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ENVIRONMENT

Mr. Robert Hill
Maryland Department of Environment
Oil Control Program
1800 Washington Boulevard
Baltimore, MD 21230

Subject:
Third Quarter 2011 Remedial Action Progress Report
Former Exxon Facility #25553
143 Frederick Road
Thurmont, Frederick County, Maryland
MDE Case No. 9-0923-FR (1989-0923-FR)

Date:
18 October 2011

Contact:
Rusty Kahl

Phone:
410-923-7813

Email:
rusty.kahl@arcadis-us.com

Our ref:
B0085851.0003

Dear Mr. Hill:

ARCADIS U.S. (ARCADIS), on behalf of ExxonMobil Environmental Services Company (EMESC) on behalf ExxonMobil Corporation (ExxonMobil), is pleased to submit the Remedial Action Progress Report (RAPR) for activities performed during Third Quarter 2011 at the Former Exxon Station #25553, located at 143 Frederick Road, Thurmont, Maryland.

Please contact Rusty Kahl of ARCADIS at (410) 923-7813 if you have any questions regarding the information contained in this report.

Sincerely,

ARCADIS U.S., Inc.

Sandra Kamas	William R. Kahl, P.G.	Michael Cleary
Geologist II	Project Geologist	Associate Project Manager

Copies:
Ms. Jewel Cox, ExxonMobil

Third Quarter 2011 Groundwater Monitoring Report
Former Exxon Facility #25553, 143 Frederick Road, Thurmont, MD

Regulatory Information

Regulatory Agency:	Maryland Department of the Environment (MDE)
MDE Case Number:	9-0923-FR (1989-0923-FR)
Agency Contact:	Mr. Robert Hill
Current Case Status:	Operation of soil vapor extraction (SVE) remediation system with quarterly groundwater monitoring sampling
Reporting Period:	1 July 2011 through 30 September 2011
Last Report:	First Quarter 2010 Remedial Action Progress Report, August 2011

General Site Information

ExxonMobil Contact:	Ms. Jewel Cox
Consultant Contact:	Mr. Rusty Kahl
Facility Status:	Former Exxon service station – 143 Frederick Road is now operating as a bank. Remediation System and monitoring wells are located across MD Rt. 806 and 140 Frederick Road, which is currently an active 7-11 gasoline station and truck stop, owned by Premier Associates Group
Area Property Use:	See local area map (**Figure 1**)
Monitoring Wells:	MW-5, MW-6, MW-11, MW-13A, MW-37, MW-38, MW-46, MW-48, MW-50, MW-53D, MW-53I, MW-53S, MW-54D, MW-54S, MW-55D, MW-55I, MW-55S, VW-1 through VW-3, MP-1 through MP-3, and DPE-1. (**Figure 2**)
Recovery Wells:	RW-52 (Former GWPT), MW-14A, and MW-51 (SVE)
Abandoned Wells:	MW-4, MW-7, MW-9, MW-10, MW-12, MW-13, MW-14 through MW-24, MW-30 through MW-33, MW-35, MW-39 through MW42, MW-44, MW-47, MW-49, and VMW-4

Destroyed Wells:	MW-1 through MW-3, MW-8, MW-25, MW-26, MW-27, MW-29, MW-34, MW-36, MW-43, MW-45, and VW-5 through VW-7
Site Geology:	Dense silt, clay, and cobble overlaying competent fractured limestone bedrock

Activities Completed this Period

2-3 August 2011 – Groundwater Gauging and Sampling

Wells Gauged and Sampled:	MW-5, MW-6, MW-13A, MW-14A, MW-37, MW-38, MW-46, MW-48, MW-50, MW-51, MW-53D, MW-53I, MW-53S, MW-54D, MW-55S, MW-55I, MW-55D, VW-1 through VW-3, MP-1, and DPE-1.
Wells Gauged Only:	MW-11 and MW-53S were not sampled due to insufficient water volume, MP-2 was not sampled due to the detection of 0.01 feet of product, and MP-3 was not sampled due to an obstruction.
Liquid Phase Hydrocarbon:	0.01 feet was detected in well MP-2
Minimum/Maximum Depth to Water:	7.69 feet (MW-51) / 22.08 feet (MW-48)
Groundwater Flow Direction:	Southeast

