysis deviates from the laboratory's current volume requirements. An aliquot was taken from the original sample container, extracted, analyzed, and reported accordingly.

#### Semivolatile Organics by SIM

L2522156-05: The volume of sample received for the analysis deviates from the laboratory's current volume requirements. An aliquot was taken from the original sample container, extracted, analyzed, and reported accordingly.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

*Melissa Sturgis* Melissa Sturgis

Title: Technical Director/Representative

Date: 04/23/25

# ORGANICS

# VOLATILES

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-01  
**Client ID:** HI21-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 10:30  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/18/25 09:44  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | ND     |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS****Lab ID:** L2522156-01**Date Collected:** 04/10/25 10:30**Client ID:** HI21-MWS**Date Received:** 04/10/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 121        |           | 70-130              |
| Toluene-d8            | 98         |           | 70-130              |
| 4-Bromofluorobenzene  | 97         |           | 70-130              |
| Dibromofluoromethane  | 108        |           | 70-130              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

Lab ID: L2522156-01  
 Client ID: HI21-MWS  
 Sample Location: B14

Date Collected: 04/10/25 10:30  
 Date Received: 04/10/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/18/25 09:44  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 120        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 99         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-02  
**Client ID:** HI19-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 11:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/18/25 10:37  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 2.7    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 12     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | 4.1    |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS****Lab ID:** L2522156-02**Date Collected:** 04/10/25 11:25**Client ID:** HI19-MWS**Date Received:** 04/10/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | 0.19   | J         | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | 2.0    |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | 1.0    |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | 3.0    |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 119        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 94         |           | 70-130              |
| Dibromofluoromethane  | 107        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-02  
**Client ID:** HI19-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 11:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/18/25 10:37  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 117        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 97         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-03  
**Client ID:** HI16-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:05  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/18/25 10:11  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 7.4    |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-03  
**Client ID:** HI16-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:05  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 120        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 109        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-03  
**Client ID:** HI16-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:05  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/18/25 10:11  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 118        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 99         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-04  
**Client ID:** HI15-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:50  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/18/25 11:04  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 3.2    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |





**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS****Lab ID:** L2522156-04**Date Collected:** 04/10/25 13:50**Client ID:** HI15-MWS**Date Received:** 04/10/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 119        |           | 70-130              |
| Toluene-d8            | 98         |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 111        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

Lab ID: L2522156-04  
 Client ID: HI15-MWS  
 Sample Location: B14

Date Collected: 04/10/25 13:50  
 Date Received: 04/10/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/18/25 11:04  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 118        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 99         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-05  
**Client ID:** HI14-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 14:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/18/25 11:30  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 2.2    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS****Lab ID:** L2522156-05**Date Collected:** 04/10/25 14:25**Client ID:** HI14-MWS**Date Received:** 04/10/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 120        |           | 70-130              |
| Toluene-d8            | 100        |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 109        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-05  
**Client ID:** HI14-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 14:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/18/25 11:30  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 119        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 99         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-06  
**Client ID:** TB-001  
**Sample Location:** B14

**Date Collected:** 04/10/25 00:00  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/17/25 13:10  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | ND     |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS****Lab ID:** L2522156-06**Date Collected:** 04/10/25 00:00**Client ID:** TB-001**Date Received:** 04/10/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,1,2,2-Tetrachloroethane                    | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 120        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 97         |           | 70-130              |
| Dibromofluoromethane  | 106        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

Lab ID: L2522156-06  
 Client ID: TB-001  
 Sample Location: B14

Date Collected: 04/10/25 00:00  
 Date Received: 04/10/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/17/25 13:10  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 116        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 04/17/25 06:07  
 Analyst: MCM

| Parameter                                                                         | Result | Qualifier | Units | RL   | MDL  |
|-----------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 06 Batch: WG2055118-5 |        |           |       |      |      |
| Dichlorodifluoromethane                                                           | ND     |           | ug/l  | 5.0  | 0.24 |
| Chloromethane                                                                     | ND     |           | ug/l  | 2.5  | 0.20 |
| Vinyl chloride                                                                    | ND     |           | ug/l  | 1.0  | 0.07 |
| Bromomethane                                                                      | ND     |           | ug/l  | 1.0  | 0.26 |
| Chloroethane                                                                      | ND     |           | ug/l  | 1.0  | 0.13 |
| Trichlorofluoromethane                                                            | ND     |           | ug/l  | 2.5  | 0.16 |
| 1,1-Dichloroethene                                                                | ND     |           | ug/l  | 0.50 | 0.17 |
| Carbon disulfide                                                                  | ND     |           | ug/l  | 5.0  | 0.30 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                             | ND     |           | ug/l  | 2.5  | 0.15 |
| Methylene chloride                                                                | ND     |           | ug/l  | 2.5  | 0.68 |
| Acetone                                                                           | ND     |           | ug/l  | 5.0  | 1.5  |
| trans-1,2-Dichloroethene                                                          | ND     |           | ug/l  | 0.75 | 0.16 |
| Methyl Acetate                                                                    | ND     |           | ug/l  | 2.0  | 0.23 |
| Methyl tert butyl ether                                                           | ND     |           | ug/l  | 1.0  | 0.17 |
| 1,1-Dichloroethane                                                                | ND     |           | ug/l  | 0.75 | 0.21 |
| cis-1,2-Dichloroethene                                                            | ND     |           | ug/l  | 0.50 | 0.19 |
| Cyclohexane                                                                       | ND     |           | ug/l  | 10   | 0.27 |
| Chloroform                                                                        | ND     |           | ug/l  | 0.75 | 0.22 |
| Carbon tetrachloride                                                              | ND     |           | ug/l  | 0.50 | 0.13 |
| 1,1,1-Trichloroethane                                                             | ND     |           | ug/l  | 0.50 | 0.16 |
| 2-Butanone                                                                        | ND     |           | ug/l  | 5.0  | 1.9  |
| Benzene                                                                           | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,2-Dichloroethane                                                                | ND     |           | ug/l  | 0.50 | 0.13 |
| Trichloroethene                                                                   | ND     |           | ug/l  | 0.50 | 0.18 |
| 1,2-Dichloropropane                                                               | ND     |           | ug/l  | 1.0  | 0.14 |
| Bromodichloromethane                                                              | ND     |           | ug/l  | 0.50 | 0.19 |
| cis-1,3-Dichloropropene                                                           | ND     |           | ug/l  | 0.50 | 0.14 |
| Toluene                                                                           | ND     |           | ug/l  | 0.75 | 0.20 |
| Tetrachloroethene                                                                 | ND     |           | ug/l  | 0.50 | 0.18 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 04/17/25 06:07  
 Analyst: MCM

| Parameter                                                                         | Result | Qualifier | Units | RL   | MDL  |
|-----------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 06 Batch: WG2055118-5 |        |           |       |      |      |
| 4-Methyl-2-pentanone                                                              | ND     |           | ug/l  | 5.0  | 0.42 |
| trans-1,3-Dichloropropene                                                         | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,3-Dichloropropene, Total                                                        | ND     |           | ug/l  | 0.50 | 0.14 |
| 1,1,2-Trichloroethane                                                             | ND     |           | ug/l  | 0.75 | 0.14 |
| Dibromochloromethane                                                              | ND     |           | ug/l  | 0.50 | 0.15 |
| 1,2-Dibromoethane                                                                 | ND     |           | ug/l  | 2.0  | 0.19 |
| 2-Hexanone                                                                        | ND     |           | ug/l  | 5.0  | 0.52 |
| Chlorobenzene                                                                     | ND     |           | ug/l  | 0.50 | 0.18 |
| Ethylbenzene                                                                      | ND     |           | ug/l  | 0.50 | 0.17 |
| p/m-Xylene                                                                        | ND     |           | ug/l  | 1.0  | 0.33 |
| o-Xylene                                                                          | ND     |           | ug/l  | 1.0  | 0.39 |
| Xylenes, Total                                                                    | ND     |           | ug/l  | 1.0  | 0.33 |
| Styrene                                                                           | ND     |           | ug/l  | 1.0  | 0.36 |
| Bromoform                                                                         | ND     |           | ug/l  | 2.0  | 0.25 |
| Isopropylbenzene                                                                  | ND     |           | ug/l  | 0.50 | 0.19 |
| 1,1,2,2-Tetrachloroethane                                                         | ND     |           | ug/l  | 0.50 | 0.17 |
| 1,3-Dichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,4-Dichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,2-Dichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.18 |
| 1,2-Dibromo-3-chloropropane                                                       | ND     |           | ug/l  | 2.5  | 0.35 |
| 1,2,4-Trichlorobenzene                                                            | ND     |           | ug/l  | 2.5  | 0.22 |
| 1,2,3-Trichlorobenzene                                                            | ND     |           | ug/l  | 2.5  | 0.23 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
Analytical Date: 04/17/25 06:07  
Analyst: MCM

| Parameter                                                                         | Result | Qualifier | Units | RL | MDL |
|-----------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 06 Batch: WG2055118-5 |        |           |       |    |     |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 119       |           | 70-130                 |
| Toluene-d8            | 99        |           | 70-130                 |
| 4-Bromofluorobenzene  | 98        |           | 70-130                 |
| Dibromofluoromethane  | 106       |           | 70-130                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D-SIM(M)  
Analytical Date: 04/17/25 06:07  
Analyst: MCM

| Parameter                                                                             | Result | Qualifier | Units | RL    | MDL   |
|---------------------------------------------------------------------------------------|--------|-----------|-------|-------|-------|
| Volatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 06 Batch: WG2055119-5 |        |           |       |       |       |
| 1,1,1,2-Tetrachloroethane                                                             | ND     |           | ug/l  | 0.050 | 0.006 |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 115       |           | 70-130                 |
| 4-Bromofluorobenzene  | 96        |           | 70-130                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 04/18/25 06:13  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2056423-5 |        |           |       |      |      |
| Dichlorodifluoromethane                                                              | ND     |           | ug/l  | 5.0  | 0.24 |
| Chloromethane                                                                        | ND     |           | ug/l  | 2.5  | 0.20 |
| Vinyl chloride                                                                       | ND     |           | ug/l  | 1.0  | 0.07 |
| Bromomethane                                                                         | ND     |           | ug/l  | 1.0  | 0.26 |
| Chloroethane                                                                         | ND     |           | ug/l  | 1.0  | 0.13 |
| Trichlorofluoromethane                                                               | ND     |           | ug/l  | 2.5  | 0.16 |
| 1,1-Dichloroethene                                                                   | ND     |           | ug/l  | 0.50 | 0.17 |
| Carbon disulfide                                                                     | ND     |           | ug/l  | 5.0  | 0.30 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                | ND     |           | ug/l  | 2.5  | 0.15 |
| Methylene chloride                                                                   | ND     |           | ug/l  | 2.5  | 0.68 |
| Acetone                                                                              | ND     |           | ug/l  | 5.0  | 1.5  |
| trans-1,2-Dichloroethene                                                             | ND     |           | ug/l  | 0.75 | 0.16 |
| Methyl Acetate                                                                       | ND     |           | ug/l  | 2.0  | 0.23 |
| Methyl tert butyl ether                                                              | ND     |           | ug/l  | 1.0  | 0.17 |
| 1,1-Dichloroethane                                                                   | ND     |           | ug/l  | 0.75 | 0.21 |
| cis-1,2-Dichloroethene                                                               | ND     |           | ug/l  | 0.50 | 0.19 |
| Cyclohexane                                                                          | ND     |           | ug/l  | 10   | 0.27 |
| Chloroform                                                                           | ND     |           | ug/l  | 0.75 | 0.22 |
| Carbon tetrachloride                                                                 | ND     |           | ug/l  | 0.50 | 0.13 |
| 1,1,1-Trichloroethane                                                                | ND     |           | ug/l  | 0.50 | 0.16 |
| 2-Butanone                                                                           | ND     |           | ug/l  | 5.0  | 1.9  |
| Benzene                                                                              | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,2-Dichloroethane                                                                   | ND     |           | ug/l  | 0.50 | 0.13 |
| Trichloroethene                                                                      | ND     |           | ug/l  | 0.50 | 0.18 |
| 1,2-Dichloropropane                                                                  | ND     |           | ug/l  | 1.0  | 0.14 |
| Bromodichloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.19 |
| cis-1,3-Dichloropropene                                                              | ND     |           | ug/l  | 0.50 | 0.14 |
| Toluene                                                                              | ND     |           | ug/l  | 0.75 | 0.20 |
| Tetrachloroethene                                                                    | ND     |           | ug/l  | 0.50 | 0.18 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8260D  
**Analytical Date:** 04/18/25 06:13  
**Analyst:** MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2056423-5 |        |           |       |      |      |
| 4-Methyl-2-pentanone                                                                 | ND     |           | ug/l  | 5.0  | 0.42 |
| trans-1,3-Dichloropropene                                                            | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,3-Dichloropropene, Total                                                           | ND     |           | ug/l  | 0.50 | 0.14 |
| 1,1,2-Trichloroethane                                                                | ND     |           | ug/l  | 0.75 | 0.14 |
| Dibromochloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.15 |
| 1,2-Dibromoethane                                                                    | ND     |           | ug/l  | 2.0  | 0.19 |
| 2-Hexanone                                                                           | ND     |           | ug/l  | 5.0  | 0.52 |
| Chlorobenzene                                                                        | ND     |           | ug/l  | 0.50 | 0.18 |
| Ethylbenzene                                                                         | ND     |           | ug/l  | 0.50 | 0.17 |
| p/m-Xylene                                                                           | ND     |           | ug/l  | 1.0  | 0.33 |
| o-Xylene                                                                             | ND     |           | ug/l  | 1.0  | 0.39 |
| Xylenes, Total                                                                       | ND     |           | ug/l  | 1.0  | 0.33 |
| Styrene                                                                              | ND     |           | ug/l  | 1.0  | 0.36 |
| Bromoform                                                                            | ND     |           | ug/l  | 2.0  | 0.25 |
| Isopropylbenzene                                                                     | ND     |           | ug/l  | 0.50 | 0.19 |
| 1,1,2,2-Tetrachloroethane                                                            | ND     |           | ug/l  | 0.50 | 0.17 |
| 1,3-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,4-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,2-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.18 |
| 1,2-Dibromo-3-chloropropane                                                          | ND     |           | ug/l  | 2.5  | 0.35 |
| 1,2,4-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.22 |
| 1,2,3-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.23 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
Analytical Date: 04/18/25 06:13  
Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL | MDL |
|--------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2056423-5 |        |           |       |    |     |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 120       |           | 70-130                 |
| Toluene-d8            | 98        |           | 70-130                 |
| 4-Bromofluorobenzene  | 98        |           | 70-130                 |
| Dibromofluoromethane  | 108       |           | 70-130                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D-SIM(M)  
Analytical Date: 04/18/25 06:13  
Analyst: MCM

| Parameter                                                                                | Result | Qualifier | Units | RL    | MDL   |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-------|-------|
| Volatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-05 Batch: WG2056429-5 |        |           |       |       |       |
| 1,1,1,2-Tetrachloroethane                                                                | ND     |           | ug/l  | 0.050 | 0.006 |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 116       |           | 70-130                 |
| 4-Bromofluorobenzene  | 94        |           | 70-130                 |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522156

**Report Date:** 04/23/25

| Parameter                                                                                            | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 06 Batch: WG2055118-3 WG2055118-4 |                  |      |                   |      |                     |     |      |               |
| Dichlorodifluoromethane                                                                              | 110              |      | 100               |      | 36-147              | 10  |      | 20            |
| Chloromethane                                                                                        | 97               |      | 95                |      | 64-130              | 2   |      | 20            |
| Vinyl chloride                                                                                       | 110              |      | 100               |      | 55-140              | 10  |      | 20            |
| Bromomethane                                                                                         | 64               |      | 66                |      | 39-139              | 3   |      | 20            |
| Chloroethane                                                                                         | 120              |      | 120               |      | 55-138              | 0   |      | 20            |
| Trichlorofluoromethane                                                                               | 110              |      | 110               |      | 62-150              | 0   |      | 20            |
| 1,1-Dichloroethene                                                                                   | 100              |      | 98                |      | 61-145              | 2   |      | 20            |
| Carbon disulfide                                                                                     | 110              |      | 100               |      | 51-130              | 10  |      | 20            |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                                | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Methylene chloride                                                                                   | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| Acetone                                                                                              | 120              |      | 110               |      | 58-148              | 9   |      | 20            |
| trans-1,2-Dichloroethene                                                                             | 100              |      | 97                |      | 70-130              | 3   |      | 20            |
| Methyl Acetate                                                                                       | 120              |      | 120               |      | 70-130              | 0   |      | 20            |
| Methyl tert butyl ether                                                                              | 100              |      | 99                |      | 63-130              | 1   |      | 20            |
| 1,1-Dichloroethane                                                                                   | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| cis-1,2-Dichloroethene                                                                               | 98               |      | 96                |      | 70-130              | 2   |      | 20            |
| Cyclohexane                                                                                          | 120              |      | 110               |      | 70-130              | 9   |      | 20            |
| Chloroform                                                                                           | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Carbon tetrachloride                                                                                 | 100              |      | 100               |      | 63-132              | 0   |      | 20            |
| 1,1,1-Trichloroethane                                                                                | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| 2-Butanone                                                                                           | 96               |      | 93                |      | 63-138              | 3   |      | 20            |
| Benzene                                                                                              | 100              |      | 96                |      | 70-130              | 4   |      | 20            |
| 1,2-Dichloroethane                                                                                   | 110              |      | 110               |      | 70-130              | 0   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522156

**Report Date:** 04/23/25

| Parameter                                                                                            | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 06 Batch: WG2055118-3 WG2055118-4 |                  |      |                   |      |                     |     |      |               |
| Trichloroethene                                                                                      | 98               |      | 94                |      | 70-130              | 4   |      | 20            |
| 1,2-Dichloropropane                                                                                  | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Bromodichloromethane                                                                                 | 99               |      | 95                |      | 67-130              | 4   |      | 20            |
| cis-1,3-Dichloropropene                                                                              | 94               |      | 92                |      | 70-130              | 2   |      | 20            |
| Toluene                                                                                              | 98               |      | 96                |      | 70-130              | 2   |      | 20            |
| Tetrachloroethene                                                                                    | 99               |      | 95                |      | 70-130              | 4   |      | 20            |
| 4-Methyl-2-pentanone                                                                                 | 89               |      | 88                |      | 59-130              | 1   |      | 20            |
| trans-1,3-Dichloropropene                                                                            | 99               |      | 99                |      | 70-130              | 0   |      | 20            |
| 1,1,2-Trichloroethane                                                                                | 100              |      | 98                |      | 70-130              | 2   |      | 20            |
| Dibromochloromethane                                                                                 | 92               |      | 92                |      | 63-130              | 0   |      | 20            |
| 1,2-Dibromoethane                                                                                    | 96               |      | 94                |      | 70-130              | 2   |      | 20            |
| 2-Hexanone                                                                                           | 95               |      | 91                |      | 57-130              | 4   |      | 20            |
| Chlorobenzene                                                                                        | 96               |      | 94                |      | 75-130              | 2   |      | 20            |
| Ethylbenzene                                                                                         | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| p/m-Xylene                                                                                           | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| o-Xylene                                                                                             | 100              |      | 95                |      | 70-130              | 5   |      | 20            |
| Styrene                                                                                              | 95               |      | 95                |      | 70-130              | 0   |      | 20            |
| Bromoform                                                                                            | 89               |      | 88                |      | 54-136              | 1   |      | 20            |
| Isopropylbenzene                                                                                     | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| 1,1,2,2-Tetrachloroethane                                                                            | 100              |      | 96                |      | 67-130              | 4   |      | 20            |
| 1,3-Dichlorobenzene                                                                                  | 100              |      | 97                |      | 70-130              | 3   |      | 20            |
| 1,4-Dichlorobenzene                                                                                  | 99               |      | 95                |      | 70-130              | 4   |      | 20            |
| 1,2-Dichlorobenzene                                                                                  | 98               |      | 96                |      | 70-130              | 2   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522156

**Project Number:** 21010214

**Report Date:** 04/23/25

| <b>Parameter</b>                                                                                     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|------------------------------------------------------------------------------------------------------|--------------------------|-------------|--------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 06 Batch: WG2055118-3 WG2055118-4 |                          |             |                          |             |                             |            |             |                       |
| 1,2-Dibromo-3-chloropropane                                                                          | 82                       |             | 80                       |             | 41-144                      | 2          |             | 20                    |
| 1,2,4-Trichlorobenzene                                                                               | 94                       |             | 89                       |             | 70-130                      | 5          |             | 20                    |
| 1,2,3-Trichlorobenzene                                                                               | 92                       |             | 91                       |             | 70-130                      | 1          |             | 20                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|--------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 115                      |             | 112                      |             | 70-130                         |
| Toluene-d8            | 101                      |             | 100                      |             | 70-130                         |
| 4-Bromofluorobenzene  | 102                      |             | 100                      |             | 70-130                         |
| Dibromofluoromethane  | 103                      |             | 103                      |             | 70-130                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522156

**Project Number:** 21010214

**Report Date:** 04/23/25

| <b>Parameter</b>                                                                                         | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|----------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 06 Batch: WG2055119-3 WG2055119-4 |                          |             |                           |             |                             |            |             |                       |
| 1,1,2,2-Tetrachloroethane                                                                                | 100                      |             | 96                        |             | 70-130                      | 4          |             | 25                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 114                      |             | 115                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 98                       |             | 97                        |             | 70-130                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522156

**Report Date:** 04/23/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2056423-3 WG2056423-4 |                  |      |                   |      |                     |     |      |               |
| Dichlorodifluoromethane                                                                                 | 110              |      | 110               |      | 36-147              | 0   |      | 20            |
| Chloromethane                                                                                           | 110              |      | 110               |      | 64-130              | 0   |      | 20            |
| Vinyl chloride                                                                                          | 110              |      | 110               |      | 55-140              | 0   |      | 20            |
| Bromomethane                                                                                            | 58               |      | 62                |      | 39-139              | 7   |      | 20            |
| Chloroethane                                                                                            | 120              |      | 120               |      | 55-138              | 0   |      | 20            |
| Trichlorofluoromethane                                                                                  | 110              |      | 110               |      | 62-150              | 0   |      | 20            |
| 1,1-Dichloroethene                                                                                      | 100              |      | 100               |      | 61-145              | 0   |      | 20            |
| Carbon disulfide                                                                                        | 110              |      | 110               |      | 51-130              | 0   |      | 20            |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                                   | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Methylene chloride                                                                                      | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Acetone                                                                                                 | 120              |      | 120               |      | 58-148              | 0   |      | 20            |
| trans-1,2-Dichloroethene                                                                                | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Methyl Acetate                                                                                          | 130              |      | 130               |      | 70-130              | 0   |      | 20            |
| Methyl tert butyl ether                                                                                 | 98               |      | 99                |      | 63-130              | 1   |      | 20            |
| 1,1-Dichloroethane                                                                                      | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| cis-1,2-Dichloroethene                                                                                  | 98               |      | 98                |      | 70-130              | 0   |      | 20            |
| Cyclohexane                                                                                             | 120              |      | 120               |      | 70-130              | 0   |      | 20            |
| Chloroform                                                                                              | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Carbon tetrachloride                                                                                    | 100              |      | 100               |      | 63-132              | 0   |      | 20            |
| 1,1,1-Trichloroethane                                                                                   | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| 2-Butanone                                                                                              | 100              |      | 100               |      | 63-138              | 0   |      | 20            |
| Benzene                                                                                                 | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| 1,2-Dichloroethane                                                                                      | 110              |      | 110               |      | 70-130              | 0   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522156

**Report Date:** 04/23/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2056423-3 WG2056423-4 |                  |      |                   |      |                     |     |      |               |
| Trichloroethene                                                                                         | 98               |      | 97                |      | 70-130              | 1   |      | 20            |
| 1,2-Dichloropropane                                                                                     | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Bromodichloromethane                                                                                    | 98               |      | 97                |      | 67-130              | 1   |      | 20            |
| cis-1,3-Dichloropropene                                                                                 | 92               |      | 92                |      | 70-130              | 0   |      | 20            |
| Toluene                                                                                                 | 98               |      | 98                |      | 70-130              | 0   |      | 20            |
| Tetrachloroethene                                                                                       | 97               |      | 97                |      | 70-130              | 0   |      | 20            |
| 4-Methyl-2-pentanone                                                                                    | 87               |      | 90                |      | 59-130              | 3   |      | 20            |
| trans-1,3-Dichloropropene                                                                               | 97               |      | 99                |      | 70-130              | 2   |      | 20            |
| 1,1,2-Trichloroethane                                                                                   | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Dibromochloromethane                                                                                    | 92               |      | 92                |      | 63-130              | 0   |      | 20            |
| 1,2-Dibromoethane                                                                                       | 95               |      | 95                |      | 70-130              | 0   |      | 20            |
| 2-Hexanone                                                                                              | 93               |      | 96                |      | 57-130              | 3   |      | 20            |
| Chlorobenzene                                                                                           | 96               |      | 96                |      | 75-130              | 0   |      | 20            |
| Ethylbenzene                                                                                            | 99               |      | 100               |      | 70-130              | 1   |      | 20            |
| p/m-Xylene                                                                                              | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| o-Xylene                                                                                                | 95               |      | 95                |      | 70-130              | 0   |      | 20            |
| Styrene                                                                                                 | 95               |      | 95                |      | 70-130              | 0   |      | 20            |
| Bromoform                                                                                               | 86               |      | 86                |      | 54-136              | 0   |      | 20            |
| Isopropylbenzene                                                                                        | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| 1,1,2,2-Tetrachloroethane                                                                               | 98               |      | 100               |      | 67-130              | 2   |      | 20            |
| 1,3-Dichlorobenzene                                                                                     | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| 1,4-Dichlorobenzene                                                                                     | 98               |      | 98                |      | 70-130              | 0   |      | 20            |
| 1,2-Dichlorobenzene                                                                                     | 95               |      | 98                |      | 70-130              | 3   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522156