Groundwater samples were collected 2 August through 3 August 2011 and submitted to TestAmerica Laboratories of Nashville, Tennessee for analysis of full list volatile organic compounds (VOCs) and fuel oxygenates using Environmental Protection Agency (EPA) Method 8260B, and total petroleum hydrocarbons – gasoline range organics (TPH-GRO) and total petroleum hydrocarbons – diesel range organics (TPH-DRO) by EPA Method 8015. Monitoring well gauging and sampling data are summarized on **Table 1** and depicted on **Figure 3**. During this reporting period liquid phase hydrocarbon (LPH) was detected at one well (MP-2), at a thickness of 0.01 feet. Quality Assurance/Quality Control (QA/QC) samples collected included two duplicate samples (collected from MW-14A and MW-51), a daily rinsate blank, and trip blanks. The TestAmerica laboratory groundwater analytical report is attached as **Appendix A**.

Groundwater Sampling Results

Benzene, toluene, ethylbenzene, methyl tertiary butyl ether (MTBE), TPH-DRO, and TPH-GRO were detected at concentrations exceeding MDE Groundwater (GW) Clean-up Standards in samples collected from one or more wells:

- Benzene was detected at concentrations exceeding the MDE GW standard (5 µg/L) in 15 samples (MW-13A, MW-14A, MW-38, MW-51, MW-53I, MW-53S, MW-53D, MW-54D, MW-55I, MW-55S, MW-55D, RW-52, DPE-1, MP-1, and VW-1) with a maximum concentration of 14,900 µg/L in the sample collected from MW-53S.

- Toluene was detected at concentrations exceeding the MDE GW standard (1,000 µg/L) in four samples (MW-53S, MW-55I, DPE-1, and MP-1) with a maximum concentration of 14,300 µg/L in the sample collected from DPE-1.

- Ethylbenzene was detected at concentrations exceeding the MDE GW standard (700 µg/L) in six samples (MW-53S, MW-54D, MW-55I, DPE-1, MP-1, and VW-1) with a maximum concentration of 1,830 µg/L in the sample collected from DPE-1.

- MTBE was detected at concentrations exceeding the MDE GW standard (20 µg/L) in eight samples (MW-13A, MW-14A, MW-53I, MW-53S, MW-54D, DPE-1, MP-1, and VW-1) with a maximum concentration of 527 µg/L in the sample collected from MW-53S.

- TPH-GRO was detected above the MDE GW standard (47 µg/L) in 17 samples (MW-6, MW-13A, MW-14A, MW-38, MW-50, MW-51, MW-53I, MW-53S, MW-54D, MW-55S, MW-55I, MW-55D, RW-52, DPE-1, MP-1, VW-1, and VW-2) with a maximum concentration of 118,000 µg/L in the sample collected from MW-53S.

- TPH-DRO was detected above the MDE GW standard (47 µg/L) in 20 samples (MW-5, MW-6, MW-13A, MW-14A, MW-38, MW-46, MW-48, MW-50, MW-51, MW-53S, MW-53I, MW-54D, MW-55S, MW-55I, MW-55D, DPE-1, MP-1, RW-52, VW-1, and VW-2) with a maximum concentration of 3,280 µg/L in the sample collected from MW-54D.

Activities Planned for Next Period (Fourth Quarter 2011)

Activities proposed for the Fourth Quarter 2011 include quarterly groundwater gauging and sampling to support remedial action to be conducted during the second half of 2011. Groundwater samples will be collected from wells MW-5, MW-6, MW-11, MW-13A, MW-14A, MW-46, MW-48, MW-50, MW-51, MW-53D, MW-53I, MW-53S, MW-54S, MW-54D, MW-55D, MW-55I, MW-55S, VW-1 through VW-3, MP-1 through MP-3, DPE-1, and RW-52 and analyzed for BTEX AND MTBE EPA Method 8260B, and TPH-GRO and TPH-DRO using EPA Method 8015.

Figures and Tables

Table 1: Groundwater Monitoring and Analytical Data
Figure 1: Local Area Map
Figure 2: Site Map
Figure 3: Groundwater Elevation Contour Map and Analytical Data – Shallow (2 August 2011)
Figure 4: Groundwater Elevation Contour Map and Analytical Data – Bedrock (2 August 2011)

Appendices
Appendix A: TestAmerica Laboratory Analytical Report (16 August 2011)

Tables

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-1	4/12/2005	467.19	0.00	ND	ND	467.19	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/26/2005	467.19	0.00	ND	ND	467.19	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	4/20/2006	467.19	7.19	ND	ND	460.00	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.94 J	NA	NA	
	11/7/2006	467.19	1.40	ND	ND	465.79	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	4/26/2007	467.19	0.00	ND	ND	467.19	ND(1.0)	1.3	0.61 J	0.77 J	2.7 J	ND(1.0)	NA	NA	
	10/11/2007	467.19	0.00	ND	ND	467.19	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/21/2008	467.19	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	467.19	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	467.19	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well destroyed
MW-2	4/12/2005	466.92	0.00	ND	ND	466.92	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/26/2005	466.92	1.79	ND	ND	465.13	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	4/20/2006	466.92	5.47	ND	ND	461.45	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	11/7/2006	466.92	0.00	ND	ND	466.92	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/14/2006	466.92	1.10	ND	ND	465.82	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	466.92	0.00	ND	ND	466.92	0.70 J	3.8	1.8	6.1	12.4 J	ND(1.0)	NA	NA	
	10/11/2007	466.92	2.80	ND	ND	464.12	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/24/2007	466.92	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well destroyed

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		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-5	4/12/2005	465.31	10.31	ND	ND	455.00	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/26/2005	465.31	20.35	ND	ND	444.96	2.0	8.1	1.2	6.0	17.3	0.84 J	NA	NA	
	4/20/2006	465.31	12.39	ND	ND	452.92	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	11/7/2006	465.31	16.60	ND	ND	448.71	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/14/2006	465.31	14.17	ND	ND	451.14	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	465.31	17.60	ND	ND	447.71	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/21/2008	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	465.31	5.90	ND	ND	459.41	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	465.31	12.21	ND	ND	453.10	3.2	1.2	0.64 J	0.76 J	5.8 J	ND(1.0)	NA	NA	
	1/13/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	465.31	9.43	ND	ND	455.88	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	5/5/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	465.31	12.16	ND	ND	453.15	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/9/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	465.31	6.51	ND	ND	458.80	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	ND(200)	
	4/30/2010	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

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Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-5	5/16/2010	465.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	465.31	10.59	ND	ND	454.72	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	9/2/2010	465.31	13.09	ND	ND	452.22	0.33 J	ND(1.0)	ND(1.0)	ND(1.0)	0.33 J	ND(1.0)	ND(100)	ND(200)	
	12/8/2010	465.31	11.28	ND	ND	454.03	0.81 J	0.39 J	ND(1.0)	ND(1.0)	1.20 J	ND(1.0)	0.189	ND(200)	
	1/12/2011	465.31	11.92	ND	ND	453.39	**10.2**	5.43	6.44	23.6	45.67	<1	**278**	**227**	
	5/9/2011	465.31	5.71	ND	ND	459.60	<1	<1	<1	<3	BRL	<1	**155**	<100	
	8/2/2011	465.31	11.59	ND	ND	453.72	4.22	<1	<1	<3	4.22	<1	**310**	<100	

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Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-6	4/12/2005	466.34	8.38	ND	ND	457.96	**43.2**	1.6	0.79 J	0.71 J	46.3 J	**49.6**	NA	NA	
	10/26/2005	466.34	16.01	ND	ND	450.33	**186**	19.1	34.3	73.5	313	**28.3**	NA	NA	
	4/20/2006	466.34	15.71	ND	ND	450.63	**186**	11.5	10.5	2.7	211	16.9	NA	NA	
	11/7/2006	466.34	14.75	ND	ND	451.59	**397**	33.7	20.5	110	561	11.0	NA	NA	
	12/14/2006	466.34	12.66	ND	ND	453.68	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	466.34	7.95	ND	ND	458.39	4.4	ND(1.0)	ND(1.0)	ND(1.0)	4.4	0.55 J	NA	NA	
	10/11/2007	466.34	19.75	ND	ND	446.59	**69.8**	4.7	0.98 J	15.3	90.8 J	9.2	NA	NA	
	1/21/2008	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	466.34	4.57	ND	ND	461.77	**316**	20.0	10.1	54.1	400	4.3	NA	NA	Gauged on 5/19/08
	10/28/2008	466.34	8.36	ND	ND	457.98	**9.0**	0.24 J	ND(1.0)	ND(1.0)	9.2 J	0.60 J	NA	NA	
	1/13/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	466.34	6.35	ND	ND	459.99	1.1	ND(1.0)	ND(1.0)	ND(1.0)	1.1	ND(1.0)	NA	NA	
	5/5/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	466.34	10.02	ND	ND	456.32	4.8	0.96 J	3.7	2.0	11.5 J	ND(1.0)	NA	NA	
	12/9/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	466.34	6.36	ND	ND	459.98	5.0	0.85 J	1.6	0.39 J	7.8 J	ND(1.0)	NA	**558**	