**Project Number:** 21010214

**Report Date:** 04/23/25

| <b>Parameter</b>                                                                                        | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|---------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2056423-3 WG2056423-4 |                          |             |                           |             |                             |            |             |                       |
| 1,2-Dibromo-3-chloropropane                                                                             | 81                       |             | 87                        |             | 41-144                      | 7          |             | 20                    |
| 1,2,4-Trichlorobenzene                                                                                  | 90                       |             | 93                        |             | 70-130                      | 3          |             | 20                    |
| 1,2,3-Trichlorobenzene                                                                                  | 93                       |             | 94                        |             | 70-130                      | 1          |             | 20                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 115                      |             | 112                       |             | 70-130                         |
| Toluene-d8            | 100                      |             | 100                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 95                       |             | 99                        |             | 70-130                         |
| Dibromofluoromethane  | 102                      |             | 101                       |             | 70-130                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522156

**Project Number:** 21010214

**Report Date:** 04/23/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-05 Batch: WG2056429-3 WG2056429-4 |                          |             |                           |             |                             |            |             |                       |
| 1,1,1,2-Tetrachloroethane                                                                                   | 104                      |             | 96                        |             | 70-130                      | 8          |             | 25                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 114                      |             | 116                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 94                       |             | 92                        |             | 70-130                         |



# SEMIVOLATILES

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-01  
**Client ID:** HI21-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 10:30  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 22:21  
**Analyst:** ALS

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 18:56

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS**

Lab ID: L2522156-01

Date Collected: 04/10/25 10:30

Client ID: HI21-MWS

Date Received: 04/10/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 48         |           | 21-120              |
| Phenol-d6            | 37         |           | 10-120              |
| Nitrobenzene-d5      | 72         |           | 23-120              |
| 2-Fluorobiphenyl     | 66         |           | 15-120              |
| 2,4,6-Tribromophenol | 67         |           | 10-120              |
| 4-Terphenyl-d14      | 73         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-01  
**Client ID:** HI21-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 10:30  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/13/25 14:01  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 18:56

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.05   | J         | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 123        | Q         | 23-120              |
| 2-Fluorobiphenyl | 90         |           | 15-120              |
| 4-Terphenyl-d14  | 97         |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-02  
**Client ID:** HI19-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 11:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 22:44  
**Analyst:** ALS

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 18:56

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | 4.2    | J         | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | 3.2    | J         | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | 57     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | 150    |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 44     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | 3.1    |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | 0.76   | J         | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS**

Lab ID: L2522156-02

Date Collected: 04/10/25 11:25

Client ID: HI19-MWS

Date Received: 04/10/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | 1.4    | J         | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | 1.7    | J         | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 51         |           | 21-120              |
| Phenol-d6            | 40         |           | 10-120              |
| Nitrobenzene-d5      | 83         |           | 23-120              |
| 2-Fluorobiphenyl     | 76         |           | 15-120              |
| 2,4,6-Tribromophenol | 81         |           | 10-120              |
| 4-Terphenyl-d14      | 81         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-02      **D**  
**Client ID:** HI19-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 11:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/15/25 11:55  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 18:56

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 52     |           | ug/l  | 0.20 | 0.05 | 2               |
| 2-Methylnaphthalene                                  | 3.7    |           | ug/l  | 0.20 | 0.06 | 2               |
| Acenaphthylene                                       | 0.56   |           | ug/l  | 0.20 | 0.04 | 2               |
| Acenaphthene                                         | 0.90   |           | ug/l  | 0.20 | 0.05 | 2               |
| Fluorene                                             | 1.0    |           | ug/l  | 0.20 | 0.05 | 2               |
| Phenanthrene                                         | 1.6    |           | ug/l  | 0.10 | 0.08 | 2               |
| Anthracene                                           | 0.30   |           | ug/l  | 0.20 | 0.05 | 2               |
| Fluoranthene                                         | 0.37   |           | ug/l  | 0.20 | 0.05 | 2               |
| Pyrene                                               | 0.24   |           | ug/l  | 0.20 | 0.09 | 2               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.10 | 0.06 | 2               |
| Chrysene                                             | ND     |           | ug/l  | 0.20 | 0.06 | 2               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.05 | 2               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.20 | 0.07 | 2               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.20 | 0.05 | 2               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.20 | 0.04 | 2               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.10 | 0.05 | 2               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.20 | 0.05 | 2               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 117        |           | 23-120              |
| 2-Fluorobiphenyl | 100        |           | 15-120              |
| 4-Terphenyl-d14  | 101        |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-03  
**Client ID:** HI16-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:05  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 11:05  
**Analyst:** JG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 21:31

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 1.6    | J         | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |





**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS**

Lab ID: L2522156-03

Date Collected: 04/10/25 13:05

Client ID: HI16-MWS

Date Received: 04/10/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | 1.6    | J         | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 70         |           | 21-120              |
| Phenol-d6            | 53         |           | 10-120              |
| Nitrobenzene-d5      | 68         |           | 23-120              |
| 2-Fluorobiphenyl     | 92         |           | 15-120              |
| 2,4,6-Tribromophenol | 74         |           | 10-120              |
| 4-Terphenyl-d14      | 104        |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-03  
**Client ID:** HI16-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:05  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/13/25 14:34  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 21:31

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 1.6    |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.08   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.08   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.06   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.05   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.35   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.24   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.23   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | 0.06   |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | 0.07   |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | 0.03   | J         | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 124        | Q         | 23-120              |
| 2-Fluorobiphenyl | 94         |           | 15-120              |
| 4-Terphenyl-d14  | 117        |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-04  
**Client ID:** HI15-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:50  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 11:28  
**Analyst:** JG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 21:31

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 0.65   | J         | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS**

Lab ID: L2522156-04

Date Collected: 04/10/25 13:50

Client ID: HI15-MWS

Date Received: 04/10/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | 8.2    |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | 1.1    | J         | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 53         |           | 21-120              |
| Phenol-d6            | 42         |           | 10-120              |
| Nitrobenzene-d5      | 46         |           | 23-120              |
| 2-Fluorobiphenyl     | 77         |           | 15-120              |
| 2,4,6-Tribromophenol | 44         |           | 10-120              |
| 4-Terphenyl-d14      | 91         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-04  
**Client ID:** HI15-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 13:50  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/13/25 14:50  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 21:31

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.64   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.03   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.06   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 98         |           | 23-120              |
| 2-Fluorobiphenyl | 75         |           | 15-120              |
| 4-Terphenyl-d14  | 101        |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-05  
**Client ID:** HI14-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 14:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/17/25 11:51  
**Analyst:** JG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 21:31

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | 0.80   | J         | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**SAMPLE RESULTS**

Lab ID: L2522156-05

Date Collected: 04/10/25 14:25

Client ID: HI14-MWS

Date Received: 04/10/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | 1.4    | J         | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | 0.81   | J         | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | 0.74   | J         | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | 1.0    | J         | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 74         |           | 21-120              |
| Phenol-d6            | 57         |           | 10-120              |
| Nitrobenzene-d5      | 58         |           | 23-120              |
| 2-Fluorobiphenyl     | 90         |           | 15-120              |
| 2,4,6-Tribromophenol | 54         |           | 10-120              |
| 4-Terphenyl-d14      | 103        |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**SAMPLE RESULTS**

**Lab ID:** L2522156-05  
**Client ID:** HI14-MWS  
**Sample Location:** B14

**Date Collected:** 04/10/25 14:25  
**Date Received:** 04/10/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/13/25 15:06  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 21:31

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.81   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | 0.05   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.05   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 1.6    |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.82   |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.85   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.19   |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.18   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.11   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 125        | Q         | 23-120              |
| 2-Fluorobiphenyl | 91         |           | 15-120              |
| 4-Terphenyl-d14  | 112        |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 18:13  
**Analyst:** LJG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 17:50

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2052805-1 |        |           |       |     |      |
| Benzaldehyde                                                                             | ND     |           | ug/l  | 5.0 | 1.1  |
| Phenol                                                                                   | ND     |           | ug/l  | 5.0 | 0.35 |
| Bis(2-chloroethyl)ether                                                                  | ND     |           | ug/l  | 2.0 | 0.39 |
| 2-Chlorophenol                                                                           | ND     |           | ug/l  | 2.0 | 0.65 |
| 2-Methylphenol                                                                           | ND     |           | ug/l  | 5.0 | 2.3  |
| Bis(2-chloroisopropyl)ether                                                              | ND     |           | ug/l  | 2.0 | 0.40 |
| Acetophenone                                                                             | ND     |           | ug/l  | 5.0 | 0.92 |
| n-Nitrosodi-n-propylamine                                                                | ND     |           | ug/l  | 5.0 | 0.91 |
| 3-Methylphenol/4-Methylphenol                                                            | ND     |           | ug/l  | 5.0 | 1.4  |
| Hexachloroethane                                                                         | ND     |           | ug/l  | 2.0 | 0.20 |
| Nitrobenzene                                                                             | ND     |           | ug/l  | 1.4 | 0.20 |
| Isophorone                                                                               | ND     |           | ug/l  | 5.0 | 0.86 |
| 2,4-Dimethylphenol                                                                       | ND     |           | ug/l  | 5.0 | 2.0  |
| Bis(2-chloroethoxy)methane                                                               | ND     |           | ug/l  | 5.0 | 0.84 |
| 2,4-Dichlorophenol                                                                       | ND     |           | ug/l  | 5.0 | 1.7  |
| Naphthalene                                                                              | ND     |           | ug/l  | 2.0 | 0.54 |
| 4-Chloroaniline                                                                          | ND     |           | ug/l  | 3.7 | 0.47 |
| Hexachlorobutadiene                                                                      | ND     |           | ug/l  | 2.0 | 0.36 |
| Caprolactam                                                                              | ND     |           | ug/l  | 10  | 1.2  |
| 2-Methylnaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.37 |
| Hexachlorocyclopentadiene                                                                | ND     |           | ug/l  | 20  | 1.2  |
| 1,2,4,5-Tetrachlorobenzene                                                               | ND     |           | ug/l  | 1.7 | 0.24 |
| 2,4,6-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| 2,4,5-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| Biphenyl                                                                                 | ND     |           | ug/l  | 2.0 | 0.20 |
| 2-Chloronaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.35 |
| 2-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.0  |
| 2,6-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.84 |
| Acenaphthylene                                                                           | ND     |           | ug/l  | 2.0 | 0.32 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/16/25 18:13  
**Analyst:** LJG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 17:50

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2052805-1 |        |           |       |     |      |
| Acenaphthene                                                                             | ND     |           | ug/l  | 2.0 | 0.40 |
| 2,4-Dinitrophenol                                                                        | ND     |           | ug/l  | 20  | 5.4  |
| 2,4-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.54 |
| 2,3,4,6-Tetrachlorophenol                                                                | ND     |           | ug/l  | 5.0 | 2.2  |
| Diethyl phthalate                                                                        | ND     |           | ug/l  | 5.0 | 0.76 |
| Fluorene                                                                                 | ND     |           | ug/l  | 2.0 | 0.44 |
| 4-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.4  |
| NDPA/DPA                                                                                 | ND     |           | ug/l  | 2.0 | 0.92 |
| Hexachlorobenzene                                                                        | ND     |           | ug/l  | 2.0 | 0.45 |
| Pentachlorophenol                                                                        | ND     |           | ug/l  | 10  | 2.5  |
| Phenanthrene                                                                             | ND     |           | ug/l  | 2.0 | 0.42 |
| Anthracene                                                                               | ND     |           | ug/l  | 2.0 | 0.47 |
| Carbazole                                                                                | ND     |           | ug/l  | 2.0 | 0.31 |
| Di-n-butylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 0.96 |
| Fluoranthene                                                                             | ND     |           | ug/l  | 2.0 | 0.41 |
| Pyrene                                                                                   | ND     |           | ug/l  | 2.0 | 0.41 |
| 3,3'-Dichlorobenzidine                                                                   | ND     |           | ug/l  | 5.0 | 1.8  |
| Benzo(a)anthracene                                                                       | ND     |           | ug/l  | 2.0 | 0.32 |
| Chrysene                                                                                 | ND     |           | ug/l  | 1.5 | 0.22 |
| Bis(2-ethylhexyl)phthalate                                                               | ND     |           | ug/l  | 3.0 | 1.4  |
| Di-n-octylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 2.3  |
| Benzo(b)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.53 |
| Benzo(k)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.62 |
| Benzo(a)pyrene                                                                           | ND     |           | ug/l  | 2.0 | 0.37 |
| Indeno(1,2,3-cd)pyrene                                                                   | ND     |           | ug/l  | 2.0 | 0.48 |
| Dibenzo(a,h)anthracene                                                                   | ND     |           | ug/l  | 2.0 | 0.29 |
| Benzo(ghi)perylene                                                                       | ND     |           | ug/l  | 2.0 | 0.37 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8270E  
Analytical Date: 04/16/25 18:13  
Analyst: LJG

Extraction Method: EPA 3510C  
Extraction Date: 04/11/25 17:50

| Parameter                                                                                | Result | Qualifier | Units | RL | MDL |
|------------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2052805-1 |        |           |       |    |     |

| Surrogate            | %Recovery | Qualifier | Acceptance<br>Criteria |
|----------------------|-----------|-----------|------------------------|
| 2-Fluorophenol       | 37        |           | 21-120                 |
| Phenol-d6            | 26        |           | 10-120                 |
| Nitrobenzene-d5      | 53        |           | 23-120                 |
| 2-Fluorobiphenyl     | 57        |           | 15-120                 |
| 2,4,6-Tribromophenol | 56        |           | 10-120                 |
| 4-Terphenyl-d14      | 66        |           | 41-149                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/13/25 12:54  
**Analyst:** JJW

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/11/25 17:50

| Parameter                                                                                    | Result | Qualifier | Units | RL   | MDL  |
|----------------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-05 Batch: WG2052806-1 |        |           |       |      |      |
| Naphthalene                                                                                  | ND     |           | ug/l  | 0.10 | 0.02 |
| 2-Methylnaphthalene                                                                          | ND     |           | ug/l  | 0.10 | 0.03 |
| Acenaphthylene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Acenaphthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluorene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Phenanthrene                                                                                 | ND     |           | ug/l  | 0.05 | 0.04 |
| Anthracene                                                                                   | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluoranthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.03 |
| Pyrene                                                                                       | ND     |           | ug/l  | 0.10 | 0.04 |
| Benzo(a)anthracene                                                                           | ND     |           | ug/l  | 0.05 | 0.03 |
| Chrysene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(b)fluoranthene                                                                         | ND     |           | ug/l  | 0.05 | 0.03 |
| Benzo(k)fluoranthene                                                                         | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(a)pyrene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Indeno(1,2,3-cd)pyrene                                                                       | ND     |           | ug/l  | 0.10 | 0.02 |
| Dibenzo(a,h)anthracene                                                                       | ND     |           | ug/l  | 0.05 | 0.02 |
| Benzo(ghi)perylene                                                                           | ND     |           | ug/l  | 0.10 | 0.02 |

| Surrogate            | %Recovery | Qualifier | Acceptance Criteria |
|----------------------|-----------|-----------|---------------------|
| 2-Fluorophenol       | 51        |           | 21-120              |
| Phenol-d6            | 38        |           | 10-120              |
| Nitrobenzene-d5      | 77        |           | 23-120              |
| 2-Fluorobiphenyl     | 67        |           | 15-120              |
| 2,4,6-Tribromophenol | 79        |           | 10-120              |
| 4-Terphenyl-d14      | 70        |           | 41-149              |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522156

**Report Date:** 04/23/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2052805-2 WG2052805-3 |                  |      |                   |      |                     |     |      |               |
| Benzaldehyde                                                                                                | 60               |      | 64                |      | 40-140              | 6   |      | 30            |
| Phenol                                                                                                      | 27               |      | 28                |      | 12-110              | 4   |      | 30            |
| Bis(2-chloroethyl)ether                                                                                     | 57               |      | 58                |      | 40-140              | 2   |      | 30            |
| 2-Chlorophenol                                                                                              | 54               |      | 55                |      | 27-123              | 2   |      | 30            |
| 2-Methylphenol                                                                                              | 52               |      | 50                |      | 30-130              | 4   |      | 30            |
| Bis(2-chloroisopropyl)ether                                                                                 | 58               |      | 60                |      | 40-140              | 3   |      | 30            |
| Acetophenone                                                                                                | 61               |      | 62                |      | 39-129              | 2   |      | 30            |
| n-Nitrosodi-n-propylamine                                                                                   | 54               |      | 56                |      | 29-132              | 4   |      | 30            |
| 3-Methylphenol/4-Methylphenol                                                                               | 53               |      | 52                |      | 30-130              | 2   |      | 30            |
| Hexachloroethane                                                                                            | 41               |      | 46                |      | 40-140              | 11  |      | 30            |
| Nitrobenzene                                                                                                | 54               |      | 57                |      | 40-140              | 5   |      | 30            |
| Isophorone                                                                                                  | 53               |      | 54                |      | 40-140              | 2   |      | 30            |
| 2,4-Dimethylphenol                                                                                          | 50               |      | 50                |      | 30-130              | 0   |      | 30            |
| Bis(2-chloroethoxy)methane                                                                                  | 54               |      | 56                |      | 40-140              | 4   |      | 30            |
| 2,4-Dichlorophenol                                                                                          | 60               |      | 61                |      | 30-130              | 2   |      | 30            |
| Naphthalene                                                                                                 | 50               |      | 55                |      | 40-140              | 10  |      | 30            |
| 4-Chloroaniline                                                                                             | 40               |      | 42                |      | 40-140              | 5   |      | 30            |
| Hexachlorobutadiene                                                                                         | 43               |      | 47                |      | 40-140              | 9   |      | 30            |
| Caprolactam                                                                                                 | 30               |      | 32                |      | 10-130              | 6   |      | 30            |
| 2-Methylnaphthalene                                                                                         | 49               |      | 55                |      | 40-140              | 12  |      | 30            |
| Hexachlorocyclopentadiene                                                                                   | 30               | Q    | 36                | Q    | 40-140              | 18  |      | 30            |
| 1,2,4,5-Tetrachlorobenzene                                                                                  | 49               |      | 55                |      | 2-134               | 12  |      | 30            |
| 2,4,6-Trichlorophenol                                                                                       | 56               |      | 63                |      | 30-130              | 12  |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522156

**Report Date:** 04/23/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2052805-2 WG2052805-3 |                  |      |                   |      |                     |     |      |               |
| 2,4,5-Trichlorophenol                                                                                       | 63               |      | 71                |      | 30-130              | 12  |      | 30            |
| Biphenyl                                                                                                    | 55               |      | 60                |      | 40-140              | 9   |      | 30            |
| 2-Chloronaphthalene                                                                                         | 51               |      | 56                |      | 40-140              | 9   |      | 30            |
| 2-Nitroaniline                                                                                              | 62               |      | 65                |      | 52-143              | 5   |      | 30            |
| 2,6-Dinitrotoluene                                                                                          | 59               |      | 69                |      | 40-140              | 16  |      | 30            |
| Acenaphthylene                                                                                              | 52               |      | 58                |      | 45-123              | 11  |      | 30            |
| Acenaphthene                                                                                                | 53               |      | 57                |      | 37-111              | 7   |      | 30            |
| 2,4-Dinitrophenol                                                                                           | 54               |      | 59                |      | 20-130              | 9   |      | 30            |
| 2,4-Dinitrotoluene                                                                                          | 58               |      | 64                |      | 48-143              | 10  |      | 30            |
| 2,3,4,6-Tetrachlorophenol                                                                                   | 63               |      | 64                |      | 54-145              | 2   |      | 30            |
| Diethyl phthalate                                                                                           | 56               |      | 62                |      | 40-140              | 10  |      | 30            |
| Fluorene                                                                                                    | 55               |      | 60                |      | 40-140              | 9   |      | 30            |
| 4-Nitroaniline                                                                                              | 54               |      | 59                |      | 51-143              | 9   |      | 30            |
| NDPA/DPA                                                                                                    | 57               |      | 47                |      | 40-140              | 19  |      | 30            |
| Hexachlorobenzene                                                                                           | 58               |      | 63                |      | 40-140              | 8   |      | 30            |
| Pentachlorophenol                                                                                           | 60               |      | 71                |      | 9-103               | 17  |      | 30            |
| Phenanthrene                                                                                                | 60               |      | 62                |      | 40-140              | 3   |      | 30            |
| Anthracene                                                                                                  | 61               |      | 63                |      | 40-140              | 3   |      | 30            |
| Carbazole                                                                                                   | 60               |      | 64                |      | 55-144              | 6   |      | 30            |
| Di-n-butylphthalate                                                                                         | 60               |      | 65                |      | 40-140              | 8   |      | 30            |
| Fluoranthene                                                                                                | 63               |      | 66                |      | 40-140              | 5   |      | 30            |
| Pyrene                                                                                                      | 62               |      | 66                |      | 26-127              | 6   |      | 30            |
| 3,3'-Dichlorobenzidine                                                                                      | 57               |      | 42                |      | 40-140              | 30  |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522156

**Project Number:** 21010214

**Report Date:** 04/23/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2052805-2 WG2052805-3 |                  |      |                   |      |                     |     |      |               |
| Benzo(a)anthracene                                                                                          | 62               |      | 63                |      | 40-140              | 2   |      | 30            |
| Chrysene                                                                                                    | 64               |      | 65                |      | 40-140              | 2   |      | 30            |
| Bis(2-ethylhexyl)phthalate                                                                                  | 61               |      | 66                |      | 40-140              | 8   |      | 30            |
| Di-n-octylphthalate                                                                                         | 61               |      | 66                |      | 40-140              | 8   |      | 30            |
| Benzo(b)fluoranthene                                                                                        | 64               |      | 66                |      | 40-140              | 3   |      | 30            |
| Benzo(k)fluoranthene                                                                                        | 67               |      | 67                |      | 40-140              | 0   |      | 30            |
| Benzo(a)pyrene                                                                                              | 65               |      | 64                |      | 40-140              | 2   |      | 30            |
| Indeno(1,2,3-cd)pyrene                                                                                      | 64               |      | 65                |      | 40-140              | 2   |      | 30            |
| Dibenzo(a,h)anthracene                                                                                      | 65               |      | 66                |      | 40-140              | 2   |      | 30            |
| Benzo(ghi)perylene                                                                                          | 62               |      | 62                |      | 40-140              | 0   |      | 30            |

| Surrogate            | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | Acceptance<br>Criteria |
|----------------------|------------------|------|-------------------|------|------------------------|
| 2-Fluorophenol       | 41               |      | 40                |      | 21-120                 |
| Phenol-d6            | 29               |      | 30                |      | 10-120                 |
| Nitrobenzene-d5      | 55               |      | 56                |      | 23-120                 |
| 2-Fluorobiphenyl     | 57               |      | 61                |      | 15-120                 |
| 2,4,6-Tribromophenol | 60               |      | 62                |      | 10-120                 |
| 4-Terphenyl-d14      | 60               |      | 64                |      | 41-149                 |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2522156

**Report Date:** 04/23/25

| Parameter                                                                                                       | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-----------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-05 Batch: WG2052806-2 WG2052806-3 |                  |      |                   |      |                     |     |      |               |
| Naphthalene                                                                                                     | 51               |      | 59                |      | 40-140              | 15  |      | 40            |
| 2-Methylnaphthalene                                                                                             | 49               |      | 57                |      | 40-140              | 15  |      | 40            |
| Acenaphthylene                                                                                                  | 63               |      | 72                |      | 40-140              | 13  |      | 40            |
| Acenaphthene                                                                                                    | 57               |      | 67                |      | 37-111              | 16  |      | 40            |
| Fluorene                                                                                                        | 63               |      | 73                |      | 40-140              | 15  |      | 40            |
| Phenanthrene                                                                                                    | 61               |      | 70                |      | 40-140              | 14  |      | 40            |
| Anthracene                                                                                                      | 64               |      | 73                |      | 40-140              | 13  |      | 40            |
| Fluoranthene                                                                                                    | 62               |      | 68                |      | 40-140              | 9   |      | 40            |
| Pyrene                                                                                                          | 59               |      | 66                |      | 26-127              | 11  |      | 40            |
| Benzo(a)anthracene                                                                                              | 67               |      | 77                |      | 40-140              | 14  |      | 40            |
| Chrysene                                                                                                        | 63               |      | 72                |      | 40-140              | 13  |      | 40            |
| Benzo(b)fluoranthene                                                                                            | 70               |      | 83                |      | 40-140              | 17  |      | 40            |
| Benzo(k)fluoranthene                                                                                            | 69               |      | 78                |      | 40-140              | 12  |      | 40            |
| Benzo(a)pyrene                                                                                                  | 71               |      | 81                |      | 40-140              | 13  |      | 40            |
| Indeno(1,2,3-cd)pyrene                                                                                          | 73               |      | 84                |      | 40-140              | 14  |      | 40            |
| Dibenzo(a,h)anthracene                                                                                          | 76               |      | 86                |      | 40-140              | 12  |      | 40            |
| Benzo(ghi)perylene                                                                                              | 74               |      | 85                |      | 40-140              | 14  |      | 40            |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2522156

**Project Number:** 21010214

**Report Date:** 04/23/25

| <b>Parameter</b>                                                                                                | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-----------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-05 Batch: WG2052806-2 WG2052806-3 |                          |             |                           |             |                             |            |             |                       |