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MW-6	4/30/2010	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	466.34	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	466.34	9.24	ND	ND	457.10	**76.7**	ND(1.0)	0.69 J	0.32 J	77.7 J	ND(1.0)	**342**	**610**	
	9/2/2010	466.34	10.12	ND	ND	456.22	2.6	ND(1.0)	2.2	ND(1.0)	4.8	ND(1.0)	**208**	**329**	
	12/8/2010	466.34	6.69	ND	ND	459.65	**95.4**	ND(1.0)	ND(1.0)	ND(1.0)	95.4	ND(1.0)	**181**	**580**	
	1/12/2011	466.34	8.19	ND	ND	458.15	**11.2 [11.3]**	3.62	5.68 [6.38]	11.8	32.3 [33.72]	<1 [<1]	**159 [110]**	**586 [700]**	
	5/9/2011	466.34	4.96	ND	ND	461.38	1.04	1.96	10.9	5.89	19.79	<1	**125**	**481**	sheen
	8/2/2011	466.34	9.54	ND	ND	456.80	<1	<1	1.06	3.22	4.28	<1	**307**	**609**	

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Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-11	4/12/2005	466.63	13.00	ND	ND	453.63	3.5	ND(1.0)	1.2	ND(1.0)	4.7	ND(1.0)	NA	NA	
	5/3/2005	466.63	16.13	ND	ND	450.50	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	466.63	14.75	ND	ND	451.88	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	466.63	14.15	ND	ND	452.48	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/20/2005	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	466.63	16.11	ND	ND	450.52	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/25/2007	466.63	11.80	ND	ND	454.83	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	466.63	11.80	ND	ND	454.83	**83.4**	6.2	96.0	33.0	218.6	11.2	NA	NA	
	7/19/2007	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-11	2/19/2008	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	466.63	7.65	ND	ND	458.98	**656**	5.8	22.0	11.9	696	**51.6**	NA	NA	
	10/28/2008	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/13/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	466.63	14.34	ND	ND	452.29	0.30 J	ND(1.0)	ND(1.0)	ND(1.0)	0.30 J	ND(1.0)	NA	NA	
	5/5/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/9/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	466.63	8.30	ND	ND	458.33	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	ND(200)	
	4/30/2010	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	466.63	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well inaccessible
	9/2/2010	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/11/2011	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/9/2011	466.63	7.92	ND	ND	458.71	**55.7**	<1	<1	<3	55.7	1.36	**1,300**	**341**	
	8/2/2011	466.63	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-13A	4/12/2005	464.13	19.80	ND	ND	444.33	**292**	71.8	381	368	1,113	**137**	NA	NA	
	5/3/2005	464.13	22.50	ND	ND	441.63	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	464.13	28.70	ND	ND	435.43	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	464.13	31.32	ND	ND	432.81	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	464.13	35.68	ND	ND	428.45	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	464.13	37.74	ND	ND	426.39	**26.0**	1.4	9.2	8.3	44.9	1.6	NA	NA	
	12/20/2005	464.13	35.68	ND	ND	428.45	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	464.13	35.42	ND	ND	428.71	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	464.13	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	464.13	32.33	ND	ND	431.80	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	464.13	34.75	ND	ND	429.38	4.6	4.8	3.4	1.4	14.2	0.88 J	NA	NA	
	5/31/2006	464.13	32.98	ND	ND	431.15	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	464.13	30.75	ND	ND	433.38	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	464.13	30.35	ND	ND	433.78	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	464.13	30.47	ND	ND	433.66	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	464.13	29.90	ND	ND	434.23	**61.6**	25.2	48.1	238	373	5.6	NA	NA	
	11/22/2006	464.13	28.08	ND	ND	436.05	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	464.13	23.44	ND	ND	440.69	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	464.13	29.81	ND	ND	434.32	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	464.13	16.28	ND	ND	447.85	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	464.13	16.02	ND	ND	448.11	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	464.13	29.76	ND	ND	434.37	NS	NS	NS	NS	NS	NS	NS	NS	
	4/25/2007	464.13	11.69	ND	ND	452.44	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	464.13	11.70	ND	ND	452.43	**85.8**	6.6	103	37.8	233	12.3	NA	NA	
	7/19/2007	464.13	19.88	ND	ND	444.25	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	464.13	28.47	ND	ND	435.66	**797**	147	360	1,060	2,364	**82.7**	NA	NA	
	1/21/2008	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	464.13	9.85	ND	ND	454.28	**143**	7.2	86.4	42.5	279	9.1	NA	NA	Gauged on 5/19/08

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-13A	10/28/2008	464.13	22.40	ND	ND	441.73	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/13/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	464.13	16.91	ND	ND	447.22	**152**	20.0	234	63.5	470	4.9	NA	NA	
	5/5/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	464.13	22.05	ND	ND	442.08	0.32 J	ND(1.0)	ND(1.0)	ND(1.0)	0.32 J	ND(1.0)	NA	NA	
	12/9/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	464.13	10.39	ND	ND	453.74	4.7	0.44 J	2.4	0.99 J	8.5 J	0.45 J	NA	ND(200)	
	4/30/2010	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	464.13	19.91	ND	ND	444.22	**15.7**	2.2	13.9	4.6	36.4	1.4	**454**	**1,390**	
	9/2/2010	464.13	25.61	ND	ND	438.52	2.4	ND(1.0)	ND(1.0)	ND(1.0)	2.4	ND(1.0)	**555**	ND(200)	
	12/8/2010	464.13	23.40	ND	ND	440.73	1.6	ND(1.0)	ND(1.0)	ND(1.0)	1.6	ND(1.0)	**299**	ND(200)	
	1/11/2011	464.13	24.33	ND	ND	439.80	**725**	75.2	443	291	1,534.2	**60.1**	**2,110**	**5,370**	
	5/9/2011	464.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	under mulch pile
	8/2/2011	464.13	21.58	ND	ND	442.55	**804**	58.5	366	203	1,431.5	**49.6**	**1,560**	**8,770**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-14A	4/12/2005	463.23	19.45	ND	ND	443.78	**956**	381	340	2,150	3,827	**239**	NA	NA	
	5/3/2005	463.23	22.81	ND	ND	440.42	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	463.23	29.25	ND	ND	433.98	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	463.23	32.00	ND	ND	431.23	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	463.23	37.55	ND	ND	425.68	NS	NS	NS	NS	NS	NS	NS	NS	Trace LPH detected <0.01 thickness
	12/20/2005	463.23	37.76	ND	ND	425.47	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	463.23	37.78	ND	ND	425.45	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	463.23	35.89	ND	ND	427.34	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	463.23	33.10	ND	ND	430.13	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	463.23	36.49	ND	ND	426.74	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	463.23	33.80	ND	ND	429.43	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	463.23	33.35	ND	ND	429.88	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	463.23	31.10	ND	ND	432.13	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	463.23	30.52	ND	ND	432.71	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	463.23	30.52	ND	ND	432.71	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	463.23	30.20	ND	ND	433.03	**521**	301	311	2,680	3,813	**122**	NA	NA	
	12/22/2006	463.23	30.15	ND	ND	433.08	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	463.23	17.11	ND	ND	446.12	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	463.23	16.83	ND	ND	446.40	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	463.23	30.09	ND	ND	433.14	NS	NS	NS	NS	NS	NS	NS	NS	
	7/19/2007	463.23	20.02	ND	ND	443.21	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	463.23	28.40	ND	ND	434.83	**667**	630	**756**	2,820	4,873	**144**	NA	NA	
	10/24/2007	463.23	28.44	ND	ND	434.79	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	463.23	26.42	ND	ND	436.81	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	463.23	20.40	ND	ND	442.83	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	463.23	9.85	ND	ND	453.38	**166**	4.3	159	172	501	9.5	NA	NA	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-14A	10/28/2008	463.23	7.00	ND	ND	456.23	**329**	170	388	1,550	2,437	**44.1**	NA	NA	
	1/13/2009	463.23	23.05	ND	ND	440.18	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	463.23	20.38	ND	ND	442.85	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	463.23	20.25	ND	ND	442.98	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	463.23	18.59	ND	ND	444.64	**166**	17.6	154	486	824	16.2	NA	NA	
	5/5/2009	463.23	18.36	ND	ND	444.87	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	463.23	18.26	ND	ND	444.97	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	463.23	18.42	ND	ND	444.81	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	463.23	18.15	ND	ND	445.08	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	463.23	18.32	ND	ND	444.91	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	463.23	18.47	ND	ND	444.76	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	463.23	22.07	ND	ND	441.16	**363**	87.8	324	1,430	2,205	**29.8**	NA	NA	
	12/9/2009	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	463.23	18.18	ND	ND	445.05	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well inaccessible
	3/23/2010	463.23	15.30	ND	ND	447.93	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	463.23	11.80	ND	ND	451.43	**10.5**	0.36 J	15.8	15.8	42.5 J	0.72 J	NA	404	
	4/30/2010	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	463.23	20.38	ND	ND	442.85	**130**	3.3	62.2	32.5	228	7.2	**1,180**	**1,710**	
	7/7/2010	463.23	20.12	ND	ND	443.11	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	463.23	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	463.23	25.39	ND	ND	437.84	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	463.23	25.04	ND	ND	438.19	**311**	164	319	909	1,703	**45.8**	**3,570**	**9,090**	
	9/16/2010	463.23	26.22	ND	ND	437.01	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	463.23	25.58	ND	ND	437.65	NS	NS	NS	NS	NS	NS	NS	NS	
	10/20/2010	463.23	18.01	ND	ND	445.22	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	463.23	24.80	ND	ND	438.42	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	463.23	24.67	ND	ND	438.56	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	463.23	23.43	ND	ND	439.80	**199**	48	159	428	834	**42.4**	**5,550**	**3,950**	
	12/21/2010	463.23	24.68	ND	ND	438.55	NS	NS	NS	NS	NS	NS	NS	NS	
	1/11/2011	463.23	24.10	ND	ND	439.13	**707**	57.2	578	239	1,581.2	**60.8**	**3,010**	**7,450**	
	5/9/2011	463.23	11.05	ND	ND	452.18	**5.33 [5.76]**	<1 [<1]	4.5 [5.14]	<3 [3.08]	9.83 [10.9]	<1 [1.06]	**646 [831]**	**134 [114]**	
	8/2/2011	463.23	21.67	ND	ND	441.56	**637 [693]**	35.7 [42.7]	425 [448]	126 [133]	1,223.7 [1,316.7]	**46 [50]**	**2,960 [2,430]**	**8,720 [42,600]**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-23	4/12/2005	460.57	22.12	ND	ND	438.45	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/26/2005	460.57	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	460.57	44.68	ND	ND	415.89	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	11/7/2006	460.57	31.72	ND	ND	428.85	ND(1.0)	ND(1.0)	ND(1.0)	0.36 J	0.36 J	ND(1.0)	NA	NA	
	4/26/2007	460.57	12.30	ND	ND	448.27	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/11/2007	460.57	30.71	ND	ND	429.86	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/21/2008	460.57	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	460.57	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	460.57	8.55	ND	ND	452.02	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	460.57	NM	NM	NM	NM	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/13/2009	460.57	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	460.57	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	460.57	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	460.57	17.97	ND	ND	442.60	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Well destroyed/abandoned