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 52                       |             | 56                        |             | 21-120                         |
| Phenol-d6            | 41                       |             | 44                        |             | 10-120                         |
| Nitrobenzene-d5      | 72                       |             | 78                        |             | 23-120                         |
| 2-Fluorobiphenyl     | 60                       |             | 65                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 77                       |             | 84                        |             | 10-120                         |
| 4-Terphenyl-d14      | 59                       |             | 62                        |             | 41-149                         |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2522156**Project Number:** 21010214**Report Date:** 04/23/25**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

**Cooler Information**

|               |                     |
|---------------|---------------------|
| <b>Cooler</b> | <b>Custody Seal</b> |
| A             | Absent              |

**Container Information**

| <b>Container ID</b> | <b>Container Type</b>   | <b>Cooler</b> | <b>Initial pH</b> | <b>Final pH</b> | <b>Temp deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen Date/Time</b> | <b>Analysis(*)</b>               |
|---------------------|-------------------------|---------------|-------------------|-----------------|-------------------|-------------|-------------|-------------------------|----------------------------------|
| L2522156-01A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-01B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-01C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-01D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522156-01E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522156-02A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-02B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-02C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-02D        | Amber 100ml unpreserved | A             | 12                | 12              | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522156-02E        | Amber 100ml unpreserved | A             | 12                | 12              | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522156-03A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-03B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-03C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-03D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522156-03E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522156-04A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-04B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-04C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-04D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522156-04E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2522156-05A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-05B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-05C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

Serial\_No:04232510:16  
**Lab Number:** L2522156  
**Report Date:** 04/23/25

**Container Information**

| <b>Container ID</b> | <b>Container Type</b>   | <b>Cooler</b> | <b>Initial pH</b> | <b>Final pH</b> | <b>Temp deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen Date/Time</b> | <b>Analysis(*)</b>               |
|---------------------|-------------------------|---------------|-------------------|-----------------|-------------------|-------------|-------------|-------------------------|----------------------------------|
| L2522156-05D        | Amber 250ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522156-05E        | Amber 250ml unpreserved | A             | 7                 | 7               | 3.8               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2522156-06A        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-06B        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-06C        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2522156-06D        | Vial HCl preserved      | A             | NA                |                 | 3.8               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |

**Container Comments**

L2522156-05D Amber-A.25  
L2522156-05E Amber-A.25

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

## GLOSSARY

### Acronyms

|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| DL       | - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                |
| EDL      | - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).                                                                                         |
| EMPC     | - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.                                                                                                                                                                                                                               |
| EPA      | - Environmental Protection Agency.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| LCS      | - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                          |
| LCSD     | - Laboratory Control Sample Duplicate: Refer to LCS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |
| LFB      | - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                         |
| LOD      | - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                                                               |
| LOQ      | - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)<br><br>Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| MDL      | - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                          |
| MS       | - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.                                                                                                                                                                                                    |
| MSD      | - Matrix Spike Sample Duplicate: Refer to MS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NA       | - Not Applicable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| NC       | - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.                                                                                                                                                                                                                                                                                                                                                                           |
| NDPA/DPA | - N-Nitrosodiphenylamine/Diphenylamine.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| NI       | - Not Ignitable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| NP       | - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NR       | - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.                                                                                                                                                                                                                                                                                                                                                                    |
| RL       | - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                                                                                                                                   |
| RPD      | - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.                                                                  |
| SRM      | - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.                                                                                                                                                                                                                                                                                                                                                                     |
| STLP     | - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| TEF      | - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.                                                                                                                                                                                                                                                                                                                                                                                             |
| TEQ      | - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.                                                                                                                                                                                                                                                                                                                                                        |
| TIC      | - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.                                                                                                                                                                                                                                                                      |

Report Format: DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

### Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

### Terms

**Analytical Method:** Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

**Chlordane:** The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

**Difference:** With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

**Final pH:** As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

**Frozen Date/Time:** With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**Initial pH:** As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

**PAH Total:** With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

**PFAS Total:** With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**Total:** With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

### Data Qualifiers

- A** - Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F** - The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The lower value for the two columns has been reported due to obvious interference.
- J** - Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

**Report Format:** DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

#### Data Qualifiers

Identified Compounds (TICs). For calculated parameters, this represents that one or more values used in the calculation were estimated.

- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND** - Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ** - Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- R** - Analytical results are from sample re-analysis.
- RE** - Analytical results are from sample re-extraction.
- S** - Analytical results are from modified screening analysis.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2522156  
**Report Date:** 04/23/25

## REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

## LIMITATION OF LIABILITIES

Pace Analytical Services performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Pace Analytical Services shall be to re-perform the work at it's own expense. In no event shall Pace Analytical Services be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Pace Analytical Services.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



**Pace Analytical Services LLC**Facility: **Northeast**Department: **Quality Assurance**Title: **Certificate/Approval Program Summary**ID No.: **17873**

Revision 27

Published Date: 01/24/2025

Page 1 of 2

**Certification Information**

The following analytes are not included in our Primary NELAP Scope of Accreditation:

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****EPA 624.1:** m/p-xylene, o-xylene, Naphthalene**EPA 625.1:** alpha-Terpineol**EPA 8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****SM 2540D:** TSS.**EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**MADEP-APH.****Nonpotable Water:** EPA RSK-175 Dissolved Gases**Biological Tissue Matrix:** EPA 3050B**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048****EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**Nonpotable Water:** EPA RSK-175 Dissolved Gases

The following test method is not included in our New Jersey Secondary NELAP Scope of Accreditation:

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Determination of Selected Perfluorinated Alkyl Substances by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry Isotope Dilution (via Alpha SOP 23528)**

The following analytes are included in our Massachusetts DEP Scope of Accreditation

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****Drinking Water****EPA 300.0:** Chloride, Nitrate-N, Fluoride, Sulfate; **EPA 353.2:** Nitrate-N, Nitrite-N; **SM4500NO3-F:** Nitrate-N, Nitrite-N; **SM4500F-C, SM4500CN-CE,****EPA 180.1, SM2130B, SM4500Cl-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B****EPA 524.2:** THMs and VOCs; **EPA 504.1:** EDB, DBCP.**Microbiology:** SM9215B; SM9223-P/A, SM9223B-Colilert-QT, SM9222D.**Non-Potable Water****SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH:** Ammonia-N and Kjeldahl-N, **EPA 350.1:**Ammonia-N, **LCHAT 10-107-06-1-B:** Ammonia-N, **EPA 351.1, SM4500NO3-F, EPA 353.2:** Nitrate-N, **SM4500P-E, SM4500P-B, E, SM4500SO4-E,****SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300:** Chloride, Sulfate, Nitrate.**EPA 624.1:** Volatile Halocarbons & Aromatics,**EPA 608.3:** Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II,

Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

**EPA 625.1:** SVOC (Acid/Base/Neutral Extractables).**Microbiology:** SM9223B-Colilert-QT; Enterolert-QT, EPA 1600, EPA 1603, SM9222D.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Drinking Water****EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1 Hg.****EPA 522, EPA 537.1.****Non-Potable Water****EPA 200.7:** Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.**EPA 200.8:** Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.**EPA 245.1 Hg.****SM2340B**



**Pace Analytical Services LLC**ID No.: **17873**Facility: **Northeast**

Revision 27

Department: **Quality Assurance**

Published Date: 01/24/2025

Title: **Certificate/Approval Program Summary**

Page 2 of 2

**Certification IDs:****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581**

CT PH-0826, IL 200077, IN C-MA-03, KY JY98045, ME MA00086, MD 348, MA M-MA086, NH 2064, NJ MA935, NY 11148, NC (DW) 25700, NC (NPW/SCM) 666, OR MA-1316, PA 68-03671, RI LAO00065, TX T104704476, VT VT-0935, VA 460195

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048**

CT PH-0825, ANAB/DoD L2474, IL 200081, IN C-MA-04, KY KY98046, LA 3090, ME MA00030, MI 9110, MN 025-999-495, NH 2062, NJ MA015, NY 11627, NC (NPW/SCM) 685, OR MA-0262, PA 68-02089, RI LAO00299, TX T-104704419, VT VT-0015, VA 460194, WA C954

**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048**

ANAB/DoD L2474, ME MA01156, MN 025-999-498, NH 2249, NJ MA025, NY 12191, OR 4203, TX T104704583, VA 460311, WA C1104.

---

For a complete listing of analytes and methods, please contact your Project Manager.





## ANALYTICAL REPORT

|                 |                                                                             |
|-----------------|-----------------------------------------------------------------------------|
| Lab Number:     | L2523111                                                                    |
| Client:         | Tradepoint Atlantic<br>1600 Sparrows Point Boulevard<br>Baltimore, MD 21219 |
| ATTN:           | Matthew Newman                                                              |
| Phone:          | (443) 649-5073                                                              |
| Project Name:   | B14 1ST HALF 2025 GW                                                        |
| Project Number: | 21010214                                                                    |
| Report Date:    | 04/29/25                                                                    |

The original project report/data package is held by Pace Analytical Services. This report/data package is paginated and should be reproduced only in its entirety. Pace Analytical Services holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0826), IL (200077), IN (C-MA-03), KY (KY98045), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), OR (MA-1316), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #525-23-122-91930A1).

---

Eight Walkup Drive, Westborough, MA 01581-1019  
508-898-9220 (Fax) 508-898-9193 800-624-9220 - [www.alphalab.com](http://www.alphalab.com)



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

| Lab<br>Sample ID | Client ID   | Matrix | Sample<br>Location | Collection<br>Date/Time | Receive Date |
|------------------|-------------|--------|--------------------|-------------------------|--------------|
| L2523111-01      | B24-002-MWS | WATER  | B14                | 04/15/25 09:50          | 04/15/25     |
| L2523111-02      | TM04-PZM006 | WATER  | B14                | 04/15/25 10:40          | 04/15/25     |
| L2523111-03      | HI13-MWS    | WATER  | B14                | 04/15/25 11:15          | 04/15/25     |
| L2523111-04      | TB-001      | WATER  | B14                | 04/15/25 00:00          | 04/15/25     |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

### Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Pace Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments and solids are reported on a dry weight basis unless otherwise noted. Tissues are reported "as received" or on a wet weight basis, unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

**HOLD POLICY** - For samples submitted on hold, Pace's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Pace Project Manager and made arrangements for Pace to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.

---

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**Case Narrative (continued)**

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature: *Tiffani Morrissey* - Tiffani Morrissey

Title: Technical Director/Representative

Date: 04/29/25

# ORGANICS

# VOLATILES



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-01  
**Client ID:** B24-002-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 09:50  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/22/25 11:34  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 8.2    |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | 0.56   | J         | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | 2.9    |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | 4.5    |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-01  
**Client ID:** B24-002-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 09:50  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | 6.9    |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | 7.9    |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 125        |           | 70-130              |
| Toluene-d8            | 100        |           | 70-130              |
| 4-Bromofluorobenzene  | 97         |           | 70-130              |
| Dibromofluoromethane  | 113        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-01  
**Client ID:** B24-002-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 09:50  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D-SIM(M)  
**Analytical Date:** 04/22/25 11:34  
**Analyst:** MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 121        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-02  
**Client ID:** TM04-PZM006  
**Sample Location:** B14

**Date Collected:** 04/15/25 10:40  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/22/25 11:07  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | 0.15   | J         | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 5.8    |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | 0.22   | J         | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 64     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | 0.36   | J         | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-02  
**Client ID:** TM04-PZM006  
**Sample Location:** B14

**Date Collected:** 04/15/25 10:40  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | 0.71   |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | 0.65   | J         | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | 0.65   | J         | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 119        |           | 70-130              |
| Toluene-d8            | 100        |           | 70-130              |
| 4-Bromofluorobenzene  | 98         |           | 70-130              |
| Dibromofluoromethane  | 106        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

Lab ID: L2523111-02  
 Client ID: TM04-PZM006  
 Sample Location: B14

Date Collected: 04/15/25 10:40  
 Date Received: 04/15/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/22/25 11:07  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 116        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-03  
**Client ID:** HI13-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 11:15  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/22/25 10:41  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 3.7    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.20   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25**SAMPLE RESULTS****Lab ID:** L2523111-03**Date Collected:** 04/15/25 11:15**Client ID:** HI13-MWS**Date Received:** 04/15/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 123        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 111        |           | 70-130              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

Lab ID: L2523111-03  
 Client ID: HI13-MWS  
 Sample Location: B14

Date Collected: 04/15/25 11:15  
 Date Received: 04/15/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/22/25 10:41  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 121        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 92         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-04  
**Client ID:** TB-001  
**Sample Location:** B14

**Date Collected:** 04/15/25 00:00  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 04/22/25 09:21  
**Analyst:** MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 1.5    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25**SAMPLE RESULTS****Lab ID:** L2523111-04**Date Collected:** 04/15/25 00:00**Client ID:** TB-001**Date Received:** 04/15/25**Sample Location:** B14**Field Prep:** Not Specified**Sample Depth:**

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 124        |           | 70-130              |
| Toluene-d8            | 99         |           | 70-130              |
| 4-Bromofluorobenzene  | 97         |           | 70-130              |
| Dibromofluoromethane  | 111        |           | 70-130              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

Lab ID: L2523111-04  
 Client ID: TB-001  
 Sample Location: B14

Date Collected: 04/15/25 00:00  
 Date Received: 04/15/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/22/25 09:21  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 121        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 93         |           | 70-130              |                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D-SIM(M)  
Analytical Date: 04/22/25 06:43  
Analyst: MCM

| Parameter                                                                                | Result | Qualifier | Units | RL    | MDL   |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-------|-------|
| Volatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-04 Batch: WG2057033-5 |        |           |       |       |       |
| 1,1,1,2-Tetrachloroethane                                                                | ND     |           | ug/l  | 0.050 | 0.006 |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 120       |           | 70-130                 |
| 4-Bromofluorobenzene  | 93        |           | 70-130                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 04/22/25 06:43  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2057035-5 |        |           |       |      |      |
| Dichlorodifluoromethane                                                              | ND     |           | ug/l  | 5.0  | 0.24 |
| Chloromethane                                                                        | ND     |           | ug/l  | 2.5  | 0.20 |
| Vinyl chloride                                                                       | ND     |           | ug/l  | 1.0  | 0.07 |
| Bromomethane                                                                         | ND     |           | ug/l  | 1.0  | 0.26 |
| Chloroethane                                                                         | ND     |           | ug/l  | 1.0  | 0.13 |
| Trichlorofluoromethane                                                               | ND     |           | ug/l  | 2.5  | 0.16 |
| 1,1-Dichloroethene                                                                   | ND     |           | ug/l  | 0.50 | 0.17 |
| Carbon disulfide                                                                     | ND     |           | ug/l  | 5.0  | 0.30 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                | ND     |           | ug/l  | 2.5  | 0.15 |
| Methylene chloride                                                                   | ND     |           | ug/l  | 2.5  | 0.68 |
| Acetone                                                                              | ND     |           | ug/l  | 5.0  | 1.5  |
| trans-1,2-Dichloroethene                                                             | ND     |           | ug/l  | 0.75 | 0.16 |
| Methyl Acetate                                                                       | ND     |           | ug/l  | 2.0  | 0.23 |
| Methyl tert butyl ether                                                              | ND     |           | ug/l  | 1.0  | 0.17 |
| 1,1-Dichloroethane                                                                   | ND     |           | ug/l  | 0.75 | 0.21 |
| cis-1,2-Dichloroethene                                                               | ND     |           | ug/l  | 0.50 | 0.19 |
| Cyclohexane                                                                          | ND     |           | ug/l  | 10   | 0.27 |
| Chloroform                                                                           | ND     |           | ug/l  | 0.75 | 0.22 |
| Carbon tetrachloride                                                                 | ND     |           | ug/l  | 0.50 | 0.13 |
| 1,1,1-Trichloroethane                                                                | ND     |           | ug/l  | 0.50 | 0.16 |
| 2-Butanone                                                                           | ND     |           | ug/l  | 5.0  | 1.9  |
| Benzene                                                                              | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,2-Dichloroethane                                                                   | ND     |           | ug/l  | 0.50 | 0.13 |
| Trichloroethene                                                                      | ND     |           | ug/l  | 0.50 | 0.18 |
| 1,2-Dichloropropane                                                                  | ND     |           | ug/l  | 1.0  | 0.14 |
| Bromodichloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.19 |
| cis-1,3-Dichloropropene                                                              | ND     |           | ug/l  | 0.50 | 0.14 |
| Toluene                                                                              | ND     |           | ug/l  | 0.75 | 0.20 |
| Tetrachloroethene                                                                    | ND     |           | ug/l  | 0.50 | 0.18 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8260D  
**Analytical Date:** 04/22/25 06:43  
**Analyst:** MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2057035-5 |        |           |       |      |      |
| 4-Methyl-2-pentanone                                                                 | ND     |           | ug/l  | 5.0  | 0.42 |
| trans-1,3-Dichloropropene                                                            | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,3-Dichloropropene, Total                                                           | ND     |           | ug/l  | 0.50 | 0.14 |
| 1,1,2-Trichloroethane                                                                | ND     |           | ug/l  | 0.75 | 0.14 |
| Dibromochloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.15 |
| 1,2-Dibromoethane                                                                    | ND     |           | ug/l  | 2.0  | 0.19 |
| 2-Hexanone                                                                           | ND     |           | ug/l  | 5.0  | 0.52 |
| Chlorobenzene                                                                        | ND     |           | ug/l  | 0.50 | 0.18 |
| Ethylbenzene                                                                         | ND     |           | ug/l  | 0.50 | 0.17 |
| p/m-Xylene                                                                           | ND     |           | ug/l  | 1.0  | 0.33 |
| o-Xylene                                                                             | ND     |           | ug/l  | 1.0  | 0.39 |
| Xylenes, Total                                                                       | ND     |           | ug/l  | 1.0  | 0.33 |
| Styrene                                                                              | ND     |           | ug/l  | 1.0  | 0.36 |
| Bromoform                                                                            | ND     |           | ug/l  | 2.0  | 0.25 |
| Isopropylbenzene                                                                     | ND     |           | ug/l  | 0.50 | 0.19 |
| 1,1,2,2-Tetrachloroethane                                                            | ND     |           | ug/l  | 0.50 | 0.17 |
| 1,3-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,4-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,2-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.18 |
| 1,2-Dibromo-3-chloropropane                                                          | ND     |           | ug/l  | 2.5  | 0.35 |
| 1,2,4-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.22 |
| 1,2,3-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.23 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
Analytical Date: 04/22/25 06:43  
Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL | MDL |
|--------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2057035-5 |        |           |       |    |     |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 123       |           | 70-130                 |
| Toluene-d8            | 100       |           | 70-130                 |
| 4-Bromofluorobenzene  | 97        |           | 70-130                 |
| Dibromofluoromethane  | 110       |           | 70-130                 |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2523111

**Project Number:** 21010214

**Report Date:** 04/29/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2057033-3 WG2057033-4 |                          |             |                           |             |                             |            |             |                       |
| 1,1,1,2-Tetrachloroethane                                                                                   | 104                      |             | 103                       |             | 70-130                      | 1          |             | 25                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 120                      |             | 122                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 92                       |             | 91                        |             | 70-130                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2523111

**Report Date:** 04/29/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2057035-3 WG2057035-4 |                  |      |                   |      |                     |     |      |               |
| Dichlorodifluoromethane                                                                                 | 97               |      | 99                |      | 36-147              | 2   |      | 20            |
| Chloromethane                                                                                           | 110              |      | 110               |      | 64-130              | 0   |      | 20            |
| Vinyl chloride                                                                                          | 110              |      | 110               |      | 55-140              | 0   |      | 20            |
| Bromomethane                                                                                            | 59               |      | 61                |      | 39-139              | 3   |      | 20            |
| Chloroethane                                                                                            | 120              |      | 120               |      | 55-138              | 0   |      | 20            |
| Trichlorofluoromethane                                                                                  | 110              |      | 110               |      | 62-150              | 0   |      | 20            |
| 1,1-Dichloroethene                                                                                      | 100              |      | 100               |      | 61-145              | 0   |      | 20            |
| Carbon disulfide                                                                                        | 110              |      | 110               |      | 51-130              | 0   |      | 20            |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                                   | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Methylene chloride                                                                                      | 100              |      | 110               |      | 70-130              | 10  |      | 20            |
| Acetone                                                                                                 | 130              |      | 130               |      | 58-148              | 0   |      | 20            |
| trans-1,2-Dichloroethene                                                                                | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Methyl Acetate                                                                                          | 140              | Q    | 140               | Q    | 70-130              | 0   |      | 20            |
| Methyl tert butyl ether                                                                                 | 100              |      | 100               |      | 63-130              | 0   |      | 20            |
| 1,1-Dichloroethane                                                                                      | 120              |      | 120               |      | 70-130              | 0   |      | 20            |
| cis-1,2-Dichloroethene                                                                                  | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| Cyclohexane                                                                                             | 120              |      | 120               |      | 70-130              | 0   |      | 20            |
| Chloroform                                                                                              | 100              |      | 110               |      | 70-130              | 10  |      | 20            |
| Carbon tetrachloride                                                                                    | 100              |      | 110               |      | 63-132              | 10  |      | 20            |
| 1,1,1-Trichloroethane                                                                                   | 100              |      | 110               |      | 67-130              | 10  |      | 20            |
| 2-Butanone                                                                                              | 110              |      | 110               |      | 63-138              | 0   |      | 20            |
| Benzene                                                                                                 | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| 1,2-Dichloroethane                                                                                      | 110              |      | 120               |      | 70-130              | 9   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2523111

**Report Date:** 04/29/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2057035-3 WG2057035-4 |                  |      |                   |      |                     |     |      |               |
| Trichloroethene                                                                                         | 99               |      | 100               |      | 70-130              | 1   |      | 20            |
| 1,2-Dichloropropane                                                                                     | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Bromodichloromethane                                                                                    | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| cis-1,3-Dichloropropene                                                                                 | 94               |      | 97                |      | 70-130              | 3   |      | 20            |
| Toluene                                                                                                 | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| Tetrachloroethene                                                                                       | 96               |      | 99                |      | 70-130              | 3   |      | 20            |
| 4-Methyl-2-pentanone                                                                                    | 89               |      | 92                |      | 59-130              | 3   |      | 20            |
| trans-1,3-Dichloropropene                                                                               | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| 1,1,2-Trichloroethane                                                                                   | 100              |      | 110               |      | 70-130              | 10  |      | 20            |
| Dibromochloromethane                                                                                    | 93               |      | 97                |      | 63-130              | 4   |      | 20            |
| 1,2-Dibromoethane                                                                                       | 97               |      | 100               |      | 70-130              | 3   |      | 20            |
| 2-Hexanone                                                                                              | 100              |      | 100               |      | 57-130              | 0   |      | 20            |
| Chlorobenzene                                                                                           | 95               |      | 99                |      | 75-130              | 4   |      | 20            |
| Ethylbenzene                                                                                            | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| p/m-Xylene                                                                                              | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| o-Xylene                                                                                                | 95               |      | 100               |      | 70-130              | 5   |      | 20            |
| Styrene                                                                                                 | 95               |      | 100               |      | 70-130              | 5   |      | 20            |
| Bromoform                                                                                               | 88               |      | 91                |      | 54-136              | 3   |      | 20            |
| Isopropylbenzene                                                                                        | 97               |      | 100               |      | 70-130              | 3   |      | 20            |
| 1,1,2,2-Tetrachloroethane                                                                               | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| 1,3-Dichlorobenzene                                                                                     | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| 1,4-Dichlorobenzene                                                                                     | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| 1,2-Dichlorobenzene                                                                                     | 98               |      | 100               |      | 70-130              | 2   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2523111

**Project Number:** 21010214

**Report Date:** 04/29/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCS<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2057035-3 WG2057035-4 |                  |      |                  |      |                     |     |      |               |
| 1,2-Dibromo-3-chloropropane                                                                             | 81               |      | 87               |      | 41-144              | 7   |      | 20            |
| 1,2,4-Trichlorobenzene                                                                                  | 90               |      | 92               |      | 70-130              | 2   |      | 20            |
| 1,2,3-Trichlorobenzene                                                                                  | 92               |      | 94               |      | 70-130              | 2   |      | 20            |

| Surrogate             | LCS<br>%Recovery | Qual | LCS<br>%Recovery | Qual | Acceptance<br>Criteria |
|-----------------------|------------------|------|------------------|------|------------------------|
| 1,2-Dichloroethane-d4 | 117              |      | 117              |      | 70-130                 |
| Toluene-d8            | 101              |      | 101              |      | 70-130                 |
| 4-Bromofluorobenzene  | 96               |      | 96               |      | 70-130                 |
| Dibromofluoromethane  | 106              |      | 105              |      | 70-130                 |

# SEMIVOLATILES

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-01  
**Client ID:** B24-002-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 09:50  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/23/25 04:43  
**Analyst:** SMZ

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523111-01  
 Client ID: B24-002-MWS  
 Sample Location: B14

Date Collected: 04/15/25 09:50  
 Date Received: 04/15/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 72         |           | 21-120              |
| Phenol-d6            | 52         |           | 10-120              |
| Nitrobenzene-d5      | 108        |           | 23-120              |
| 2-Fluorobiphenyl     | 108        |           | 15-120              |
| 2,4,6-Tribromophenol | 119        |           | 10-120              |
| 4-Terphenyl-d14      | 126        |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-01  
**Client ID:** B24-002-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 09:50  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/22/25 13:05  
**Analyst:** LJG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.08   | J         | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.03   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.05   | J         | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 105        |           | 23-120              |
| 2-Fluorobiphenyl | 75         |           | 15-120              |
| 4-Terphenyl-d14  | 94         |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-02  
**Client ID:** TM04-PZM006  
**Sample Location:** B14