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-37	4/12/2005	451.04	4.00	ND	ND	447.04	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/26/2005	451.04	14.02	ND	ND	437.02	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	4/20/2006	451.04	9.90	ND	ND	441.14	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	11/7/2006	451.04	12.05	ND	ND	438.99	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	4/26/2007	451.04	3.70	ND	ND	447.34	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/11/2007	451.04	12.70	ND	ND	438.34	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/21/2008	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	451.04	2.99	ND	ND	448.05	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	451.04	9.72	ND	ND	441.32	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/13/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	451.04	3.49	ND	ND	447.55	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	5/5/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	451.04	10.13	ND	ND	440.91	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/9/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	451.04	3.05	ND	ND	447.99	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	ND(200)	
	4/30/2010	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-37	5/16/2010	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	451.04	7.78	ND	ND	443.26	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	9/2/2010	451.04	11.65	ND	ND	439.39	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	12/8/2010	451.04	9.00	ND	ND	442.04	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	1/11/2011	451.04	9.25	ND	ND	441.79	<1	<1	<1	<3	BRL	<1	<102	<100	
	5/9/2011	451.04	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/2/2011	451.04	10.51	NM	NM	440.53	<1	<1	<1	<3	BRL	<1	<111	<100	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-38	4/12/2005	454.65	9.81	ND	ND	444.84	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/26/2005	454.65	23.31	ND	ND	431.34	1.9	ND(1.0)	ND(1.0)	ND(1.0)	1.9	1.8	NA	NA	
	4/20/2006	454.65	20.29	ND	ND	434.36	**164**	0.47 J	1.3	ND(1.0)	166 J	**37.8**	NA	NA	
	11/7/2006	454.65	18.79	ND	ND	435.86	**108**	0.52 J	1.1	ND(1.0)	110 J	15.9	NA	NA	
	4/26/2007	454.65	5.20	ND	ND	449.45	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/11/2007	454.65	19.27	ND	ND	435.38	**90.3**	1.6	13.4	0.44 J	105.7 J	**22.1**	NA	NA	
	1/21/2008	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	454.65	3.19	ND	ND	451.46	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	454.65	14.48	ND	ND	440.17	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.55 J	NA	NA	
	1/13/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	454.65	7.43	ND	ND	447.22	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	5/5/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	454.65	14.31	ND	ND	440.34	**5.8**	ND(1.0)	2.2	ND(1.0)	8.0	7.1	NA	NA	
	12/9/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	454.65	3.98	ND	ND	450.67	0.56 J	ND(1.0)	ND(1.0)	ND(1.0)	0.56 J	0.47 J	NA	ND(200)	
	4/30/2010	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-38	5/16/2010	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	454.65	12.22	ND	ND	442.43	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	9/2/2010	454.65	17.69	ND	ND	436.96	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	12/8/2010	454.65	15.55	ND	ND	439.10	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	1/11/2011	454.65	14.15	ND	ND	440.50	**56.5**	2.76	70.1	<3	129.36	9.03	**172**	**2,660**	
	5/9/2011	454.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/2/2011	454.65	13.84	NM	NM	440.81	**46.4**	<1	43.2	3	92.6	7.7	**123**	**2,560**	
MW-40	10/26/2005	446.69	24.00	ND	ND	422.69	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	446.69	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well abandoned
MW-43	10/26/2005	439.13	6.30	ND	ND	432.83	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	4/20/2006	439.13	15.23	ND	ND	423.90	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	11/7/2006	439.13	14.98	ND	ND	424.15	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	4/26/2007	439.13	7.00	ND	ND	432.13	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/11/2007	439.13	23.50	ND	ND	415.63	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/21/2008	439.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	439.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	439.13	4.49	ND	ND	434.64	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	NM	NM	NM	NM	NM	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/13/2009	439.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	439.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	439.13	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	439.13	4.41	ND	ND	434.72	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	