**Date Collected:** 04/15/25 10:40  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/23/25 05:06  
**Analyst:** SMZ

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523111-02  
 Client ID: TM04-PZM006  
 Sample Location: B14

Date Collected: 04/15/25 10:40  
 Date Received: 04/15/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 69         |           | 21-120              |
| Phenol-d6            | 50         |           | 10-120              |
| Nitrobenzene-d5      | 110        |           | 23-120              |
| 2-Fluorobiphenyl     | 100        |           | 15-120              |
| 2,4,6-Tribromophenol | 119        |           | 10-120              |
| 4-Terphenyl-d14      | 115        |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-02  
**Client ID:** TM04-PZM006  
**Sample Location:** B14

**Date Collected:** 04/15/25 10:40  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/22/25 13:21  
**Analyst:** LJG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.12   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | 0.07   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.08   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.05   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.06   | J         | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 62         |           | 21-120              |
| Phenol-d6            | 50         |           | 10-120              |
| Nitrobenzene-d5      | 128        | Q         | 23-120              |
| 2-Fluorobiphenyl     | 89         |           | 15-120              |
| 2,4,6-Tribromophenol | 117        |           | 10-120              |
| 4-Terphenyl-d14      | 104        |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-03  
**Client ID:** HI13-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 11:15  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E  
**Analytical Date:** 04/23/25 05:29  
**Analyst:** SMZ

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523111-03

Date Collected: 04/15/25 11:15

Client ID: HI13-MWS

Date Received: 04/15/25

Sample Location: B14

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 77         |           | 21-120              |
| Phenol-d6            | 60         |           | 10-120              |
| Nitrobenzene-d5      | 116        |           | 23-120              |
| 2-Fluorobiphenyl     | 111        |           | 15-120              |
| 2,4,6-Tribromophenol | 115        |           | 10-120              |
| 4-Terphenyl-d14      | 129        |           | 41-149              |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

**Lab ID:** L2523111-03  
**Client ID:** HI13-MWS  
**Sample Location:** B14

**Date Collected:** 04/15/25 11:15  
**Date Received:** 04/15/25  
**Field Prep:** Not Specified

**Sample Depth:**

**Matrix:** Water  
**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/22/25 13:37  
**Analyst:** LJG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | 0.21   |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | ND     |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate        | % Recovery | Qualifier | Acceptance Criteria |
|------------------|------------|-----------|---------------------|
| Nitrobenzene-d5  | 134        | Q         | 23-120              |
| 2-Fluorobiphenyl | 92         |           | 15-120              |
| 4-Terphenyl-d14  | 115        |           | 41-149              |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/22/25 23:46  
**Analyst:** SMZ

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-03 Batch: WG2056799-1 |        |           |       |     |      |
| Benzaldehyde                                                                             | ND     |           | ug/l  | 5.0 | 1.1  |
| Phenol                                                                                   | ND     |           | ug/l  | 5.0 | 0.35 |
| Bis(2-chloroethyl)ether                                                                  | ND     |           | ug/l  | 2.0 | 0.39 |
| 2-Chlorophenol                                                                           | ND     |           | ug/l  | 2.0 | 0.65 |
| 2-Methylphenol                                                                           | ND     |           | ug/l  | 5.0 | 2.3  |
| Bis(2-chloroisopropyl)ether                                                              | ND     |           | ug/l  | 2.0 | 0.40 |
| Acetophenone                                                                             | ND     |           | ug/l  | 5.0 | 0.92 |
| n-Nitrosodi-n-propylamine                                                                | ND     |           | ug/l  | 5.0 | 0.91 |
| 3-Methylphenol/4-Methylphenol                                                            | ND     |           | ug/l  | 5.0 | 1.4  |
| Hexachloroethane                                                                         | ND     |           | ug/l  | 2.0 | 0.20 |
| Nitrobenzene                                                                             | ND     |           | ug/l  | 1.4 | 0.20 |
| Isophorone                                                                               | ND     |           | ug/l  | 5.0 | 0.86 |
| 2,4-Dimethylphenol                                                                       | ND     |           | ug/l  | 5.0 | 2.0  |
| Bis(2-chloroethoxy)methane                                                               | ND     |           | ug/l  | 5.0 | 0.84 |
| 2,4-Dichlorophenol                                                                       | ND     |           | ug/l  | 5.0 | 1.7  |
| Naphthalene                                                                              | ND     |           | ug/l  | 2.0 | 0.54 |
| 4-Chloroaniline                                                                          | ND     |           | ug/l  | 3.7 | 0.47 |
| Hexachlorobutadiene                                                                      | ND     |           | ug/l  | 2.0 | 0.36 |
| Caprolactam                                                                              | ND     |           | ug/l  | 10  | 1.2  |
| 2-Methylnaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.37 |
| Hexachlorocyclopentadiene                                                                | ND     |           | ug/l  | 20  | 1.2  |
| 1,2,4,5-Tetrachlorobenzene                                                               | ND     |           | ug/l  | 1.7 | 0.24 |
| 2,4,6-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| 2,4,5-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| Biphenyl                                                                                 | ND     |           | ug/l  | 2.0 | 0.20 |
| 2-Chloronaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.35 |
| 2-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.0  |
| 2,6-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.84 |
| Acenaphthylene                                                                           | ND     |           | ug/l  | 2.0 | 0.32 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E  
**Analytical Date:** 04/22/25 23:46  
**Analyst:** SMZ

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-03 Batch: WG2056799-1 |        |           |       |     |      |
| Acenaphthene                                                                             | ND     |           | ug/l  | 2.0 | 0.40 |
| 2,4-Dinitrophenol                                                                        | ND     |           | ug/l  | 20  | 5.4  |
| 2,4-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.54 |
| 2,3,4,6-Tetrachlorophenol                                                                | ND     |           | ug/l  | 5.0 | 2.2  |
| Diethyl phthalate                                                                        | ND     |           | ug/l  | 5.0 | 0.76 |
| Fluorene                                                                                 | ND     |           | ug/l  | 2.0 | 0.44 |
| 4-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.4  |
| NDPA/DPA                                                                                 | ND     |           | ug/l  | 2.0 | 0.92 |
| Hexachlorobenzene                                                                        | ND     |           | ug/l  | 2.0 | 0.45 |
| Pentachlorophenol                                                                        | ND     |           | ug/l  | 10  | 2.5  |
| Phenanthrene                                                                             | ND     |           | ug/l  | 2.0 | 0.42 |
| Anthracene                                                                               | ND     |           | ug/l  | 2.0 | 0.47 |
| Carbazole                                                                                | ND     |           | ug/l  | 2.0 | 0.31 |
| Di-n-butylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 0.96 |
| Fluoranthene                                                                             | ND     |           | ug/l  | 2.0 | 0.41 |
| Pyrene                                                                                   | ND     |           | ug/l  | 2.0 | 0.41 |
| 3,3'-Dichlorobenzidine                                                                   | ND     |           | ug/l  | 5.0 | 1.8  |
| Benzo(a)anthracene                                                                       | ND     |           | ug/l  | 2.0 | 0.32 |
| Chrysene                                                                                 | ND     |           | ug/l  | 1.5 | 0.22 |
| Bis(2-ethylhexyl)phthalate                                                               | ND     |           | ug/l  | 3.0 | 1.4  |
| Di-n-octylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 2.3  |
| Benzo(b)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.53 |
| Benzo(k)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.62 |
| Benzo(a)pyrene                                                                           | ND     |           | ug/l  | 2.0 | 0.37 |
| Indeno(1,2,3-cd)pyrene                                                                   | ND     |           | ug/l  | 2.0 | 0.48 |
| Dibenzo(a,h)anthracene                                                                   | ND     |           | ug/l  | 2.0 | 0.29 |
| Benzo(ghi)perylene                                                                       | ND     |           | ug/l  | 2.0 | 0.37 |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8270E  
Analytical Date: 04/22/25 23:46  
Analyst: SMZ

Extraction Method: EPA 3510C  
Extraction Date: 04/22/25 00:12

| Parameter                                                                                | Result | Qualifier | Units | RL | MDL |
|------------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-03 Batch: WG2056799-1 |        |           |       |    |     |

| Surrogate            | %Recovery | Qualifier | Acceptance<br>Criteria |
|----------------------|-----------|-----------|------------------------|
| 2-Fluorophenol       | 68        |           | 21-120                 |
| Phenol-d6            | 50        |           | 10-120                 |
| Nitrobenzene-d5      | 100       |           | 23-120                 |
| 2-Fluorobiphenyl     | 103       |           | 15-120                 |
| 2,4,6-Tribromophenol | 110       |           | 10-120                 |
| 4-Terphenyl-d14      | 127       |           | 41-149                 |

**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/22/25 12:48  
**Analyst:** LJG

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 00:12

| Parameter                                                                                    | Result | Qualifier | Units | RL   | MDL  |
|----------------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-03 Batch: WG2056800-1 |        |           |       |      |      |
| Naphthalene                                                                                  | ND     |           | ug/l  | 0.10 | 0.02 |
| 2-Methylnaphthalene                                                                          | ND     |           | ug/l  | 0.10 | 0.03 |
| Acenaphthylene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Acenaphthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluorene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Phenanthrene                                                                                 | ND     |           | ug/l  | 0.05 | 0.04 |
| Anthracene                                                                                   | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluoranthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.03 |
| Pyrene                                                                                       | ND     |           | ug/l  | 0.10 | 0.04 |
| Benzo(a)anthracene                                                                           | ND     |           | ug/l  | 0.05 | 0.03 |
| Chrysene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(b)fluoranthene                                                                         | ND     |           | ug/l  | 0.05 | 0.03 |
| Benzo(k)fluoranthene                                                                         | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(a)pyrene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Indeno(1,2,3-cd)pyrene                                                                       | ND     |           | ug/l  | 0.10 | 0.02 |
| Dibenzo(a,h)anthracene                                                                       | ND     |           | ug/l  | 0.05 | 0.02 |
| Benzo(ghi)perylene                                                                           | ND     |           | ug/l  | 0.10 | 0.02 |

| Surrogate            | %Recovery | Qualifier | Acceptance Criteria |
|----------------------|-----------|-----------|---------------------|
| 2-Fluorophenol       | 68        |           | 21-120              |
| Phenol-d6            | 55        |           | 10-120              |
| Nitrobenzene-d5      | 135       | Q         | 23-120              |
| 2-Fluorobiphenyl     | 94        |           | 15-120              |
| 2,4,6-Tribromophenol | 113       |           | 10-120              |
| 4-Terphenyl-d14      | 123       |           | 41-149              |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2523111

**Report Date:** 04/29/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-03 Batch: WG2056799-2 WG2056799-3 |                  |      |                   |      |                     |     |      |               |
| Benzaldehyde                                                                                                | 113              |      | 113               |      | 40-140              | 0   |      | 30            |
| Phenol                                                                                                      | 55               |      | 56                |      | 12-110              | 2   |      | 30            |
| Bis(2-chloroethyl)ether                                                                                     | 111              |      | 108               |      | 40-140              | 3   |      | 30            |
| 2-Chlorophenol                                                                                              | 104              |      | 102               |      | 27-123              | 2   |      | 30            |
| 2-Methylphenol                                                                                              | 96               |      | 100               |      | 30-130              | 4   |      | 30            |
| Bis(2-chloroisopropyl)ether                                                                                 | 109              |      | 114               |      | 40-140              | 4   |      | 30            |
| Acetophenone                                                                                                | 112              |      | 119               |      | 39-129              | 6   |      | 30            |
| n-Nitrosodi-n-propylamine                                                                                   | 108              |      | 104               |      | 29-132              | 4   |      | 30            |
| 3-Methylphenol/4-Methylphenol                                                                               | 100              |      | 97                |      | 30-130              | 3   |      | 30            |
| Hexachloroethane                                                                                            | 81               |      | 90                |      | 40-140              | 11  |      | 30            |
| Nitrobenzene                                                                                                | 104              |      | 107               |      | 40-140              | 3   |      | 30            |
| Isophorone                                                                                                  | 105              |      | 104               |      | 40-140              | 1   |      | 30            |
| 2,4-Dimethylphenol                                                                                          | 86               |      | 90                |      | 30-130              | 5   |      | 30            |
| Bis(2-chloroethoxy)methane                                                                                  | 107              |      | 110               |      | 40-140              | 3   |      | 30            |
| 2,4-Dichlorophenol                                                                                          | 122              |      | 118               |      | 30-130              | 3   |      | 30            |
| Naphthalene                                                                                                 | 103              |      | 101               |      | 40-140              | 2   |      | 30            |
| 4-Chloroaniline                                                                                             | 48               |      | 50                |      | 40-140              | 4   |      | 30            |
| Hexachlorobutadiene                                                                                         | 82               |      | 88                |      | 40-140              | 7   |      | 30            |
| Caprolactam                                                                                                 | 44               |      | 40                |      | 10-130              | 10  |      | 30            |
| 2-Methylnaphthalene                                                                                         | 106              |      | 104               |      | 40-140              | 2   |      | 30            |
| Hexachlorocyclopentadiene                                                                                   | 77               |      | 80                |      | 40-140              | 4   |      | 30            |
| 1,2,4,5-Tetrachlorobenzene                                                                                  | 107              |      | 106               |      | 2-134               | 1   |      | 30            |
| 2,4,6-Trichlorophenol                                                                                       | 125              |      | 114               |      | 30-130              | 9   |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2523111

**Report Date:** 04/29/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-03 Batch: WG2056799-2 WG2056799-3 |                  |      |                   |      |                     |     |      |               |
| 2,4,5-Trichlorophenol                                                                                       | 136              | Q    | 128               |      | 30-130              | 6   |      | 30            |
| Biphenyl                                                                                                    | 117              |      | 113               |      | 40-140              | 3   |      | 30            |
| 2-Chloronaphthalene                                                                                         | 111              |      | 107               |      | 40-140              | 4   |      | 30            |
| 2-Nitroaniline                                                                                              | 128              |      | 120               |      | 52-143              | 6   |      | 30            |
| 2,6-Dinitrotoluene                                                                                          | 131              |      | 116               |      | 40-140              | 12  |      | 30            |
| Acenaphthylene                                                                                              | 115              |      | 111               |      | 45-123              | 4   |      | 30            |
| Acenaphthene                                                                                                | 112              | Q    | 110               |      | 37-111              | 2   |      | 30            |
| 2,4-Dinitrophenol                                                                                           | 86               |      | 83                |      | 20-130              | 4   |      | 30            |
| 2,4-Dinitrotoluene                                                                                          | 123              |      | 113               |      | 48-143              | 8   |      | 30            |
| 2,3,4,6-Tetrachlorophenol                                                                                   | 126              |      | 121               |      | 54-145              | 4   |      | 30            |
| Diethyl phthalate                                                                                           | 116              |      | 111               |      | 40-140              | 4   |      | 30            |
| Fluorene                                                                                                    | 118              |      | 114               |      | 40-140              | 3   |      | 30            |
| 4-Nitroaniline                                                                                              | 113              |      | 107               |      | 51-143              | 5   |      | 30            |
| NDPA/DPA                                                                                                    | 123              |      | 117               |      | 40-140              | 5   |      | 30            |
| Hexachlorobenzene                                                                                           | 120              |      | 116               |      | 40-140              | 3   |      | 30            |
| Pentachlorophenol                                                                                           | 126              | Q    | 114               | Q    | 9-103               | 10  |      | 30            |
| Phenanthrene                                                                                                | 122              |      | 117               |      | 40-140              | 4   |      | 30            |
| Anthracene                                                                                                  | 129              |      | 123               |      | 40-140              | 5   |      | 30            |
| Carbazole                                                                                                   | 128              |      | 121               |      | 55-144              | 6   |      | 30            |
| Di-n-butylphthalate                                                                                         | 130              |      | 119               |      | 40-140              | 9   |      | 30            |
| Fluoranthene                                                                                                | 132              |      | 124               |      | 40-140              | 6   |      | 30            |
| Pyrene                                                                                                      | 132              | Q    | 125               |      | 26-127              | 5   |      | 30            |
| 3,3'-Dichlorobenzidine                                                                                      | 108              |      | 98                |      | 40-140              | 10  |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2523111

**Project Number:** 21010214

**Report Date:** 04/29/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-03 Batch: WG2056799-2 WG2056799-3 |                  |      |                   |      |                     |     |      |               |
| Benzo(a)anthracene                                                                                          | 127              |      | 119               |      | 40-140              | 7   |      | 30            |
| Chrysene                                                                                                    | 128              |      | 123               |      | 40-140              | 4   |      | 30            |
| Bis(2-ethylhexyl)phthalate                                                                                  | 137              |      | 126               |      | 40-140              | 8   |      | 30            |
| Di-n-octylphthalate                                                                                         | 144              | Q    | 130               |      | 40-140              | 10  |      | 30            |
| Benzo(b)fluoranthene                                                                                        | 134              |      | 126               |      | 40-140              | 6   |      | 30            |
| Benzo(k)fluoranthene                                                                                        | 135              |      | 125               |      | 40-140              | 8   |      | 30            |
| Benzo(a)pyrene                                                                                              | 135              |      | 125               |      | 40-140              | 8   |      | 30            |
| Indeno(1,2,3-cd)pyrene                                                                                      | 127              |      | 118               |      | 40-140              | 7   |      | 30            |
| Dibenzo(a,h)anthracene                                                                                      | 128              |      | 125               |      | 40-140              | 2   |      | 30            |
| Benzo(ghi)perylene                                                                                          | 123              |      | 121               |      | 40-140              | 2   |      | 30            |

| Surrogate            | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | Acceptance<br>Criteria |
|----------------------|------------------|------|-------------------|------|------------------------|
| 2-Fluorophenol       | 71               |      | 73                |      | 21-120                 |
| Phenol-d6            | 56               |      | 56                |      | 10-120                 |
| Nitrobenzene-d5      | 96               |      | 100               |      | 23-120                 |
| 2-Fluorobiphenyl     | 106              |      | 99                |      | 15-120                 |
| 2,4,6-Tribromophenol | 112              |      | 108               |      | 10-120                 |
| 4-Terphenyl-d14      | 120              |      | 113               |      | 41-149                 |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Project Number:** 21010214

**Lab Number:** L2523111

**Report Date:** 04/29/25

| Parameter                                                                                                       | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-----------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-03 Batch: WG2056800-2 WG2056800-3 |                  |      |                   |      |                     |     |      |               |
| Naphthalene                                                                                                     | 44               |      | 67                |      | 40-140              | 41  | Q    | 40            |
| 2-Methylnaphthalene                                                                                             | 45               |      | 69                |      | 40-140              | 42  | Q    | 40            |
| Acenaphthylene                                                                                                  | 56               |      | 82                |      | 40-140              | 38  |      | 40            |
| Acenaphthene                                                                                                    | 46               |      | 68                |      | 37-111              | 39  |      | 40            |
| Fluorene                                                                                                        | 52               |      | 76                |      | 40-140              | 38  |      | 40            |
| Phenanthrene                                                                                                    | 51               |      | 74                |      | 40-140              | 37  |      | 40            |
| Anthracene                                                                                                      | 55               |      | 79                |      | 40-140              | 36  |      | 40            |
| Fluoranthene                                                                                                    | 59               |      | 83                |      | 40-140              | 34  |      | 40            |
| Pyrene                                                                                                          | 58               |      | 83                |      | 26-127              | 35  |      | 40            |
| Benzo(a)anthracene                                                                                              | 47               |      | 70                |      | 40-140              | 39  |      | 40            |
| Chrysene                                                                                                        | 51               |      | 76                |      | 40-140              | 39  |      | 40            |
| Benzo(b)fluoranthene                                                                                            | 50               |      | 74                |      | 40-140              | 39  |      | 40            |
| Benzo(k)fluoranthene                                                                                            | 49               |      | 75                |      | 40-140              | 42  | Q    | 40            |
| Benzo(a)pyrene                                                                                                  | 52               |      | 79                |      | 40-140              | 41  | Q    | 40            |
| Indeno(1,2,3-cd)pyrene                                                                                          | 56               |      | 84                |      | 40-140              | 40  |      | 40            |
| Dibenzo(a,h)anthracene                                                                                          | 56               |      | 84                |      | 40-140              | 40  |      | 40            |
| Benzo(ghi)perylene                                                                                              | 53               |      | 79                |      | 40-140              | 39  |      | 40            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** B14 1ST HALF 2025 GW

**Lab Number:** L2523111

**Project Number:** 21010214

**Report Date:** 04/29/25

| <b>Parameter</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
|------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-03 Batch: WG2056800-2 WG2056800-3

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 36                       |             | 53                        |             | 21-120                         |
| Phenol-d6            | 30                       |             | 44                        |             | 10-120                         |
| Nitrobenzene-d5      | 68                       |             | 101                       |             | 23-120                         |
| 2-Fluorobiphenyl     | 47                       |             | 69                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 58                       |             | 84                        |             | 10-120                         |
| 4-Terphenyl-d14      | 53                       |             | 77                        |             | 41-149                         |

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

**Cooler Information**

|               |                     |
|---------------|---------------------|
| <b>Cooler</b> | <b>Custody Seal</b> |
| A             | Absent              |

**Container Information**

| <b>Container ID</b> | <b>Container Type</b>   | <b>Cooler</b> | <b>Initial pH</b> | <b>Final pH</b> | <b>Temp deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen Date/Time</b> | <b>Analysis(*)</b>               |
|---------------------|-------------------------|---------------|-------------------|-----------------|-------------------|-------------|-------------|-------------------------|----------------------------------|
| L2523111-01A        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-01B        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-01C        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-01D        | Amber 100ml unpreserved | A             | 12                | 12              | 2.1               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2523111-01E        | Amber 100ml unpreserved | A             | 12                | 12              | 2.1               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2523111-02A        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-02B        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-02C        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-02D        | Amber 100ml unpreserved | A             | 9                 | 9               | 2.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523111-02E        | Amber 100ml unpreserved | A             | 9                 | 9               | 2.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523111-03A        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-03B        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-03C        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-03D        | Amber 100ml unpreserved | A             | 11                | 11              | 2.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523111-03E        | Amber 100ml unpreserved | A             | 11                | 11              | 2.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523111-04A        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-04B        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-04C        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523111-04D        | Vial HCl preserved      | A             | NA                |                 | 2.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

## GLOSSARY

### Acronyms

|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| DL       | - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                |
| EDL      | - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).                                                                                         |
| EMPC     | - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.                                                                                                                                                                                                                               |
| EPA      | - Environmental Protection Agency.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| LCS      | - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                          |
| LCSD     | - Laboratory Control Sample Duplicate: Refer to LCS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |
| LFB      | - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                         |
| LOD      | - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                                                               |
| LOQ      | - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)<br><br>Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| MDL      | - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                          |
| MS       | - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.                                                                                                                                                                                                    |
| MSD      | - Matrix Spike Sample Duplicate: Refer to MS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NA       | - Not Applicable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| NC       | - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.                                                                                                                                                                                                                                                                                                                                                                           |
| NDPA/DPA | - N-Nitrosodiphenylamine/Diphenylamine.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| NI       | - Not Ignitable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| NP       | - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NR       | - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.                                                                                                                                                                                                                                                                                                                                                                    |
| RL       | - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                                                                                                                                   |
| RPD      | - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.                                                                  |
| SRM      | - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.                                                                                                                                                                                                                                                                                                                                                                     |
| STLP     | - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| TEF      | - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.                                                                                                                                                                                                                                                                                                                                                                                             |
| TEQ      | - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.                                                                                                                                                                                                                                                                                                                                                        |
| TIC      | - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.                                                                                                                                                                                                                                                                      |

Report Format: DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW  
**Project Number:** 21010214

**Lab Number:** L2523111  
**Report Date:** 04/29/25

### Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

### Terms

**Analytical Method:** Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

**Chlordane:** The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

**Difference:** With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

**Final pH:** As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

**Frozen Date/Time:** With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**Initial pH:** As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

**PAH Total:** With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

**PFAS Total:** With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**Total:** With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

### Data Qualifiers

- A** - Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F** - The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The lower value for the two columns has been reported due to obvious interference.
- J** - Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

**Report Format:** DU Report with 'J' Qualifiers



**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25**Data Qualifiers**

Identified Compounds (TICs). For calculated parameters, this represents that one or more values used in the calculation were estimated.

- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND** - Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ** - Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- R** - Analytical results are from sample re-analysis.
- RE** - Analytical results are from sample re-extraction.
- S** - Analytical results are from modified screening analysis.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

**Project Name:** B14 1ST HALF 2025 GW**Lab Number:** L2523111**Project Number:** 21010214**Report Date:** 04/29/25

## REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

## LIMITATION OF LIABILITIES

Pace Analytical Services performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Pace Analytical Services shall be to re-perform the work at it's own expense. In no event shall Pace Analytical Services be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Pace Analytical Services.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



**Pace Analytical Services LLC**Facility: **Northeast**Department: **Quality Assurance**Title: **Certificate/Approval Program Summary**ID No.: **17873**

Revision 27

Published Date: 01/24/2025

Page 1 of 2

**Certification Information**

The following analytes are not included in our Primary NELAP Scope of Accreditation:

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****EPA 624.1:** m/p-xylene, o-xylene, Naphthalene**EPA 625.1:** alpha-Terpineol**EPA 8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****SM 2540D:** TSS.**EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**MADEP-APH.****Nonpotable Water:** EPA RSK-175 Dissolved Gases**Biological Tissue Matrix:** EPA 3050B**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048****EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**Nonpotable Water:** EPA RSK-175 Dissolved Gases

The following test method is not included in our New Jersey Secondary NELAP Scope of Accreditation:

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Determination of Selected Perfluorinated Alkyl Substances by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry Isotope Dilution (via Alpha SOP 23528)**

The following analytes are included in our Massachusetts DEP Scope of Accreditation

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****Drinking Water****EPA 300.0:** Chloride, Nitrate-N, Fluoride, Sulfate; **EPA 353.2:** Nitrate-N, Nitrite-N; **SM4500NO3-F:** Nitrate-N, Nitrite-N; **SM4500F-C, SM4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B****EPA 524.2:** THMs and VOCs; **EPA 504.1:** EDB, DBCP.**Microbiology:** SM9215B; SM9223-P/A, SM9223B-Colilert-QT, SM9222D.**Non-Potable Water****SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH:** Ammonia-N and Kjeldahl-N, **EPA 350.1:** Ammonia-N, **LCHAT 10-107-06-1-B:** Ammonia-N, **EPA 351.1, SM4500NO3-F, EPA 353.2:** Nitrate-N, **SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300:** Chloride, Sulfate, Nitrate.**EPA 624.1:** Volatile Halocarbons & Aromatics,**EPA 608.3:** Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs**EPA 625.1:** SVOC (Acid/Base/Neutral Extractables).**Microbiology:** SM9223B-Colilert-QT; Enterolert-QT, EPA 1600, EPA 1603, SM9222D.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Drinking Water****EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1 Hg. EPA 522, EPA 537.1.****Non-Potable Water****EPA 200.7:** Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.**EPA 200.8:** Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.**EPA 245.1 Hg.****SM2340B**

**Pace Analytical Services LLC**ID No.: **17873**Facility: **Northeast**

Revision 27

Department: **Quality Assurance**

Published Date: 01/24/2025

Title: **Certificate/Approval Program Summary**

Page 2 of 2

**Certification IDs:****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581**

CT PH-0826, IL 200077, IN C-MA-03, KY JY98045, ME MA00086, MD 348, MA M-MA086, NH 2064, NJ MA935, NY 11148, NC (DW) 25700, NC (NPW/SCM) 666, OR MA-1316, PA 68-03671, RI LAO00065, TX T104704476, VT VT-0935, VA 460195

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048**

CT PH-0825, ANAB/DoD L2474, IL 200081, IN C-MA-04, KY KY98046, LA 3090, ME MA00030, MI 9110, MN 025-999-495, NH 2062, NJ MA015, NY 11627, NC (NPW/SCM) 685, OR MA-0262, PA 68-02089, RI LAO00299, TX T-104704419, VT VT-0015, VA 460194, WA C954

**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048**

ANAB/DoD L2474, ME MA01156, MN 025-999-498, NH 2249, NJ MA025, NY 12191, OR 4203, TX T104704583, VA 460311, WA C1104.

---

For a complete listing of analytes and methods, please contact your Project Manager.





## ANALYTICAL REPORT

|                 |                                                                             |
|-----------------|-----------------------------------------------------------------------------|
| Lab Number:     | L2523470                                                                    |
| Client:         | Tradepoint Atlantic<br>1600 Sparrows Point Boulevard<br>Baltimore, MD 21219 |
| ATTN:           | Matthew Newman                                                              |
| Phone:          | (443) 649-5073                                                              |
| Project Name:   | TMC SURFACE WATER SAMPLING                                                  |
| Project Number: | Not Specified                                                               |
| Report Date:    | 04/29/25                                                                    |

The original project report/data package is held by Pace Analytical Services. This report/data package is paginated and should be reproduced only in its entirety. Pace Analytical Services holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0826), IL (200077), IN (C-MA-03), KY (KY98045), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), OR (MA-1316), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #525-23-122-91930A1).

---

Eight Walkup Drive, Westborough, MA 01581-1019  
508-898-9220 (Fax) 508-898-9193 800-624-9220 - [www.alphalab.com](http://www.alphalab.com)





**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

| Lab<br>Sample ID | Client ID       | Matrix | Sample<br>Location | Collection<br>Date/Time | Receive Date |
|------------------|-----------------|--------|--------------------|-------------------------|--------------|
| L2523470-01      | TMC-OUTLET      | WATER  | B16                | 04/16/25 09:45          | 04/16/25     |
| L2523470-02      | TMC-TM04        | WATER  | B16                | 04/16/25 10:15          | 04/16/25     |
| L2523470-03      | TMC-BEND        | WATER  | B16                | 04/16/25 11:45          | 04/16/25     |
| L2523470-04      | TMC-RAIL-BRIDGE | WATER  | B16                | 04/16/25 12:00          | 04/16/25     |
| L2523470-05      | TB-001          | WATER  | B16                | 04/16/25 00:00          | 04/16/25     |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

### Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Pace Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments and solids are reported on a dry weight basis unless otherwise noted. Tissues are reported "as received" or on a wet weight basis, unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

**HOLD POLICY** - For samples submitted on hold, Pace's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Pace Project Manager and made arrangements for Pace to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.

---

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

**Case Narrative (continued)**

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:



Caitlin Walukevich

Title: Technical Director/Representative

Date: 04/29/25

# ORGANICS

# VOLATILES

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-01

Date Collected: 04/16/25 09:45

Client ID: TMC-OUTLET

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 04/24/25 11:20

Analyst: MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | 3.7    | J         | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 5.2    |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | 0.42   | J         | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.42   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-01

Date Collected: 04/16/25 09:45

Client ID: TMC-OUTLET

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 119        |           | 70-130              |
| Toluene-d8            | 107        |           | 70-130              |
| 4-Bromofluorobenzene  | 97         |           | 70-130              |
| Dibromofluoromethane  | 106        |           | 70-130              |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

Lab ID: L2523470-01  
 Client ID: TMC-OUTLET  
 Sample Location: B16

Date Collected: 04/16/25 09:45  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/24/25 11:20  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 114        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 92         |           | 70-130              |                 |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-02

Date Collected: 04/16/25 10:15

Client ID: TMC-TM04

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 04/24/25 10:57

Analyst: MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 4.2    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.30   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | 0.20   | J         | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-02

Date Collected: 04/16/25 10:15

Client ID: TMC-TM04

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 117        |           | 70-130              |
| Toluene-d8            | 108        |           | 70-130              |
| 4-Bromofluorobenzene  | 95         |           | 70-130              |
| Dibromofluoromethane  | 107        |           | 70-130              |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

**SAMPLE RESULTS**

Lab ID: L2523470-02  
 Client ID: TMC-TM04  
 Sample Location: B16

Date Collected: 04/16/25 10:15  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/24/25 10:57  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 112        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 91         |           | 70-130              |                 |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-03

Date Collected: 04/16/25 11:45

Client ID: TMC-BEND

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 04/24/25 10:33

Analyst: MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 4.1    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.17   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-03

Date Collected: 04/16/25 11:45

Client ID: TMC-BEND

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 117        |           | 70-130              |
| Toluene-d8            | 97         |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 108        |           | 70-130              |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-03

Date Collected: 04/16/25 11:45

Client ID: TMC-BEND

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D-SIM(M)

Analytical Date: 04/24/25 10:33

Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 112        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 92         |           | 70-130              |                 |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-04  
 Client ID: TMC-RAIL-BRIDGE  
 Sample Location: B16

Date Collected: 04/16/25 12:00  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:

Matrix: Water  
 Analytical Method: 1,8260D  
 Analytical Date: 04/24/25 10:09  
 Analyst: MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | 3.1    | J         | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | 0.39   | J         | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | 0.29   | J         | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-04  
 Client ID: TMC-RAIL-BRIDGE  
 Sample Location: B16

Date Collected: 04/16/25 12:00  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 118        |           | 70-130              |
| Toluene-d8            | 108        |           | 70-130              |
| 4-Bromofluorobenzene  | 96         |           | 70-130              |
| Dibromofluoromethane  | 107        |           | 70-130              |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-04  
 Client ID: TMC-RAIL-BRIDGE  
 Sample Location: B16

Date Collected: 04/16/25 12:00  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:

Matrix: Water  
 Analytical Method: 1,8260D-SIM(M)  
 Analytical Date: 04/24/25 10:09  
 Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 112        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 91         |           | 70-130              |                 |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-05

Date Collected: 04/16/25 00:00

Client ID: TB-001

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D

Analytical Date: 04/24/25 09:46

Analyst: MCM

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Dichlorodifluoromethane                      | ND     |           | ug/l  | 5.0  | 0.24 | 1               |
| Chloromethane                                | ND     |           | ug/l  | 2.5  | 0.20 | 1               |
| Vinyl chloride                               | ND     |           | ug/l  | 1.0  | 0.07 | 1               |
| Bromomethane                                 | ND     |           | ug/l  | 1.0  | 0.26 | 1               |
| Chloroethane                                 | ND     |           | ug/l  | 1.0  | 0.13 | 1               |
| Trichlorofluoromethane                       | ND     |           | ug/l  | 2.5  | 0.16 | 1               |
| 1,1-Dichloroethene                           | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| Carbon disulfide                             | ND     |           | ug/l  | 5.0  | 0.30 | 1               |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane        | ND     |           | ug/l  | 2.5  | 0.15 | 1               |
| Methylene chloride                           | ND     |           | ug/l  | 2.5  | 0.68 | 1               |
| Acetone                                      | ND     |           | ug/l  | 5.0  | 1.5  | 1               |
| trans-1,2-Dichloroethene                     | ND     |           | ug/l  | 0.75 | 0.16 | 1               |
| Methyl Acetate                               | ND     |           | ug/l  | 2.0  | 0.23 | 1               |
| Methyl tert butyl ether                      | ND     |           | ug/l  | 1.0  | 0.17 | 1               |
| 1,1-Dichloroethane                           | ND     |           | ug/l  | 0.75 | 0.21 | 1               |
| cis-1,2-Dichloroethene                       | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| Cyclohexane                                  | ND     |           | ug/l  | 10   | 0.27 | 1               |
| Chloroform                                   | ND     |           | ug/l  | 0.75 | 0.22 | 1               |
| Carbon tetrachloride                         | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| 1,1,1-Trichloroethane                        | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 2-Butanone                                   | ND     |           | ug/l  | 5.0  | 1.9  | 1               |
| Benzene                                      | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,2-Dichloroethane                           | ND     |           | ug/l  | 0.50 | 0.13 | 1               |
| Trichloroethene                              | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 1,2-Dichloropropane                          | ND     |           | ug/l  | 1.0  | 0.14 | 1               |
| Bromodichloromethane                         | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| cis-1,3-Dichloropropene                      | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| Toluene                                      | ND     |           | ug/l  | 0.75 | 0.20 | 1               |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-05

Date Collected: 04/16/25 00:00

Client ID: TB-001

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

| Parameter                                    | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|----------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab |        |           |       |      |      |                 |
| Tetrachloroethene                            | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| 4-Methyl-2-pentanone                         | ND     |           | ug/l  | 5.0  | 0.42 | 1               |
| trans-1,3-Dichloropropene                    | ND     |           | ug/l  | 0.50 | 0.16 | 1               |
| 1,3-Dichloropropene, Total                   | ND     |           | ug/l  | 0.50 | 0.14 | 1               |
| 1,1,2-Trichloroethane                        | ND     |           | ug/l  | 0.75 | 0.14 | 1               |
| Dibromochloromethane                         | ND     |           | ug/l  | 0.50 | 0.15 | 1               |
| 1,2-Dibromoethane                            | ND     |           | ug/l  | 2.0  | 0.19 | 1               |
| 2-Hexanone                                   | ND     |           | ug/l  | 5.0  | 0.52 | 1               |
| Chlorobenzene                                | ND     |           | ug/l  | 0.50 | 0.18 | 1               |
| Ethylbenzene                                 | ND     |           | ug/l  | 0.50 | 0.17 | 1               |
| p/m-Xylene                                   | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| o-Xylene                                     | ND     |           | ug/l  | 1.0  | 0.39 | 1               |
| Xylenes, Total                               | ND     |           | ug/l  | 1.0  | 0.33 | 1               |
| Styrene                                      | ND     |           | ug/l  | 1.0  | 0.36 | 1               |
| Bromoform                                    | ND     |           | ug/l  | 2.0  | 0.25 | 1               |
| Isopropylbenzene                             | ND     |           | ug/l  | 0.50 | 0.19 | 1               |
| 1,3-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,4-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.19 | 1               |
| 1,2-Dichlorobenzene                          | ND     |           | ug/l  | 2.5  | 0.18 | 1               |
| 1,2-Dibromo-3-chloropropane                  | ND     |           | ug/l  | 2.5  | 0.35 | 1               |
| 1,2,4-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.22 | 1               |
| 1,2,3-Trichlorobenzene                       | ND     |           | ug/l  | 2.5  | 0.23 | 1               |

| Surrogate             | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 118        |           | 70-130              |
| Toluene-d8            | 97         |           | 70-130              |
| 4-Bromofluorobenzene  | 95         |           | 70-130              |
| Dibromofluoromethane  | 109        |           | 70-130              |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-05

Date Collected: 04/16/25 00:00

Client ID: TB-001

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Analytical Method: 1,8260D-SIM(M)

Analytical Date: 04/24/25 09:46

Analyst: MCM

| Parameter                                        | Result | Qualifier | Units      | RL        | MDL                 | Dilution Factor |
|--------------------------------------------------|--------|-----------|------------|-----------|---------------------|-----------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab |        |           |            |           |                     |                 |
| 1,1,2,2-Tetrachloroethane                        | ND     |           | ug/l       | 0.050     | 0.006               | 1               |
| Surrogate                                        |        |           | % Recovery | Qualifier | Acceptance Criteria |                 |
| 1,2-Dichloroethane-d4                            |        |           | 112        |           | 70-130              |                 |
| 4-Bromofluorobenzene                             |        |           | 91         |           | 70-130              |                 |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D  
 Analytical Date: 04/24/25 05:49  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2059503-5 |        |           |       |      |      |
| Dichlorodifluoromethane                                                              | ND     |           | ug/l  | 5.0  | 0.24 |
| Chloromethane                                                                        | ND     |           | ug/l  | 2.5  | 0.20 |
| Vinyl chloride                                                                       | ND     |           | ug/l  | 1.0  | 0.07 |
| Bromomethane                                                                         | ND     |           | ug/l  | 1.0  | 0.26 |
| Chloroethane                                                                         | ND     |           | ug/l  | 1.0  | 0.13 |
| Trichlorofluoromethane                                                               | ND     |           | ug/l  | 2.5  | 0.16 |
| 1,1-Dichloroethene                                                                   | ND     |           | ug/l  | 0.50 | 0.17 |
| Carbon disulfide                                                                     | ND     |           | ug/l  | 5.0  | 0.30 |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                | ND     |           | ug/l  | 2.5  | 0.15 |
| Methylene chloride                                                                   | ND     |           | ug/l  | 2.5  | 0.68 |
| Acetone                                                                              | ND     |           | ug/l  | 5.0  | 1.5  |
| trans-1,2-Dichloroethene                                                             | ND     |           | ug/l  | 0.75 | 0.16 |
| Methyl Acetate                                                                       | ND     |           | ug/l  | 2.0  | 0.23 |
| Methyl tert butyl ether                                                              | ND     |           | ug/l  | 1.0  | 0.17 |
| 1,1-Dichloroethane                                                                   | ND     |           | ug/l  | 0.75 | 0.21 |
| cis-1,2-Dichloroethene                                                               | ND     |           | ug/l  | 0.50 | 0.19 |
| Cyclohexane                                                                          | ND     |           | ug/l  | 10   | 0.27 |
| Chloroform                                                                           | ND     |           | ug/l  | 0.75 | 0.22 |
| Carbon tetrachloride                                                                 | ND     |           | ug/l  | 0.50 | 0.13 |
| 1,1,1-Trichloroethane                                                                | ND     |           | ug/l  | 0.50 | 0.16 |
| 2-Butanone                                                                           | ND     |           | ug/l  | 5.0  | 1.9  |
| Benzene                                                                              | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,2-Dichloroethane                                                                   | ND     |           | ug/l  | 0.50 | 0.13 |
| Trichloroethene                                                                      | ND     |           | ug/l  | 0.50 | 0.18 |
| 1,2-Dichloropropane                                                                  | ND     |           | ug/l  | 1.0  | 0.14 |
| Bromodichloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.19 |
| cis-1,3-Dichloropropene                                                              | ND     |           | ug/l  | 0.50 | 0.14 |
| Toluene                                                                              | ND     |           | ug/l  | 0.75 | 0.20 |
| Tetrachloroethene                                                                    | ND     |           | ug/l  | 0.50 | 0.18 |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D  
 Analytical Date: 04/24/25 05:49  
 Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL   | MDL  |
|--------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2059503-5 |        |           |       |      |      |
| 4-Methyl-2-pentanone                                                                 | ND     |           | ug/l  | 5.0  | 0.42 |
| trans-1,3-Dichloropropene                                                            | ND     |           | ug/l  | 0.50 | 0.16 |
| 1,3-Dichloropropene, Total                                                           | ND     |           | ug/l  | 0.50 | 0.14 |
| 1,1,2-Trichloroethane                                                                | ND     |           | ug/l  | 0.75 | 0.14 |
| Dibromochloromethane                                                                 | ND     |           | ug/l  | 0.50 | 0.15 |
| 1,2-Dibromoethane                                                                    | ND     |           | ug/l  | 2.0  | 0.19 |
| 2-Hexanone                                                                           | ND     |           | ug/l  | 5.0  | 0.52 |
| Chlorobenzene                                                                        | ND     |           | ug/l  | 0.50 | 0.18 |
| Ethylbenzene                                                                         | ND     |           | ug/l  | 0.50 | 0.17 |
| p/m-Xylene                                                                           | ND     |           | ug/l  | 1.0  | 0.33 |
| o-Xylene                                                                             | ND     |           | ug/l  | 1.0  | 0.39 |
| Xylenes, Total                                                                       | ND     |           | ug/l  | 1.0  | 0.33 |
| Styrene                                                                              | ND     |           | ug/l  | 1.0  | 0.36 |
| Bromoform                                                                            | ND     |           | ug/l  | 2.0  | 0.25 |
| Isopropylbenzene                                                                     | ND     |           | ug/l  | 0.50 | 0.19 |
| 1,1,2,2-Tetrachloroethane                                                            | ND     |           | ug/l  | 0.50 | 0.17 |
| 1,3-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,4-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.19 |
| 1,2-Dichlorobenzene                                                                  | ND     |           | ug/l  | 2.5  | 0.18 |
| 1,2-Dibromo-3-chloropropane                                                          | ND     |           | ug/l  | 2.5  | 0.35 |
| 1,2,4-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.22 |
| 1,2,3-Trichlorobenzene                                                               | ND     |           | ug/l  | 2.5  | 0.23 |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
Analytical Date: 04/24/25 05:49  
Analyst: MCM

| Parameter                                                                            | Result | Qualifier | Units | RL | MDL |
|--------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG2059503-5 |        |           |       |    |     |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 108       |           | 70-130                 |
| Toluene-d8            | 108       |           | 70-130                 |
| 4-Bromofluorobenzene  | 96        |           | 70-130                 |
| Dibromofluoromethane  | 104       |           | 70-130                 |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D-SIM(M)  
Analytical Date: 04/24/25 05:49  
Analyst: MCM

| Parameter                                                                                | Result | Qualifier | Units | RL    | MDL   |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-------|-------|
| Volatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-05 Batch: WG2059509-5 |        |           |       |       |       |
| 1,1,1,2-Tetrachloroethane                                                                | ND     |           | ug/l  | 0.050 | 0.006 |

| Surrogate             | %Recovery | Qualifier | Acceptance<br>Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 110       |           | 70-130                 |
| 4-Bromofluorobenzene  | 92        |           | 70-130                 |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Project Number:** Not Specified

**Lab Number:** L2523470

**Report Date:** 04/29/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2059503-3 WG2059503-4 |                  |      |                   |      |                     |     |      |               |
| Dichlorodifluoromethane                                                                                 | 110              |      | 120               |      | 36-147              | 9   |      | 20            |
| Chloromethane                                                                                           | 110              |      | 110               |      | 64-130              | 0   |      | 20            |
| Vinyl chloride                                                                                          | 100              |      | 100               |      | 55-140              | 0   |      | 20            |
| Bromomethane                                                                                            | 100              |      | 100               |      | 39-139              | 0   |      | 20            |
| Chloroethane                                                                                            | 110              |      | 110               |      | 55-138              | 0   |      | 20            |
| Trichlorofluoromethane                                                                                  | 100              |      | 100               |      | 62-150              | 0   |      | 20            |
| 1,1-Dichloroethene                                                                                      | 97               |      | 100               |      | 61-145              | 3   |      | 20            |
| Carbon disulfide                                                                                        | 100              |      | 100               |      | 51-130              | 0   |      | 20            |
| 1,1,2-Trichloro-1,2,2-Trifluoroethane                                                                   | 100              |      | 110               |      | 70-130              | 10  |      | 20            |
| Methylene chloride                                                                                      | 100              |      | 99                |      | 70-130              | 1   |      | 20            |
| Acetone                                                                                                 | 120              |      | 120               |      | 58-148              | 0   |      | 20            |
| trans-1,2-Dichloroethene                                                                                | 98               |      | 100               |      | 70-130              | 2   |      | 20            |
| Methyl Acetate                                                                                          | 110              |      | 110               |      | 70-130              | 0   |      | 20            |
| Methyl tert butyl ether                                                                                 | 100              |      | 110               |      | 63-130              | 10  |      | 20            |
| 1,1-Dichloroethane                                                                                      | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| cis-1,2-Dichloroethene                                                                                  | 94               |      | 97                |      | 70-130              | 3   |      | 20            |
| Cyclohexane                                                                                             | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Chloroform                                                                                              | 100              |      | 110               |      | 70-130              | 10  |      | 20            |
| Carbon tetrachloride                                                                                    | 110              |      | 100               |      | 63-132              | 10  |      | 20            |
| 1,1,1-Trichloroethane                                                                                   | 100              |      | 110               |      | 67-130              | 10  |      | 20            |
| 2-Butanone                                                                                              | 110              |      | 100               |      | 63-138              | 10  |      | 20            |
| Benzene                                                                                                 | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| 1,2-Dichloroethane                                                                                      | 110              |      | 110               |      | 70-130              | 0   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Project Number:** Not Specified

**Lab Number:** L2523470

**Report Date:** 04/29/25

| Parameter                                                                                               | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|---------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2059503-3 WG2059503-4 |                  |      |                   |      |                     |     |      |               |
| Trichloroethene                                                                                         | 98               |      | 99                |      | 70-130              | 1   |      | 20            |
| 1,2-Dichloropropane                                                                                     | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| Bromodichloromethane                                                                                    | 100              |      | 100               |      | 67-130              | 0   |      | 20            |
| cis-1,3-Dichloropropene                                                                                 | 89               |      | 90                |      | 70-130              | 1   |      | 20            |
| Toluene                                                                                                 | 110              |      | 100               |      | 70-130              | 10  |      | 20            |
| Tetrachloroethene                                                                                       | 120              |      | 110               |      | 70-130              | 9   |      | 20            |
| 4-Methyl-2-pentanone                                                                                    | 100              |      | 94                |      | 59-130              | 6   |      | 20            |
| trans-1,3-Dichloropropene                                                                               | 100              |      | 90                |      | 70-130              | 11  |      | 20            |
| 1,1,2-Trichloroethane                                                                                   | 120              |      | 100               |      | 70-130              | 18  |      | 20            |
| Dibromochloromethane                                                                                    | 110              |      | 100               |      | 63-130              | 10  |      | 20            |
| 1,2-Dibromoethane                                                                                       | 110              |      | 100               |      | 70-130              | 10  |      | 20            |
| 2-Hexanone                                                                                              | 110              |      | 110               |      | 57-130              | 0   |      | 20            |
| Chlorobenzene                                                                                           | 110              |      | 98                |      | 75-130              | 12  |      | 20            |
| Ethylbenzene                                                                                            | 110              |      | 100               |      | 70-130              | 10  |      | 20            |
| p/m-Xylene                                                                                              | 110              |      | 100               |      | 70-130              | 10  |      | 20            |
| o-Xylene                                                                                                | 105              |      | 95                |      | 70-130              | 10  |      | 20            |
| Styrene                                                                                                 | 110              |      | 100               |      | 70-130              | 10  |      | 20            |
| Bromoform                                                                                               | 100              |      | 100               |      | 54-136              | 0   |      | 20            |
| Isopropylbenzene                                                                                        | 94               |      | 88                |      | 70-130              | 7   |      | 20            |
| 1,1,2,2-Tetrachloroethane                                                                               | 98               |      | 97                |      | 67-130              | 1   |      | 20            |
| 1,3-Dichlorobenzene                                                                                     | 98               |      | 98                |      | 70-130              | 0   |      | 20            |
| 1,4-Dichlorobenzene                                                                                     | 100              |      | 100               |      | 70-130              | 0   |      | 20            |
| 1,2-Dichlorobenzene                                                                                     | 100              |      | 100               |      | 70-130              | 0   |      | 20            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Lab Number:** L2523470

**Project Number:** Not Specified

**Report Date:** 04/29/25

| <b>Parameter</b>                                                                                        | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|---------------------------------------------------------------------------------------------------------|--------------------------|-------------|--------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-05 Batch: WG2059503-3 WG2059503-4 |                          |             |                          |             |                             |            |             |                       |
| 1,2-Dibromo-3-chloropropane                                                                             | 98                       |             | 100                      |             | 41-144                      | 2          |             | 20                    |
| 1,2,4-Trichlorobenzene                                                                                  | 94                       |             | 96                       |             | 70-130                      | 2          |             | 20                    |
| 1,2,3-Trichlorobenzene                                                                                  | 100                      |             | 100                      |             | 70-130                      | 0          |             | 20                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|--------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 111                      |             | 112                      |             | 70-130                         |
| Toluene-d8            | 110                      |             | 98                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 92                       |             | 93                       |             | 70-130                         |
| Dibromofluoromethane  | 98                       |             | 98                       |             | 70-130                         |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Lab Number:** L2523470

**Project Number:** Not Specified

**Report Date:** 04/29/25

| <b>Parameter</b>                                                                                            | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Volatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-05 Batch: WG2059509-3 WG2059509-4 |                          |             |                           |             |                             |            |             |                       |
| 1,1,1,2-Tetrachloroethane                                                                                   | 97                       |             | 105                       |             | 70-130                      | 8          |             | 25                    |

| <b>Surrogate</b>      | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|-----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 1,2-Dichloroethane-d4 | 106                      |             | 109                       |             | 70-130                         |
| 4-Bromofluorobenzene  | 91                       |             | 90                        |             | 70-130                         |

# SEMIVOLATILES

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-01

Date Collected: 04/16/25 09:45

Client ID: TMC-OUTLET

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Extraction Method: EPA 3510C

Analytical Method: 1,8270E

Extraction Date: 04/22/25 02:15

Analytical Date: 04/23/25 06:15

Analyst: JG

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-01

Date Collected: 04/16/25 09:45

Client ID: TMC-OUTLET

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 16         | Q         | 21-120              |
| Phenol-d6            | 14         |           | 10-120              |
| Nitrobenzene-d5      | 60         |           | 23-120              |
| 2-Fluorobiphenyl     | 59         |           | 15-120              |
| 2,4,6-Tribromophenol | 23         |           | 10-120              |
| 4-Terphenyl-d14      | 65         |           | 41-149              |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-01

Date Collected: 04/16/25 09:45

Client ID: TMC-OUTLET

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Extraction Method: EPA 3510C

Analytical Method: 1,8270E-SIM

Extraction Date: 04/22/25 02:15

Analytical Date: 04/22/25 11:40

Analyst: SLR

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.06   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.08   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.16   |           | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | 0.15   |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | 0.06   |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | 0.05   | J         | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | 0.03   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | 0.03   | J         | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 12         | Q         | 21-120              |
| Phenol-d6            | 12         |           | 10-120              |
| Nitrobenzene-d5      | 56         |           | 23-120              |
| 2-Fluorobiphenyl     | 44         |           | 15-120              |
| 2,4,6-Tribromophenol | 23         |           | 10-120              |
| 4-Terphenyl-d14      | 45         |           | 41-149              |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-02

Date Collected: 04/16/25 10:15

Client ID: TMC-TM04

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Extraction Method: EPA 3510C

Analytical Method: 1,8270E

Extraction Date: 04/22/25 02:15

Analytical Date: 04/23/25 06:38

Analyst: JG

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-02

Date Collected: 04/16/25 10:15

Client ID: TMC-TM04

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 16         | Q         | 21-120              |
| Phenol-d6            | 17         |           | 10-120              |
| Nitrobenzene-d5      | 82         |           | 23-120              |
| 2-Fluorobiphenyl     | 85         |           | 15-120              |
| 2,4,6-Tribromophenol | 27         |           | 10-120              |
| 4-Terphenyl-d14      | 96         |           | 41-149              |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-02

Date Collected: 04/16/25 10:15

Client ID: TMC-TM04

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Extraction Method: EPA 3510C

Analytical Method: 1,8270E-SIM

Extraction Date: 04/22/25 02:15

Analytical Date: 04/22/25 11:57

Analyst: SLR

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.04   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.05   | J         | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.03   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | 0.03   | J         | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 14         | Q         | 21-120              |
| Phenol-d6            | 16         |           | 10-120              |
| Nitrobenzene-d5      | 81         |           | 23-120              |
| 2-Fluorobiphenyl     | 66         |           | 15-120              |
| 2,4,6-Tribromophenol | 30         |           | 10-120              |
| 4-Terphenyl-d14      | 68         |           | 41-149              |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-03

Date Collected: 04/16/25 11:45

Client ID: TMC-BEND

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Extraction Method: EPA 3510C

Analytical Method: 1,8270E

Extraction Date: 04/22/25 02:15

Analytical Date: 04/23/25 07:00

Analyst: JG

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-03

Date Collected: 04/16/25 11:45

Client ID: TMC-BEND

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 15         | Q         | 21-120              |
| Phenol-d6            | 20         |           | 10-120              |
| Nitrobenzene-d5      | 92         |           | 23-120              |
| 2-Fluorobiphenyl     | 94         |           | 15-120              |
| 2,4,6-Tribromophenol | 24         |           | 10-120              |
| 4-Terphenyl-d14      | 106        |           | 41-149              |



**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-03

Date Collected: 04/16/25 11:45

Client ID: TMC-BEND

Date Received: 04/16/25

Sample Location: B16

Field Prep: Not Specified

Sample Depth:

Matrix: Water

Extraction Method: EPA 3510C

Analytical Method: 1,8270E-SIM

Extraction Date: 04/22/25 02:15

Analytical Date: 04/22/25 12:14

Analyst: SLR

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.03   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.04   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.04   | J         | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.03   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 13         | Q         | 21-120              |
| Phenol-d6            | 16         |           | 10-120              |
| Nitrobenzene-d5      | 85         |           | 23-120              |
| 2-Fluorobiphenyl     | 69         |           | 15-120              |
| 2,4,6-Tribromophenol | 25         |           | 10-120              |
| 4-Terphenyl-d14      | 68         |           | 41-149              |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-04  
 Client ID: TMC-RAIL-BRIDGE  
 Sample Location: B16

Date Collected: 04/16/25 12:00  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:

Matrix: Water  
 Analytical Method: 1,8270E  
 Analytical Date: 04/23/25 07:24  
 Analyst: JG

Extraction Method: EPA 3510C  
 Extraction Date: 04/22/25 02:15

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Benzaldehyde                                     | ND     |           | ug/l  | 5.0 | 1.1  | 1               |
| Phenol                                           | ND     |           | ug/l  | 5.0 | 0.35 | 1               |
| Bis(2-chloroethyl)ether                          | ND     |           | ug/l  | 2.0 | 0.39 | 1               |
| 2-Chlorophenol                                   | ND     |           | ug/l  | 2.0 | 0.65 | 1               |
| 2-Methylphenol                                   | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Bis(2-chloroisopropyl)ether                      | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| Acetophenone                                     | ND     |           | ug/l  | 5.0 | 0.92 | 1               |
| n-Nitrosodi-n-propylamine                        | ND     |           | ug/l  | 5.0 | 0.91 | 1               |
| 3-Methylphenol/4-Methylphenol                    | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| Hexachloroethane                                 | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| Nitrobenzene                                     | ND     |           | ug/l  | 1.4 | 0.20 | 1               |
| Isophorone                                       | ND     |           | ug/l  | 5.0 | 0.86 | 1               |
| 2,4-Dimethylphenol                               | ND     |           | ug/l  | 5.0 | 2.0  | 1               |
| Bis(2-chloroethoxy)methane                       | ND     |           | ug/l  | 5.0 | 0.84 | 1               |
| 2,4-Dichlorophenol                               | ND     |           | ug/l  | 5.0 | 1.7  | 1               |
| Naphthalene                                      | ND     |           | ug/l  | 2.0 | 0.54 | 1               |
| 4-Chloroaniline                                  | ND     |           | ug/l  | 3.7 | 0.47 | 1               |
| Hexachlorobutadiene                              | ND     |           | ug/l  | 2.0 | 0.36 | 1               |
| Caprolactam                                      | ND     |           | ug/l  | 10  | 1.2  | 1               |
| 2-Methylnaphthalene                              | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Hexachlorocyclopentadiene                        | ND     |           | ug/l  | 20  | 1.2  | 1               |
| 1,2,4,5-Tetrachlorobenzene                       | ND     |           | ug/l  | 1.7 | 0.24 | 1               |
| 2,4,6-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| 2,4,5-Trichlorophenol                            | ND     |           | ug/l  | 5.0 | 2.1  | 1               |
| Biphenyl                                         | ND     |           | ug/l  | 2.0 | 0.20 | 1               |
| 2-Chloronaphthalene                              | ND     |           | ug/l  | 2.0 | 0.35 | 1               |
| 2-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.0  | 1               |
| 2,6-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.84 | 1               |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-04  
 Client ID: TMC-RAIL-BRIDGE  
 Sample Location: B16

Date Collected: 04/16/25 12:00  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:

| Parameter                                        | Result | Qualifier | Units | RL  | MDL  | Dilution Factor |
|--------------------------------------------------|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab |        |           |       |     |      |                 |
| Acenaphthylene                                   | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Acenaphthene                                     | ND     |           | ug/l  | 2.0 | 0.40 | 1               |
| 2,4-Dinitrophenol                                | ND     |           | ug/l  | 20  | 5.4  | 1               |
| 2,4-Dinitrotoluene                               | ND     |           | ug/l  | 5.0 | 0.54 | 1               |
| 2,3,4,6-Tetrachlorophenol                        | ND     |           | ug/l  | 5.0 | 2.2  | 1               |
| Diethyl phthalate                                | ND     |           | ug/l  | 5.0 | 0.76 | 1               |
| Fluorene                                         | ND     |           | ug/l  | 2.0 | 0.44 | 1               |
| 4-Nitroaniline                                   | ND     |           | ug/l  | 5.0 | 1.4  | 1               |
| NDPA/DPA                                         | ND     |           | ug/l  | 2.0 | 0.92 | 1               |
| Hexachlorobenzene                                | ND     |           | ug/l  | 2.0 | 0.45 | 1               |
| Pentachlorophenol                                | ND     |           | ug/l  | 10  | 2.5  | 1               |
| Phenanthrene                                     | ND     |           | ug/l  | 2.0 | 0.42 | 1               |
| Anthracene                                       | ND     |           | ug/l  | 2.0 | 0.47 | 1               |
| Carbazole                                        | ND     |           | ug/l  | 2.0 | 0.31 | 1               |
| Di-n-butylphthalate                              | ND     |           | ug/l  | 5.0 | 0.96 | 1               |
| Fluoranthene                                     | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| Pyrene                                           | ND     |           | ug/l  | 2.0 | 0.41 | 1               |
| 3,3'-Dichlorobenzidine                           | ND     |           | ug/l  | 5.0 | 1.8  | 1               |
| Benzo(a)anthracene                               | ND     |           | ug/l  | 2.0 | 0.32 | 1               |
| Chrysene                                         | ND     |           | ug/l  | 1.5 | 0.22 | 1               |
| Bis(2-ethylhexyl)phthalate                       | ND     |           | ug/l  | 3.0 | 1.4  | 1               |
| Di-n-octylphthalate                              | ND     |           | ug/l  | 5.0 | 2.3  | 1               |
| Benzo(b)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.53 | 1               |
| Benzo(k)fluoranthene                             | ND     |           | ug/l  | 2.0 | 0.62 | 1               |
| Benzo(a)pyrene                                   | ND     |           | ug/l  | 2.0 | 0.37 | 1               |
| Indeno(1,2,3-cd)pyrene                           | ND     |           | ug/l  | 2.0 | 0.48 | 1               |
| Dibenzo(a,h)anthracene                           | ND     |           | ug/l  | 2.0 | 0.29 | 1               |
| Benzo(ghi)perylene                               | ND     |           | ug/l  | 2.0 | 0.37 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 22         |           | 21-120              |
| Phenol-d6            | 23         |           | 10-120              |
| Nitrobenzene-d5      | 82         |           | 23-120              |
| 2-Fluorobiphenyl     | 86         |           | 15-120              |
| 2,4,6-Tribromophenol | 39         |           | 10-120              |
| 4-Terphenyl-d14      | 103        |           | 41-149              |





**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**SAMPLE RESULTS**

Lab ID: L2523470-04  
 Client ID: TMC-RAIL-BRIDGE  
 Sample Location: B16

Date Collected: 04/16/25 12:00  
 Date Received: 04/16/25  
 Field Prep: Not Specified

Sample Depth:

Matrix: Water  
 Analytical Method: 1,8270E-SIM  
 Analytical Date: 04/22/25 12:31  
 Analyst: SLR

Extraction Method: EPA 3510C  
 Extraction Date: 04/22/25 02:15

| Parameter                                            | Result | Qualifier | Units | RL   | MDL  | Dilution Factor |
|------------------------------------------------------|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab |        |           |       |      |      |                 |
| Naphthalene                                          | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| 2-Methylnaphthalene                                  | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Acenaphthylene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Acenaphthene                                         | 0.09   | J         | ug/l  | 0.10 | 0.02 | 1               |
| Fluorene                                             | 0.10   |           | ug/l  | 0.10 | 0.03 | 1               |
| Phenanthrene                                         | 0.10   |           | ug/l  | 0.05 | 0.04 | 1               |
| Anthracene                                           | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Fluoranthene                                         | 0.06   | J         | ug/l  | 0.10 | 0.03 | 1               |
| Pyrene                                               | ND     |           | ug/l  | 0.10 | 0.04 | 1               |
| Benzo(a)anthracene                                   | 0.03   | J         | ug/l  | 0.05 | 0.03 | 1               |
| Chrysene                                             | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(b)fluoranthene                                 | ND     |           | ug/l  | 0.05 | 0.03 | 1               |
| Benzo(k)fluoranthene                                 | ND     |           | ug/l  | 0.10 | 0.03 | 1               |
| Benzo(a)pyrene                                       | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Indeno(1,2,3-cd)pyrene                               | ND     |           | ug/l  | 0.10 | 0.02 | 1               |
| Dibenzo(a,h)anthracene                               | ND     |           | ug/l  | 0.05 | 0.02 | 1               |
| Benzo(ghi)perylene                                   | ND     |           | ug/l  | 0.10 | 0.02 | 1               |

| Surrogate            | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol       | 20         | Q         | 21-120              |
| Phenol-d6            | 21         |           | 10-120              |
| Nitrobenzene-d5      | 81         |           | 23-120              |
| 2-Fluorobiphenyl     | 67         |           | 15-120              |
| 2,4,6-Tribromophenol | 43         |           | 10-120              |
| 4-Terphenyl-d14      | 69         |           | 41-149              |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270E  
 Analytical Date: 04/23/25 00:54  
 Analyst: JG

Extraction Method: EPA 3510C  
 Extraction Date: 04/22/25 02:15

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2056815-1 |        |           |       |     |      |
| Benzaldehyde                                                                             | ND     |           | ug/l  | 5.0 | 1.1  |
| Phenol                                                                                   | ND     |           | ug/l  | 5.0 | 0.35 |
| Bis(2-chloroethyl)ether                                                                  | ND     |           | ug/l  | 2.0 | 0.39 |
| 2-Chlorophenol                                                                           | ND     |           | ug/l  | 2.0 | 0.65 |
| 2-Methylphenol                                                                           | ND     |           | ug/l  | 5.0 | 2.3  |
| Bis(2-chloroisopropyl)ether                                                              | ND     |           | ug/l  | 2.0 | 0.40 |
| Acetophenone                                                                             | ND     |           | ug/l  | 5.0 | 0.92 |
| n-Nitrosodi-n-propylamine                                                                | ND     |           | ug/l  | 5.0 | 0.91 |
| 3-Methylphenol/4-Methylphenol                                                            | ND     |           | ug/l  | 5.0 | 1.4  |
| Hexachloroethane                                                                         | ND     |           | ug/l  | 2.0 | 0.20 |
| Nitrobenzene                                                                             | ND     |           | ug/l  | 1.4 | 0.20 |
| Isophorone                                                                               | ND     |           | ug/l  | 5.0 | 0.86 |
| 2,4-Dimethylphenol                                                                       | ND     |           | ug/l  | 5.0 | 2.0  |
| Bis(2-chloroethoxy)methane                                                               | ND     |           | ug/l  | 5.0 | 0.84 |
| 2,4-Dichlorophenol                                                                       | ND     |           | ug/l  | 5.0 | 1.7  |
| Naphthalene                                                                              | ND     |           | ug/l  | 2.0 | 0.54 |
| 4-Chloroaniline                                                                          | ND     |           | ug/l  | 3.7 | 0.47 |
| Hexachlorobutadiene                                                                      | ND     |           | ug/l  | 2.0 | 0.36 |
| Caprolactam                                                                              | ND     |           | ug/l  | 10  | 1.2  |
| 2-Methylnaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.37 |
| Hexachlorocyclopentadiene                                                                | ND     |           | ug/l  | 20  | 1.2  |
| 1,2,4,5-Tetrachlorobenzene                                                               | ND     |           | ug/l  | 1.7 | 0.24 |
| 2,4,6-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| 2,4,5-Trichlorophenol                                                                    | ND     |           | ug/l  | 5.0 | 2.1  |
| Biphenyl                                                                                 | ND     |           | ug/l  | 2.0 | 0.20 |
| 2-Chloronaphthalene                                                                      | ND     |           | ug/l  | 2.0 | 0.35 |
| 2-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.0  |
| 2,6-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.84 |
| Acenaphthylene                                                                           | ND     |           | ug/l  | 2.0 | 0.32 |



**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270E  
 Analytical Date: 04/23/25 00:54  
 Analyst: JG

Extraction Method: EPA 3510C  
 Extraction Date: 04/22/25 02:15

| Parameter                                                                                | Result | Qualifier | Units | RL  | MDL  |
|------------------------------------------------------------------------------------------|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2056815-1 |        |           |       |     |      |
| Acenaphthene                                                                             | ND     |           | ug/l  | 2.0 | 0.40 |
| 2,4-Dinitrophenol                                                                        | ND     |           | ug/l  | 20  | 5.4  |
| 2,4-Dinitrotoluene                                                                       | ND     |           | ug/l  | 5.0 | 0.54 |
| 2,3,4,6-Tetrachlorophenol                                                                | ND     |           | ug/l  | 5.0 | 2.2  |
| Diethyl phthalate                                                                        | ND     |           | ug/l  | 5.0 | 0.76 |
| Fluorene                                                                                 | ND     |           | ug/l  | 2.0 | 0.44 |
| 4-Nitroaniline                                                                           | ND     |           | ug/l  | 5.0 | 1.4  |
| NDPA/DPA                                                                                 | ND     |           | ug/l  | 2.0 | 0.92 |
| Hexachlorobenzene                                                                        | ND     |           | ug/l  | 2.0 | 0.45 |
| Pentachlorophenol                                                                        | ND     |           | ug/l  | 10  | 2.5  |
| Phenanthrene                                                                             | ND     |           | ug/l  | 2.0 | 0.42 |
| Anthracene                                                                               | ND     |           | ug/l  | 2.0 | 0.47 |
| Carbazole                                                                                | ND     |           | ug/l  | 2.0 | 0.31 |
| Di-n-butylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 0.96 |
| Fluoranthene                                                                             | ND     |           | ug/l  | 2.0 | 0.41 |
| Pyrene                                                                                   | ND     |           | ug/l  | 2.0 | 0.41 |
| 3,3'-Dichlorobenzidine                                                                   | ND     |           | ug/l  | 5.0 | 1.8  |
| Benzo(a)anthracene                                                                       | ND     |           | ug/l  | 2.0 | 0.32 |
| Chrysene                                                                                 | ND     |           | ug/l  | 1.5 | 0.22 |
| Bis(2-ethylhexyl)phthalate                                                               | ND     |           | ug/l  | 3.0 | 1.4  |
| Di-n-octylphthalate                                                                      | ND     |           | ug/l  | 5.0 | 2.3  |
| Benzo(b)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.53 |
| Benzo(k)fluoranthene                                                                     | ND     |           | ug/l  | 2.0 | 0.62 |
| Benzo(a)pyrene                                                                           | ND     |           | ug/l  | 2.0 | 0.37 |
| Indeno(1,2,3-cd)pyrene                                                                   | ND     |           | ug/l  | 2.0 | 0.48 |
| Dibenzo(a,h)anthracene                                                                   | ND     |           | ug/l  | 2.0 | 0.29 |
| Benzo(ghi)perylene                                                                       | ND     |           | ug/l  | 2.0 | 0.37 |



**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8270E  
Analytical Date: 04/23/25 00:54  
Analyst: JG

Extraction Method: EPA 3510C  
Extraction Date: 04/22/25 02:15

| Parameter                                                                                | Result | Qualifier | Units | RL | MDL |
|------------------------------------------------------------------------------------------|--------|-----------|-------|----|-----|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01-04 Batch: WG2056815-1 |        |           |       |    |     |

| Surrogate            | %Recovery | Qualifier | Acceptance<br>Criteria |
|----------------------|-----------|-----------|------------------------|
| 2-Fluorophenol       | 55        |           | 21-120                 |
| Phenol-d6            | 40        |           | 10-120                 |
| Nitrobenzene-d5      | 78        |           | 23-120                 |
| 2-Fluorobiphenyl     | 85        |           | 15-120                 |
| 2,4,6-Tribromophenol | 79        |           | 10-120                 |
| 4-Terphenyl-d14      | 90        |           | 41-149                 |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

### Method Blank Analysis Batch Quality Control

**Analytical Method:** 1,8270E-SIM  
**Analytical Date:** 04/22/25 10:49  
**Analyst:** SLR

**Extraction Method:** EPA 3510C  
**Extraction Date:** 04/22/25 02:15

| Parameter                                                                                    | Result | Qualifier | Units | RL   | MDL  |
|----------------------------------------------------------------------------------------------|--------|-----------|-------|------|------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01-04 Batch: WG2056816-1 |        |           |       |      |      |
| Naphthalene                                                                                  | ND     |           | ug/l  | 0.10 | 0.02 |
| 2-Methylnaphthalene                                                                          | ND     |           | ug/l  | 0.10 | 0.03 |
| Acenaphthylene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Acenaphthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluorene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Phenanthrene                                                                                 | ND     |           | ug/l  | 0.05 | 0.04 |
| Anthracene                                                                                   | ND     |           | ug/l  | 0.10 | 0.02 |
| Fluoranthene                                                                                 | ND     |           | ug/l  | 0.10 | 0.03 |
| Pyrene                                                                                       | ND     |           | ug/l  | 0.10 | 0.04 |
| Benzo(a)anthracene                                                                           | ND     |           | ug/l  | 0.05 | 0.03 |
| Chrysene                                                                                     | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(b)fluoranthene                                                                         | ND     |           | ug/l  | 0.05 | 0.03 |
| Benzo(k)fluoranthene                                                                         | ND     |           | ug/l  | 0.10 | 0.03 |
| Benzo(a)pyrene                                                                               | ND     |           | ug/l  | 0.10 | 0.02 |
| Indeno(1,2,3-cd)pyrene                                                                       | ND     |           | ug/l  | 0.10 | 0.02 |
| Dibenzo(a,h)anthracene                                                                       | ND     |           | ug/l  | 0.05 | 0.02 |
| Benzo(ghi)perylene                                                                           | ND     |           | ug/l  | 0.10 | 0.02 |

| Surrogate            | %Recovery | Qualifier | Acceptance Criteria |
|----------------------|-----------|-----------|---------------------|
| 2-Fluorophenol       | 31        |           | 21-120              |
| Phenol-d6            | 23        |           | 10-120              |
| Nitrobenzene-d5      | 51        |           | 23-120              |
| 2-Fluorobiphenyl     | 43        |           | 15-120              |
| 2,4,6-Tribromophenol | 49        |           | 10-120              |
| 4-Terphenyl-d14      | 43        |           | 41-149              |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Project Number:** Not Specified

**Lab Number:** L2523470

**Report Date:** 04/29/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2056815-2 WG2056815-3 |                  |      |                   |      |                     |     |      |               |
| Benzaldehyde                                                                                                | 85               |      | 96                |      | 40-140              | 12  |      | 30            |
| Phenol                                                                                                      | 38               |      | 45                |      | 12-110              | 17  |      | 30            |
| Bis(2-chloroethyl)ether                                                                                     | 80               |      | 94                |      | 40-140              | 16  |      | 30            |
| 2-Chlorophenol                                                                                              | 73               |      | 90                |      | 27-123              | 21  |      | 30            |
| 2-Methylphenol                                                                                              | 73               |      | 83                |      | 30-130              | 13  |      | 30            |
| Bis(2-chloroisopropyl)ether                                                                                 | 80               |      | 91                |      | 40-140              | 13  |      | 30            |
| Acetophenone                                                                                                | 82               |      | 97                |      | 39-129              | 17  |      | 30            |
| n-Nitrosodi-n-propylamine                                                                                   | 79               |      | 87                |      | 29-132              | 10  |      | 30            |
| 3-Methylphenol/4-Methylphenol                                                                               | 78               |      | 86                |      | 30-130              | 10  |      | 30            |
| Hexachloroethane                                                                                            | 73               |      | 82                |      | 40-140              | 12  |      | 30            |
| Nitrobenzene                                                                                                | 79               |      | 90                |      | 40-140              | 13  |      | 30            |
| Isophorone                                                                                                  | 77               |      | 86                |      | 40-140              | 11  |      | 30            |
| 2,4-Dimethylphenol                                                                                          | 62               |      | 71                |      | 30-130              | 14  |      | 30            |
| Bis(2-chloroethoxy)methane                                                                                  | 78               |      | 91                |      | 40-140              | 15  |      | 30            |
| 2,4-Dichlorophenol                                                                                          | 84               |      | 98                |      | 30-130              | 15  |      | 30            |
| Naphthalene                                                                                                 | 80               |      | 87                |      | 40-140              | 8   |      | 30            |
| 4-Chloroaniline                                                                                             | 74               |      | 77                |      | 40-140              | 4   |      | 30            |
| Hexachlorobutadiene                                                                                         | 79               |      | 86                |      | 40-140              | 8   |      | 30            |
| Caprolactam                                                                                                 | 41               |      | 44                |      | 10-130              | 7   |      | 30            |
| 2-Methylnaphthalene                                                                                         | 84               |      | 92                |      | 40-140              | 9   |      | 30            |
| Hexachlorocyclopentadiene                                                                                   | 74               |      | 81                |      | 40-140              | 9   |      | 30            |
| 1,2,4,5-Tetrachlorobenzene                                                                                  | 86               |      | 94                |      | 2-134               | 9   |      | 30            |
| 2,4,6-Trichlorophenol                                                                                       | 78               |      | 94                |      | 30-130              | 19  |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Project Number:** Not Specified

**Lab Number:** L2523470

**Report Date:** 04/29/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2056815-2 WG2056815-3 |                  |      |                   |      |                     |     |      |               |
| 2,4,5-Trichlorophenol                                                                                       | 93               |      | 99                |      | 30-130              | 6   |      | 30            |
| Biphenyl                                                                                                    | 90               |      | 98                |      | 40-140              | 9   |      | 30            |
| 2-Chloronaphthalene                                                                                         | 84               |      | 94                |      | 40-140              | 11  |      | 30            |
| 2-Nitroaniline                                                                                              | 90               |      | 102               |      | 52-143              | 13  |      | 30            |
| 2,6-Dinitrotoluene                                                                                          | 91               |      | 96                |      | 40-140              | 5   |      | 30            |
| Acenaphthylene                                                                                              | 84               |      | 93                |      | 45-123              | 10  |      | 30            |
| Acenaphthene                                                                                                | 82               |      | 90                |      | 37-111              | 9   |      | 30            |
| 2,4-Dinitrophenol                                                                                           | 44               |      | 62                |      | 20-130              | 34  | Q    | 30            |
| 2,4-Dinitrotoluene                                                                                          | 86               |      | 92                |      | 48-143              | 7   |      | 30            |
| 2,3,4,6-Tetrachlorophenol                                                                                   | 84               |      | 94                |      | 54-145              | 11  |      | 30            |
| Diethyl phthalate                                                                                           | 84               |      | 87                |      | 40-140              | 4   |      | 30            |
| Fluorene                                                                                                    | 86               |      | 92                |      | 40-140              | 7   |      | 30            |
| 4-Nitroaniline                                                                                              | 85               |      | 91                |      | 51-143              | 7   |      | 30            |
| NDPA/DPA                                                                                                    | 89               |      | 93                |      | 40-140              | 4   |      | 30            |
| Hexachlorobenzene                                                                                           | 87               |      | 92                |      | 40-140              | 6   |      | 30            |
| Pentachlorophenol                                                                                           | 77               |      | 90                |      | 9-103               | 16  |      | 30            |
| Phenanthrene                                                                                                | 87               |      | 91                |      | 40-140              | 4   |      | 30            |
| Anthracene                                                                                                  | 90               |      | 95                |      | 40-140              | 5   |      | 30            |
| Carbazole                                                                                                   | 90               |      | 96                |      | 55-144              | 6   |      | 30            |
| Di-n-butylphthalate                                                                                         | 92               |      | 96                |      | 40-140              | 4   |      | 30            |
| Fluoranthene                                                                                                | 91               |      | 96                |      | 40-140              | 5   |      | 30            |
| Pyrene                                                                                                      | 92               |      | 97                |      | 26-127              | 5   |      | 30            |
| 3,3'-Dichlorobenzidine                                                                                      | 85               |      | 93                |      | 40-140              | 9   |      | 30            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Project Number:** Not Specified

**Lab Number:** L2523470

**Report Date:** 04/29/25

| Parameter                                                                                                   | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-04 Batch: WG2056815-2 WG2056815-3 |                  |      |                   |      |                     |     |      |               |
| Benzo(a)anthracene                                                                                          | 90               |      | 93                |      | 40-140              | 3   |      | 30            |
| Chrysene                                                                                                    | 90               |      | 94                |      | 40-140              | 4   |      | 30            |
| Bis(2-ethylhexyl)phthalate                                                                                  | 96               |      | 98                |      | 40-140              | 2   |      | 30            |
| Di-n-octylphthalate                                                                                         | 97               |      | 101               |      | 40-140              | 4   |      | 30            |
| Benzo(b)fluoranthene                                                                                        | 94               |      | 97                |      | 40-140              | 3   |      | 30            |
| Benzo(k)fluoranthene                                                                                        | 93               |      | 98                |      | 40-140              | 5   |      | 30            |
| Benzo(a)pyrene                                                                                              | 94               |      | 99                |      | 40-140              | 5   |      | 30            |
| Indeno(1,2,3-cd)pyrene                                                                                      | 92               |      | 94                |      | 40-140              | 2   |      | 30            |
| Dibenzo(a,h)anthracene                                                                                      | 89               |      | 94                |      | 40-140              | 5   |      | 30            |
| Benzo(ghi)perylene                                                                                          | 87               |      | 91                |      | 40-140              | 4   |      | 30            |

| Surrogate            | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | Acceptance<br>Criteria |
|----------------------|------------------|------|-------------------|------|------------------------|
| 2-Fluorophenol       | 53               |      | 66                |      | 21-120                 |
| Phenol-d6            | 41               |      | 49                |      | 10-120                 |
| Nitrobenzene-d5      | 81               |      | 91                |      | 23-120                 |
| 2-Fluorobiphenyl     | 85               |      | 94                |      | 15-120                 |
| 2,4,6-Tribromophenol | 86               |      | 97                |      | 10-120                 |
| 4-Terphenyl-d14      | 91               |      | 98                |      | 41-149                 |



# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Project Number:** Not Specified

**Lab Number:** L2523470

**Report Date:** 04/29/25

| Parameter                                                                                                       | LCS<br>%Recovery | Qual | LCSD<br>%Recovery | Qual | %Recovery<br>Limits | RPD | Qual | RPD<br>Limits |
|-----------------------------------------------------------------------------------------------------------------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2056816-2 WG2056816-3 |                  |      |                   |      |                     |     |      |               |
| Naphthalene                                                                                                     | 53               |      | 58                |      | 40-140              | 9   |      | 40            |
| 2-Methylnaphthalene                                                                                             | 55               |      | 60                |      | 40-140              | 9   |      | 40            |
| Acenaphthylene                                                                                                  | 64               |      | 70                |      | 40-140              | 9   |      | 40            |
| Acenaphthene                                                                                                    | 53               |      | 58                |      | 37-111              | 9   |      | 40            |
| Fluorene                                                                                                        | 58               |      | 62                |      | 40-140              | 7   |      | 40            |
| Phenanthrene                                                                                                    | 53               |      | 55                |      | 40-140              | 4   |      | 40            |
| Anthracene                                                                                                      | 56               |      | 59                |      | 40-140              | 5   |      | 40            |
| Fluoranthene                                                                                                    | 56               |      | 59                |      | 40-140              | 5   |      | 40            |
| Pyrene                                                                                                          | 54               |      | 57                |      | 26-127              | 5   |      | 40            |
| Benzo(a)anthracene                                                                                              | 53               |      | 56                |      | 40-140              | 6   |      | 40            |
| Chrysene                                                                                                        | 54               |      | 56                |      | 40-140              | 4   |      | 40            |
| Benzo(b)fluoranthene                                                                                            | 56               |      | 58                |      | 40-140              | 4   |      | 40            |
| Benzo(k)fluoranthene                                                                                            | 57               |      | 61                |      | 40-140              | 7   |      | 40            |
| Benzo(a)pyrene                                                                                                  | 60               |      | 64                |      | 40-140              | 6   |      | 40            |
| Indeno(1,2,3-cd)pyrene                                                                                          | 66               |      | 70                |      | 40-140              | 6   |      | 40            |
| Dibenzo(a,h)anthracene                                                                                          | 64               |      | 68                |      | 40-140              | 6   |      | 40            |
| Benzo(ghi)perylene                                                                                              | 62               |      | 65                |      | 40-140              | 5   |      | 40            |

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** TMC SURFACE WATER SAMPLING

**Lab Number:** L2523470

**Project Number:** Not Specified

**Report Date:** 04/29/25

| <b>Parameter</b>                                                                                                | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>%Recovery<br/>Limits</b> | <b>RPD</b> | <b>Qual</b> | <b>RPD<br/>Limits</b> |
|-----------------------------------------------------------------------------------------------------------------|--------------------------|-------------|---------------------------|-------------|-----------------------------|------------|-------------|-----------------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01-04 Batch: WG2056816-2 WG2056816-3 |                          |             |                           |             |                             |            |             |                       |

| <b>Surrogate</b>     | <b>LCS<br/>%Recovery</b> | <b>Qual</b> | <b>LCSD<br/>%Recovery</b> | <b>Qual</b> | <b>Acceptance<br/>Criteria</b> |
|----------------------|--------------------------|-------------|---------------------------|-------------|--------------------------------|
| 2-Fluorophenol       | 42                       |             | 48                        |             | 21-120                         |
| Phenol-d6            | 33                       |             | 38                        |             | 10-120                         |
| Nitrobenzene-d5      | 75                       |             | 81                        |             | 23-120                         |
| 2-Fluorobiphenyl     | 58                       |             | 64                        |             | 15-120                         |
| 2,4,6-Tribromophenol | 69                       |             | 78                        |             | 10-120                         |
| 4-Terphenyl-d14      | 50                       |             | 53                        |             | 41-149                         |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

**Cooler Information**

|               |                     |
|---------------|---------------------|
| <b>Cooler</b> | <b>Custody Seal</b> |
| A             | Absent              |

**Container Information**

| <b>Container ID</b> | <b>Container Type</b>   | <b>Cooler</b> | <b>Initial pH</b> | <b>Final pH</b> | <b>Temp deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen Date/Time</b> | <b>Analysis(*)</b>               |
|---------------------|-------------------------|---------------|-------------------|-----------------|-------------------|-------------|-------------|-------------------------|----------------------------------|
| L2523470-01A        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-01B        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-01C        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-01D        | Amber 100ml unpreserved | A             | 8                 | 8               | 3.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523470-01E        | Amber 100ml unpreserved | A             | 8                 | 8               | 3.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523470-02A        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-02B        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-02C        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-02D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523470-02E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523470-03A        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-03B        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-03C        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-03D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523470-03E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.1               | Y           | Absent      |                         | PA-8270-RVT(7),PA-8270SIM-RVT(7) |
| L2523470-04A        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-04B        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-04C        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-04D        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.1               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2523470-04E        | Amber 100ml unpreserved | A             | 7                 | 7               | 3.1               | Y           | Absent      |                         | PA-8270SIM-RVT(7),PA-8270-RVT(7) |
| L2523470-05A        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-05B        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |
| L2523470-05C        | Vial HCl preserved      | A             | NA                |                 | 3.1               | Y           | Absent      |                         | PA-8260-SIM(14),PA-8260(14)      |

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

Serial\_No:04292518:27  
**Lab Number:** L2523470  
**Report Date:** 04/29/25

**Container Information**

| <b>Container ID</b> | <b>Container Type</b> | <b>Cooler</b> | <b>Initial<br/>pH</b> | <b>Final<br/>pH</b> | <b>Temp<br/>deg C</b> | <b>Pres</b> | <b>Seal</b> | <b>Frozen<br/>Date/Time</b> | <b>Analysis(*)</b>          |
|---------------------|-----------------------|---------------|-----------------------|---------------------|-----------------------|-------------|-------------|-----------------------------|-----------------------------|
| L2523470-05D        | Vial HCl preserved    | A             | NA                    |                     | 3.1                   | Y           | Absent      |                             | PA-8260-SIM(14),PA-8260(14) |

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25

## GLOSSARY

### Acronyms

|          |                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            |
|----------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| DL       | - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                |
| EDL      | - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME).                                                                                         |
| EMPC     | - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration.                                                                                                                                                                                                                               |
| EPA      | - Environmental Protection Agency.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| LCS      | - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                          |
| LCSD     | - Laboratory Control Sample Duplicate: Refer to LCS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                       |
| LFB      | - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes.                                                                                                                                                                                                                                                                                                                         |
| LOD      | - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)                                                                                                                                                                                                               |
| LOQ      | - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)<br><br>Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| MDL      | - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                          |
| MS       | - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.                                                                                                                                                                                                    |
| MSD      | - Matrix Spike Sample Duplicate: Refer to MS.                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NA       | - Not Applicable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                          |
| NC       | - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.                                                                                                                                                                                                                                                                                                                                                                           |
| NDPA/DPA | - N-Nitrosodiphenylamine/Diphenylamine.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| NI       | - Not Ignitable.                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           |
| NP       | - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| NR       | - No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.                                                                                                                                                                                                                                                                                                                                                                    |
| RL       | - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.                                                                                                                                                                                                                                                                                                   |
| RPD      | - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.                                                                  |
| SRM      | - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.                                                                                                                                                                                                                                                                                                                                                                     |
| STLP     | - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.                                                                                                                                                                                                                                                                                                                                                                                                                                                                |
| TEF      | - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.                                                                                                                                                                                                                                                                                                                                                                                             |
| TEQ      | - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.                                                                                                                                                                                                                                                                                                                                                        |
| TIC      | - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.                                                                                                                                                                                                                                                                      |

*Report Format: DU Report with 'J' Qualifiers*

**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

### Footnotes

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

### Terms

**Analytical Method:** Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

**Chlordane:** The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

**Difference:** With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

**Final pH:** As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

**Frozen Date/Time:** With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**Initial pH:** As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

**PAH Total:** With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

**PFAS Total:** With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. In addition, the 'PFAS, Total (6)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**Total:** With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

### Data Qualifiers

- A** - Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- F** - The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The lower value for the two columns has been reported due to obvious interference.
- J** - Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively

**Report Format:** DU Report with 'J' Qualifiers



**Project Name:** TMC SURFACE WATER SAMPLING  
**Project Number:** Not Specified

**Lab Number:** L2523470  
**Report Date:** 04/29/25

#### Data Qualifiers

Identified Compounds (TICs). For calculated parameters, this represents that one or more values used in the calculation were estimated.

- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND** - Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ** - Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- R** - Analytical results are from sample re-analysis.
- RE** - Analytical results are from sample re-extraction.
- S** - Analytical results are from modified screening analysis.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

**Project Name:** TMC SURFACE WATER SAMPLING**Lab Number:** L2523470**Project Number:** Not Specified**Report Date:** 04/29/25

## REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

## LIMITATION OF LIABILITIES

Pace Analytical Services performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Pace Analytical Services shall be to re-perform the work at it's own expense. In no event shall Pace Analytical Services be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Pace Analytical Services.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.





**Pace Analytical Services LLC**Facility: **Northeast**Department: **Quality Assurance**Title: **Certificate/Approval Program Summary**ID No.: **17873**

Revision 27

Published Date: 01/24/2025

Page 1 of 2

**Certification Information**

The following analytes are not included in our Primary NELAP Scope of Accreditation:

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****EPA 624.1:** m/p-xylene, o-xylene, Naphthalene**EPA 625.1:** alpha-Terpineol**EPA 8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****SM 2540D:** TSS.**EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**MADEP-APH.****Nonpotable Water:** EPA RSK-175 Dissolved Gases**Biological Tissue Matrix:** EPA 3050B**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048****EPA TO-15:** Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

**Nonpotable Water:** EPA RSK-175 Dissolved Gases

The following test method is not included in our New Jersey Secondary NELAP Scope of Accreditation:

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Determination of Selected Perfluorinated Alkyl Substances by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry Isotope Dilution (via Alpha SOP 23528)**

The following analytes are included in our Massachusetts DEP Scope of Accreditation

**Westborough Facility – 8 Walkup Dr. Westborough, MA 01581****Drinking Water****EPA 300.0:** Chloride, Nitrate-N, Fluoride, Sulfate; **EPA 353.2:** Nitrate-N, Nitrite-N; **SM4500NO3-F:** Nitrate-N, Nitrite-N; **SM4500F-C, SM4500CN-CE, EPA 180.1, SM2130B, SM4500Cl-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B****EPA 524.2:** THMs and VOCs; **EPA 504.1:** EDB, DBCP.**Microbiology:** SM9215B; SM9223-P/A, SM9223B-Colilert-QT, SM9222D.**Non-Potable Water****SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH:** Ammonia-N and Kjeldahl-N, **EPA 350.1:** Ammonia-N, **LACHAT 10-107-06-1-B:** Ammonia-N, **EPA 351.1, SM4500NO3-F, EPA 353.2:** Nitrate-N, **SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300:** Chloride, Sulfate, Nitrate.**EPA 624.1:** Volatile Halocarbons & Aromatics,**EPA 608.3:** Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs**EPA 625.1:** SVOC (Acid/Base/Neutral Extractables).**Microbiology:** SM9223B-Colilert-QT; Enterolert-QT, EPA 1600, EPA 1603, SM9222D.**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048****Drinking Water****EPA 200.7:** Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1 Hg. EPA 522, EPA 537.1.****Non-Potable Water****EPA 200.7:** Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.**EPA 200.8:** Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.**EPA 245.1 Hg.****SM2340B**

**Pace Analytical Services LLC**ID No.: **17873**Facility: **Northeast**

Revision 27

Department: **Quality Assurance**

Published Date: 01/24/2025

Title: **Certificate/Approval Program Summary**

Page 2 of 2

**Certification IDs:****Westborough Facility – 8 Walkup Dr. Westborough, MA 01581**

CT PH-0826, IL 200077, IN C-MA-03, KY JY98045, ME MA00086, MD 348, MA M-MA086, NH 2064, NJ MA935, NY 11148, NC (DW) 25700, NC (NPW/SCM) 666, OR MA-1316, PA 68-03671, RI LAO00065, TX T104704476, VT VT-0935, VA 460195

**Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048**

CT PH-0825, ANAB/DoD L2474, IL 200081, IN C-MA-04, KY KY98046, LA 3090, ME MA00030, MI 9110, MN 025-999-495, NH 2062, NJ MA015, NY 11627, NC (NPW/SCM) 685, OR MA-0262, PA 68-02089, RI LAO00299, TX T-104704419, VT VT-0015, VA 460194, WA C954

**Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048**

ANAB/DoD L2474, ME MA01156, MN 025-999-498, NH 2249, NJ MA025, NY 12191, OR 4203, TX T104704583, VA 460311, WA C1104.

---

For a complete listing of analytes and methods, please contact your Project Manager.



---

---

## **APPENDIX B**

---

---

## Mann-Kendall Trend Analysis

Parameter: Naphthalene

Location: TM04-PZM006

Original Data (Not Transformed)

Non-Detects Replaced with 1/2 DL

95% Confidence Level

| Xj        | Xk  | Xj - Xk | Positives | Negatives |
|-----------|-----|---------|-----------|-----------|
| 51        | 200 | -149    | 0         | 1         |
| 405       | 200 | 205     | 1         | 1         |
| 358       | 200 | 158     | 2         | 1         |
| 432 L1    | 200 | 232     | 3         | 1         |
| 140       | 200 | -60     | 3         | 2         |
| 69        | 200 | -131    | 3         | 3         |
| 240       | 200 | 40      | 4         | 3         |
| 54        | 200 | -146    | 4         | 4         |
| 230       | 200 | 30      | 5         | 4         |
| 160       | 200 | -40     | 5         | 5         |
| 180       | 200 | -20     | 5         | 6         |
| 140       | 200 | -60     | 5         | 7         |
| 210       | 200 | 10      | 6         | 7         |
| 160       | 200 | -40     | 6         | 8         |
| 210       | 200 | 10      | 7         | 8         |
| 68        | 200 | -132    | 7         | 9         |
| ND<0.05 U | 200 | -199.95 | 7         | 10        |
| 170       | 200 | -30     | 7         | 11        |
| 0.12      | 200 | -199.88 | 7         | 12        |
| 13        | 200 | -187    | 7         | 13        |
| 130       | 200 | -70     | 7         | 14        |
| 0.12      | 200 | -199.88 | 7         | 15        |
| 405       | 51  | 354     | 8         | 15        |
| 358       | 51  | 307     | 9         | 15        |
| 432 L1    | 51  | 381     | 10        | 15        |
| 140       | 51  | 89      | 11        | 15        |
| 69        | 51  | 18      | 12        | 15        |
| 240       | 51  | 189     | 13        | 15        |
| 54        | 51  | 3       | 14        | 15        |
| 230       | 51  | 179     | 15        | 15        |
| 160       | 51  | 109     | 16        | 15        |
| 180       | 51  | 129     | 17        | 15        |
| 140       | 51  | 89      | 18        | 15        |
| 210       | 51  | 159     | 19        | 15        |
| 160       | 51  | 109     | 20        | 15        |
| 210       | 51  | 159     | 21        | 15        |
| 68        | 51  | 17      | 22        | 15        |
| ND<0.05 U | 51  | -50.95  | 22        | 16        |
| 170       | 51  | 119     | 23        | 16        |
| 0.12      | 51  | -50.88  | 23        | 17        |
| 13        | 51  | -38     | 23        | 18        |
| 130       | 51  | 79      | 24        | 18        |
| 0.12      | 51  | -50.88  | 24        | 19        |
| 358       | 405 | -47     | 24        | 20        |
| 432 L1    | 405 | 27      | 25        | 20        |

|           |        |         |    |    |
|-----------|--------|---------|----|----|
| 140       | 405    | -265    | 25 | 21 |
| 69        | 405    | -336    | 25 | 22 |
| 240       | 405    | -165    | 25 | 23 |
| 54        | 405    | -351    | 25 | 24 |
| 230       | 405    | -175    | 25 | 25 |
| 160       | 405    | -245    | 25 | 26 |
| 180       | 405    | -225    | 25 | 27 |
| 140       | 405    | -265    | 25 | 28 |
| 210       | 405    | -195    | 25 | 29 |
| 160       | 405    | -245    | 25 | 30 |
| 210       | 405    | -195    | 25 | 31 |
| 68        | 405    | -337    | 25 | 32 |
| ND<0.05 U | 405    | -404.95 | 25 | 33 |
| 170       | 405    | -235    | 25 | 34 |
| 0.12      | 405    | -404.88 | 25 | 35 |
| 13        | 405    | -392    | 25 | 36 |
| 130       | 405    | -275    | 25 | 37 |
| 0.12      | 405    | -404.88 | 25 | 38 |
|           |        |         |    |    |
| 432 L1    | 358    | 74      | 26 | 38 |
| 140       | 358    | -218    | 26 | 39 |
| 69        | 358    | -289    | 26 | 40 |
| 240       | 358    | -118    | 26 | 41 |
| 54        | 358    | -304    | 26 | 42 |
| 230       | 358    | -128    | 26 | 43 |
| 160       | 358    | -198    | 26 | 44 |
| 180       | 358    | -178    | 26 | 45 |
| 140       | 358    | -218    | 26 | 46 |
| 210       | 358    | -148    | 26 | 47 |
| 160       | 358    | -198    | 26 | 48 |
| 210       | 358    | -148    | 26 | 49 |
| 68        | 358    | -290    | 26 | 50 |
| ND<0.05 U | 358    | -357.95 | 26 | 51 |
| 170       | 358    | -188    | 26 | 52 |
| 0.12      | 358    | -357.88 | 26 | 53 |
| 13        | 358    | -345    | 26 | 54 |
| 130       | 358    | -228    | 26 | 55 |
| 0.12      | 358    | -357.88 | 26 | 56 |
|           |        |         |    |    |
| 140       | 432 L1 | -292    | 26 | 57 |
| 69        | 432 L1 | -363    | 26 | 58 |
| 240       | 432 L1 | -192    | 26 | 59 |
| 54        | 432 L1 | -378    | 26 | 60 |
| 230       | 432 L1 | -202    | 26 | 61 |
| 160       | 432 L1 | -272    | 26 | 62 |
| 180       | 432 L1 | -252    | 26 | 63 |
| 140       | 432 L1 | -292    | 26 | 64 |
| 210       | 432 L1 | -222    | 26 | 65 |
| 160       | 432 L1 | -272    | 26 | 66 |
| 210       | 432 L1 | -222    | 26 | 67 |
| 68        | 432 L1 | -364    | 26 | 68 |
| ND<0.05 U | 432 L1 | -431.95 | 26 | 69 |
| 170       | 432 L1 | -262    | 26 | 70 |
| 0.12      | 432 L1 | -431.88 | 26 | 71 |
| 13        | 432 L1 | -419    | 26 | 72 |
| 130       | 432 L1 | -302    | 26 | 73 |
| 0.12      | 432 L1 | -431.88 | 26 | 74 |

|           |     |         |    |     |
|-----------|-----|---------|----|-----|
| 69        | 140 | -71     | 26 | 75  |
| 240       | 140 | 100     | 27 | 75  |
| 54        | 140 | -86     | 27 | 76  |
| 230       | 140 | 90      | 28 | 76  |
| 160       | 140 | 20      | 29 | 76  |
| 180       | 140 | 40      | 30 | 76  |
| 140       | 140 | 0       | 30 | 76  |
| 210       | 140 | 70      | 31 | 76  |
| 160       | 140 | 20      | 32 | 76  |
| 210       | 140 | 70      | 33 | 76  |
| 68        | 140 | -72     | 33 | 77  |
| ND<0.05 U | 140 | -139.95 | 33 | 78  |
| 170       | 140 | 30      | 34 | 78  |
| 0.12      | 140 | -139.88 | 34 | 79  |
| 13        | 140 | -127    | 34 | 80  |
| 130       | 140 | -10     | 34 | 81  |
| 0.12      | 140 | -139.88 | 34 | 82  |
|           |     |         |    |     |
| 240       | 69  | 171     | 35 | 82  |
| 54        | 69  | -15     | 35 | 83  |
| 230       | 69  | 161     | 36 | 83  |
| 160       | 69  | 91      | 37 | 83  |
| 180       | 69  | 111     | 38 | 83  |
| 140       | 69  | 71      | 39 | 83  |
| 210       | 69  | 141     | 40 | 83  |
| 160       | 69  | 91      | 41 | 83  |
| 210       | 69  | 141     | 42 | 83  |
| 68        | 69  | -1      | 42 | 84  |
| ND<0.05 U | 69  | -68.95  | 42 | 85  |
| 170       | 69  | 101     | 43 | 85  |
| 0.12      | 69  | -68.88  | 43 | 86  |
| 13        | 69  | -56     | 43 | 87  |
| 130       | 69  | 61      | 44 | 87  |
| 0.12      | 69  | -68.88  | 44 | 88  |
|           |     |         |    |     |
| 54        | 240 | -186    | 44 | 89  |
| 230       | 240 | -10     | 44 | 90  |
| 160       | 240 | -80     | 44 | 91  |
| 180       | 240 | -60     | 44 | 92  |
| 140       | 240 | -100    | 44 | 93  |
| 210       | 240 | -30     | 44 | 94  |
| 160       | 240 | -80     | 44 | 95  |
| 210       | 240 | -30     | 44 | 96  |
| 68        | 240 | -172    | 44 | 97  |
| ND<0.05 U | 240 | -239.95 | 44 | 98  |
| 170       | 240 | -70     | 44 | 99  |
| 0.12      | 240 | -239.88 | 44 | 100 |
| 13        | 240 | -227    | 44 | 101 |
| 130       | 240 | -110    | 44 | 102 |
| 0.12      | 240 | -239.88 | 44 | 103 |
|           |     |         |    |     |
| 230       | 54  | 176     | 45 | 103 |
| 160       | 54  | 106     | 46 | 103 |
| 180       | 54  | 126     | 47 | 103 |
| 140       | 54  | 86      | 48 | 103 |
| 210       | 54  | 156     | 49 | 103 |

|           |     |         |    |     |
|-----------|-----|---------|----|-----|
| 160       | 54  | 106     | 50 | 103 |
| 210       | 54  | 156     | 51 | 103 |
| 68        | 54  | 14      | 52 | 103 |
| ND<0.05 U | 54  | -53.95  | 52 | 104 |
| 170       | 54  | 116     | 53 | 104 |
| 0.12      | 54  | -53.88  | 53 | 105 |
| 13        | 54  | -41     | 53 | 106 |
| 130       | 54  | 76      | 54 | 106 |
| 0.12      | 54  | -53.88  | 54 | 107 |
|           |     |         |    |     |
| 160       | 230 | -70     | 54 | 108 |
| 180       | 230 | -50     | 54 | 109 |
| 140       | 230 | -90     | 54 | 110 |
| 210       | 230 | -20     | 54 | 111 |
| 160       | 230 | -70     | 54 | 112 |
| 210       | 230 | -20     | 54 | 113 |
| 68        | 230 | -162    | 54 | 114 |
| ND<0.05 U | 230 | -229.95 | 54 | 115 |
| 170       | 230 | -60     | 54 | 116 |
| 0.12      | 230 | -229.88 | 54 | 117 |
| 13        | 230 | -217    | 54 | 118 |
| 130       | 230 | -100    | 54 | 119 |
| 0.12      | 230 | -229.88 | 54 | 120 |
|           |     |         |    |     |
| 180       | 160 | 20      | 55 | 120 |
| 140       | 160 | -20     | 55 | 121 |
| 210       | 160 | 50      | 56 | 121 |
| 160       | 160 | 0       | 56 | 121 |
| 210       | 160 | 50      | 57 | 121 |
| 68        | 160 | -92     | 57 | 122 |
| ND<0.05 U | 160 | -159.95 | 57 | 123 |
| 170       | 160 | 10      | 58 | 123 |
| 0.12      | 160 | -159.88 | 58 | 124 |
| 13        | 160 | -147    | 58 | 125 |
| 130       | 160 | -30     | 58 | 126 |
| 0.12      | 160 | -159.88 | 58 | 127 |
|           |     |         |    |     |
| 140       | 180 | -40     | 58 | 128 |
| 210       | 180 | 30      | 59 | 128 |
| 160       | 180 | -20     | 59 | 129 |
| 210       | 180 | 30      | 60 | 129 |
| 68        | 180 | -112    | 60 | 130 |
| ND<0.05 U | 180 | -179.95 | 60 | 131 |
| 170       | 180 | -10     | 60 | 132 |
| 0.12      | 180 | -179.88 | 60 | 133 |
| 13        | 180 | -167    | 60 | 134 |
| 130       | 180 | -50     | 60 | 135 |
| 0.12      | 180 | -179.88 | 60 | 136 |
|           |     |         |    |     |
| 210       | 140 | 70      | 61 | 136 |
| 160       | 140 | 20      | 62 | 136 |
| 210       | 140 | 70      | 63 | 136 |
| 68        | 140 | -72     | 63 | 137 |
| ND<0.05 U | 140 | -139.95 | 63 | 138 |
| 170       | 140 | 30      | 64 | 138 |
| 0.12      | 140 | -139.88 | 64 | 139 |
| 13        | 140 | -127    | 64 | 140 |



|           |           |         |    |     |
|-----------|-----------|---------|----|-----|
| 130       | 140       | -10     | 64 | 141 |
| 0.12      | 140       | -139.88 | 64 | 142 |
| 160       | 210       | -50     | 64 | 143 |
| 210       | 210       | 0       | 64 | 143 |
| 68        | 210       | -142    | 64 | 144 |
| ND<0.05 U | 210       | -209.95 | 64 | 145 |
| 170       | 210       | -40     | 64 | 146 |
| 0.12      | 210       | -209.88 | 64 | 147 |
| 13        | 210       | -197    | 64 | 148 |
| 130       | 210       | -80     | 64 | 149 |
| 0.12      | 210       | -209.88 | 64 | 150 |
| 210       | 160       | 50      | 65 | 150 |
| 68        | 160       | -92     | 65 | 151 |
| ND<0.05 U | 160       | -159.95 | 65 | 152 |
| 170       | 160       | 10      | 66 | 152 |
| 0.12      | 160       | -159.88 | 66 | 153 |
| 13        | 160       | -147    | 66 | 154 |
| 130       | 160       | -30     | 66 | 155 |
| 0.12      | 160       | -159.88 | 66 | 156 |
| 68        | 210       | -142    | 66 | 157 |
| ND<0.05 U | 210       | -209.95 | 66 | 158 |
| 170       | 210       | -40     | 66 | 159 |
| 0.12      | 210       | -209.88 | 66 | 160 |
| 13        | 210       | -197    | 66 | 161 |
| 130       | 210       | -80     | 66 | 162 |
| 0.12      | 210       | -209.88 | 66 | 163 |
| ND<0.05 U | 68        | -67.95  | 66 | 164 |
| 170       | 68        | 102     | 67 | 164 |
| 0.12      | 68        | -67.88  | 67 | 165 |
| 13        | 68        | -55     | 67 | 166 |
| 130       | 68        | 62      | 68 | 166 |
| 0.12      | 68        | -67.88  | 68 | 167 |
| 170       | ND<0.05 U | 169.95  | 69 | 167 |
| 0.12      | ND<0.05 U | 0.07    | 70 | 167 |
| 13        | ND<0.05 U | 12.95   | 71 | 167 |
| 130       | ND<0.05 U | 129.95  | 72 | 167 |
| 0.12      | ND<0.05 U | 0.07    | 73 | 167 |
| 0.12      | 170       | -169.88 | 73 | 168 |
| 13        | 170       | -157    | 73 | 169 |
| 130       | 170       | -40     | 73 | 170 |
| 0.12      | 170       | -169.88 | 73 | 171 |
| 13        | 0.12      | 12.88   | 74 | 171 |
| 130       | 0.12      | 129.88  | 75 | 171 |
| 0.12      | 0.12      | 0       | 75 | 171 |
| 130       | 13        | 117     | 76 | 171 |
| 0.12      | 13        | -12.88  | 76 | 172 |
| 0.12      | 130       | -129.88 | 76 | 173 |

S Statistic = 76 - 173 = -97

---

| Tied Group | Value | Members |
|------------|-------|---------|
| 1          | 140   | 2       |
| 2          | 160   | 2       |
| 3          | 210   | 2       |
| 4          | 0.12  | 2       |

---

| Time Period | Observations |
|-------------|--------------|
| 12/1/2001   | 1            |
| 7/1/2004    | 1            |
| 10/13/2017  | 1            |
| 10/21/2021  | 1            |
| 11/29/2021  | 1            |
| 1/12/2022   | 1            |
| 2/18/2022   | 1            |
| 3/15/2022   | 1            |
| 4/18/2022   | 1            |
| 5/19/2022   | 1            |
| 6/15/2022   | 1            |
| 7/27/2022   | 1            |
| 8/29/2022   | 1            |
| 9/21/2022   | 1            |
| 10/26/2022  | 1            |
| 1/12/2023   | 1            |
| 4/10/2023   | 1            |
| 7/19/2023   | 1            |
| 10/4/2023   | 1            |
| 2/8/2024    | 1            |
| 6/28/2024   | 1            |
| 8/27/2024   | 1            |
| 4/15/2025   | 1            |

There are 0 time periods with multiple data

---

A = 72

B = 0

C = 0

D = 0

E = 8

F = 0

a = 25806

b = 95634

c = 1012

Group Variance = 1429.67

Z-Score = -2.53895

Comparison Level at 95% confidence level = -1.65463 (downward trend)

**-2.53895 < -1.65463 indicating a downward trend**

## Mann-Kendall Trend Analysis

Parameter: Benzene

Location: TM04-PZM006

Original Data (Not Transformed)

Non-Detects Replaced with 1/2 DL

95% Confidence Level

| Xj  | Xk   | Xj - Xk | Positives | Negatives |
|-----|------|---------|-----------|-----------|
| 610 | 1400 | -790    | 0         | 1         |
| 653 | 1400 | -747    | 0         | 2         |
| 355 | 1400 | -1045   | 0         | 3         |
| 367 | 1400 | -1033   | 0         | 4         |
| 390 | 1400 | -1010   | 0         | 5         |
| 390 | 1400 | -1010   | 0         | 6         |
| 380 | 1400 | -1020   | 0         | 7         |
| 180 | 1400 | -1220   | 0         | 8         |
| 500 | 1400 | -900    | 0         | 9         |
| 500 | 1400 | -900    | 0         | 10        |
| 460 | 1400 | -940    | 0         | 11        |
| 400 | 1400 | -1000   | 0         | 12        |
| 500 | 1400 | -900    | 0         | 13        |
| 540 | 1400 | -860    | 0         | 14        |
| 500 | 1400 | -900    | 0         | 15        |
| 370 | 1400 | -1030   | 0         | 16        |
| 340 | 1400 | -1060   | 0         | 17        |
| 490 | 1400 | -910    | 0         | 18        |
| 140 | 1400 | -1260   | 0         | 19        |
| 400 | 1400 | -1000   | 0         | 20        |
| 380 | 1400 | -1020   | 0         | 21        |
| 64  | 1400 | -1336   | 0         | 22        |
| 653 | 610  | 43      | 1         | 22        |
| 355 | 610  | -255    | 1         | 23        |
| 367 | 610  | -243    | 1         | 24        |
| 390 | 610  | -220    | 1         | 25        |
| 390 | 610  | -220    | 1         | 26        |
| 380 | 610  | -230    | 1         | 27        |
| 180 | 610  | -430    | 1         | 28        |
| 500 | 610  | -110    | 1         | 29        |
| 500 | 610  | -110    | 1         | 30        |
| 460 | 610  | -150    | 1         | 31        |
| 400 | 610  | -210    | 1         | 32        |
| 500 | 610  | -110    | 1         | 33        |
| 540 | 610  | -70     | 1         | 34        |
| 500 | 610  | -110    | 1         | 35        |
| 370 | 610  | -240    | 1         | 36        |
| 340 | 610  | -270    | 1         | 37        |
| 490 | 610  | -120    | 1         | 38        |
| 140 | 610  | -470    | 1         | 39        |
| 400 | 610  | -210    | 1         | 40        |
| 380 | 610  | -230    | 1         | 41        |
| 64  | 610  | -546    | 1         | 42        |
| 355 | 653  | -298    | 1         | 43        |
| 367 | 653  | -286    | 1         | 44        |

|     |     |      |    |    |
|-----|-----|------|----|----|
| 390 | 653 | -263 | 1  | 45 |
| 390 | 653 | -263 | 1  | 46 |
| 380 | 653 | -273 | 1  | 47 |
| 180 | 653 | -473 | 1  | 48 |
| 500 | 653 | -153 | 1  | 49 |
| 500 | 653 | -153 | 1  | 50 |
| 460 | 653 | -193 | 1  | 51 |
| 400 | 653 | -253 | 1  | 52 |
| 500 | 653 | -153 | 1  | 53 |
| 540 | 653 | -113 | 1  | 54 |
| 500 | 653 | -153 | 1  | 55 |
| 370 | 653 | -283 | 1  | 56 |
| 340 | 653 | -313 | 1  | 57 |
| 490 | 653 | -163 | 1  | 58 |
| 140 | 653 | -513 | 1  | 59 |
| 400 | 653 | -253 | 1  | 60 |
| 380 | 653 | -273 | 1  | 61 |
| 64  | 653 | -589 | 1  | 62 |
|     |     |      |    |    |
| 367 | 355 | 12   | 2  | 62 |
| 390 | 355 | 35   | 3  | 62 |
| 390 | 355 | 35   | 4  | 62 |
| 380 | 355 | 25   | 5  | 62 |
| 180 | 355 | -175 | 5  | 63 |
| 500 | 355 | 145  | 6  | 63 |
| 500 | 355 | 145  | 7  | 63 |
| 460 | 355 | 105  | 8  | 63 |
| 400 | 355 | 45   | 9  | 63 |
| 500 | 355 | 145  | 10 | 63 |
| 540 | 355 | 185  | 11 | 63 |
| 500 | 355 | 145  | 12 | 63 |
| 370 | 355 | 15   | 13 | 63 |
| 340 | 355 | -15  | 13 | 64 |
| 490 | 355 | 135  | 14 | 64 |
| 140 | 355 | -215 | 14 | 65 |
| 400 | 355 | 45   | 15 | 65 |
| 380 | 355 | 25   | 16 | 65 |
| 64  | 355 | -291 | 16 | 66 |
|     |     |      |    |    |
| 390 | 367 | 23   | 17 | 66 |
| 390 | 367 | 23   | 18 | 66 |
| 380 | 367 | 13   | 19 | 66 |
| 180 | 367 | -187 | 19 | 67 |
| 500 | 367 | 133  | 20 | 67 |
| 500 | 367 | 133  | 21 | 67 |
| 460 | 367 | 93   | 22 | 67 |
| 400 | 367 | 33   | 23 | 67 |
| 500 | 367 | 133  | 24 | 67 |
| 540 | 367 | 173  | 25 | 67 |
| 500 | 367 | 133  | 26 | 67 |
| 370 | 367 | 3    | 27 | 67 |
| 340 | 367 | -27  | 27 | 68 |
| 490 | 367 | 123  | 28 | 68 |
| 140 | 367 | -227 | 28 | 69 |
| 400 | 367 | 33   | 29 | 69 |
| 380 | 367 | 13   | 30 | 69 |
| 64  | 367 | -303 | 30 | 70 |

|     |     |      |    |    |
|-----|-----|------|----|----|
| 390 | 390 | 0    | 30 | 70 |
| 380 | 390 | -10  | 30 | 71 |
| 180 | 390 | -210 | 30 | 72 |
| 500 | 390 | 110  | 31 | 72 |
| 500 | 390 | 110  | 32 | 72 |
| 460 | 390 | 70   | 33 | 72 |
| 400 | 390 | 10   | 34 | 72 |
| 500 | 390 | 110  | 35 | 72 |
| 540 | 390 | 150  | 36 | 72 |
| 500 | 390 | 110  | 37 | 72 |
| 370 | 390 | -20  | 37 | 73 |
| 340 | 390 | -50  | 37 | 74 |
| 490 | 390 | 100  | 38 | 74 |
| 140 | 390 | -250 | 38 | 75 |
| 400 | 390 | 10   | 39 | 75 |
| 380 | 390 | -10  | 39 | 76 |
| 64  | 390 | -326 | 39 | 77 |

|     |     |      |    |    |
|-----|-----|------|----|----|
| 380 | 390 | -10  | 39 | 78 |
| 180 | 390 | -210 | 39 | 79 |
| 500 | 390 | 110  | 40 | 79 |
| 500 | 390 | 110  | 41 | 79 |
| 460 | 390 | 70   | 42 | 79 |
| 400 | 390 | 10   | 43 | 79 |
| 500 | 390 | 110  | 44 | 79 |
| 540 | 390 | 150  | 45 | 79 |
| 500 | 390 | 110  | 46 | 79 |
| 370 | 390 | -20  | 46 | 80 |
| 340 | 390 | -50  | 46 | 81 |
| 490 | 390 | 100  | 47 | 81 |
| 140 | 390 | -250 | 47 | 82 |
| 400 | 390 | 10   | 48 | 82 |
| 380 | 390 | -10  | 48 | 83 |
| 64  | 390 | -326 | 48 | 84 |

|     |     |      |    |    |
|-----|-----|------|----|----|
| 180 | 380 | -200 | 48 | 85 |
| 500 | 380 | 120  | 49 | 85 |
| 500 | 380 | 120  | 50 | 85 |
| 460 | 380 | 80   | 51 | 85 |
| 400 | 380 | 20   | 52 | 85 |
| 500 | 380 | 120  | 53 | 85 |
| 540 | 380 | 160  | 54 | 85 |
| 500 | 380 | 120  | 55 | 85 |
| 370 | 380 | -10  | 55 | 86 |
| 340 | 380 | -40  | 55 | 87 |
| 490 | 380 | 110  | 56 | 87 |
| 140 | 380 | -240 | 56 | 88 |
| 400 | 380 | 20   | 57 | 88 |
| 380 | 380 | 0    | 57 | 88 |
| 64  | 380 | -316 | 57 | 89 |

|     |     |     |    |    |
|-----|-----|-----|----|----|
| 500 | 180 | 320 | 58 | 89 |
| 500 | 180 | 320 | 59 | 89 |
| 460 | 180 | 280 | 60 | 89 |
| 400 | 180 | 220 | 61 | 89 |
| 500 | 180 | 320 | 62 | 89 |

|     |     |      |    |     |
|-----|-----|------|----|-----|
| 540 | 180 | 360  | 63 | 89  |
| 500 | 180 | 320  | 64 | 89  |
| 370 | 180 | 190  | 65 | 89  |
| 340 | 180 | 160  | 66 | 89  |
| 490 | 180 | 310  | 67 | 89  |
| 140 | 180 | -40  | 67 | 90  |
| 400 | 180 | 220  | 68 | 90  |
| 380 | 180 | 200  | 69 | 90  |
| 64  | 180 | -116 | 69 | 91  |
| 500 | 500 | 0    | 69 | 91  |
| 460 | 500 | -40  | 69 | 92  |
| 400 | 500 | -100 | 69 | 93  |
| 500 | 500 | 0    | 69 | 93  |
| 540 | 500 | 40   | 70 | 93  |
| 500 | 500 | 0    | 70 | 93  |
| 370 | 500 | -130 | 70 | 94  |
| 340 | 500 | -160 | 70 | 95  |
| 490 | 500 | -10  | 70 | 96  |
| 140 | 500 | -360 | 70 | 97  |
| 400 | 500 | -100 | 70 | 98  |
| 380 | 500 | -120 | 70 | 99  |
| 64  | 500 | -436 | 70 | 100 |
| 460 | 500 | -40  | 70 | 101 |
| 400 | 500 | -100 | 70 | 102 |
| 500 | 500 | 0    | 70 | 102 |
| 540 | 500 | 40   | 71 | 102 |
| 500 | 500 | 0    | 71 | 102 |
| 370 | 500 | -130 | 71 | 103 |
| 340 | 500 | -160 | 71 | 104 |
| 490 | 500 | -10  | 71 | 105 |
| 140 | 500 | -360 | 71 | 106 |
| 400 | 500 | -100 | 71 | 107 |
| 380 | 500 | -120 | 71 | 108 |
| 64  | 500 | -436 | 71 | 109 |
| 400 | 460 | -60  | 71 | 110 |
| 500 | 460 | 40   | 72 | 110 |
| 540 | 460 | 80   | 73 | 110 |
| 500 | 460 | 40   | 74 | 110 |
| 370 | 460 | -90  | 74 | 111 |
| 340 | 460 | -120 | 74 | 112 |
| 490 | 460 | 30   | 75 | 112 |
| 140 | 460 | -320 | 75 | 113 |
| 400 | 460 | -60  | 75 | 114 |
| 380 | 460 | -80  | 75 | 115 |
| 64  | 460 | -396 | 75 | 116 |
| 500 | 400 | 100  | 76 | 116 |
| 540 | 400 | 140  | 77 | 116 |
| 500 | 400 | 100  | 78 | 116 |
| 370 | 400 | -30  | 78 | 117 |
| 340 | 400 | -60  | 78 | 118 |
| 490 | 400 | 90   | 79 | 118 |
| 140 | 400 | -260 | 79 | 119 |
| 400 | 400 | 0    | 79 | 119 |

|     |     |      |    |     |
|-----|-----|------|----|-----|
| 380 | 400 | -20  | 79 | 120 |
| 64  | 400 | -336 | 79 | 121 |
| 540 | 500 | 40   | 80 | 121 |
| 500 | 500 | 0    | 80 | 121 |
| 370 | 500 | -130 | 80 | 122 |
| 340 | 500 | -160 | 80 | 123 |
| 490 | 500 | -10  | 80 | 124 |
| 140 | 500 | -360 | 80 | 125 |
| 400 | 500 | -100 | 80 | 126 |
| 380 | 500 | -120 | 80 | 127 |
| 64  | 500 | -436 | 80 | 128 |
| 500 | 540 | -40  | 80 | 129 |
| 370 | 540 | -170 | 80 | 130 |
| 340 | 540 | -200 | 80 | 131 |
| 490 | 540 | -50  | 80 | 132 |
| 140 | 540 | -400 | 80 | 133 |
| 400 | 540 | -140 | 80 | 134 |
| 380 | 540 | -160 | 80 | 135 |
| 64  | 540 | -476 | 80 | 136 |
| 370 | 500 | -130 | 80 | 137 |
| 340 | 500 | -160 | 80 | 138 |
| 490 | 500 | -10  | 80 | 139 |
| 140 | 500 | -360 | 80 | 140 |
| 400 | 500 | -100 | 80 | 141 |
| 380 | 500 | -120 | 80 | 142 |
| 64  | 500 | -436 | 80 | 143 |
| 340 | 370 | -30  | 80 | 144 |
| 490 | 370 | 120  | 81 | 144 |
| 140 | 370 | -230 | 81 | 145 |
| 400 | 370 | 30   | 82 | 145 |
| 380 | 370 | 10   | 83 | 145 |
| 64  | 370 | -306 | 83 | 146 |
| 490 | 340 | 150  | 84 | 146 |
| 140 | 340 | -200 | 84 | 147 |
| 400 | 340 | 60   | 85 | 147 |
| 380 | 340 | 40   | 86 | 147 |
| 64  | 340 | -276 | 86 | 148 |
| 140 | 490 | -350 | 86 | 149 |
| 400 | 490 | -90  | 86 | 150 |
| 380 | 490 | -110 | 86 | 151 |
| 64  | 490 | -426 | 86 | 152 |
| 400 | 140 | 260  | 87 | 152 |
| 380 | 140 | 240  | 88 | 152 |
| 64  | 140 | -76  | 88 | 153 |
| 380 | 400 | -20  | 88 | 154 |
| 64  | 400 | -336 | 88 | 155 |
| 64  | 380 | -316 | 88 | 156 |

S Statistic =  $88 - 156 = -68$

---

| Tied Group | Value | Members |
|------------|-------|---------|
| 1          | 390   | 2       |
| 2          | 380   | 2       |
| 3          | 500   | 4       |
| 4          | 400   | 2       |

---

| Time Period | Observations |
|-------------|--------------|
| 12/1/2001   | 1            |
| 7/1/2004    | 1            |
| 10/13/2017  | 1            |
| 10/21/2021  | 1            |
| 11/29/2021  | 1            |
| 1/12/2022   | 1            |
| 2/18/2022   | 1            |
| 3/15/2022   | 1            |
| 4/18/2022   | 1            |
| 5/19/2022   | 1            |
| 6/15/2022   | 1            |
| 7/27/2022   | 1            |
| 8/29/2022   | 1            |
| 9/21/2022   | 1            |
| 10/26/2022  | 1            |
| 1/12/2023   | 1            |
| 4/10/2023   | 1            |
| 7/19/2023   | 1            |
| 10/4/2023   | 1            |
| 2/8/2024    | 1            |
| 6/28/2024   | 1            |
| 8/27/2024   | 1            |
| 4/15/2025   | 1            |

There are 0 time periods with multiple data

---

A = 210

B = 0

C = 24

D = 0

E = 18

F = 0

a = 25806

b = 95634

c = 1012

Group Variance = 1422

Z-Score = -1.77674

Comparison Level at 95% confidence level = -1.65463 (downward trend)

**-1.77674 < -1.65463 indicating a downward trend**