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Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-45	4/12/2005	456.38	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	456.38	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	456.38	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	456.38	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	456.38	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	456.38	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	456.38	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	456.38	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	456.38	10.57	ND	ND	445.81	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	456.38	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/13/2009	456.38	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	456.38	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	456.38	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	456.38	19.65	ND	ND	436.73	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	

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Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-46	4/12/2005	467.17	13.45	ND	ND	453.72	**5.6**	9.2	2.5	11.6	28.9	9.9	NA	NA	
	5/3/2005	467.17	15.55	ND	ND	451.62	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	467.17	23.59	ND	ND	443.58	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	467.17	23.70	ND	ND	443.47	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	467.17	24.25	ND	ND	442.92	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	467.17	24.00	ND	ND	443.17	0.42 J	238	ND(1.0)	ND(1.0)	238 J	ND(1.0)	NA	NA	
	12/20/2005	467.17	24.55	ND	ND	442.62	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	467.17	25.79	ND	ND	441.38	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	467.17	24.20	ND	ND	442.97	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2006	467.17	25.49	ND	ND	441.68	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	467.17	23.31	ND	ND	443.86	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	7/19/2007	467.17	15.50	ND	ND	451.67	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	467.17	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	467.17	5.17	ND	ND	462.00	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	2.9	NA	NA	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-46	10/28/2008	467.17	22.43	ND	ND	444.74	2.0	1.0	3.7	7.3	14.0	1.3	NA	NA	
	1/13/2009	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	467.17	10.98	ND	ND	456.19	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	1.4	NA	NA	
	5/5/2009	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	NM	NM	NM	NM	NM	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.58 J	NA	NA	
	12/9/2009	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	467.17	6.50	ND	ND	460.67	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	ND(200)	
	4/30/2010	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	467.17	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	467.17	15.93	ND	ND	451.24	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.49 J	ND(100)	ND(200)	
	9/2/2010	467.17	23.10	ND	ND	444.07	ND(1.0)	ND(1.0)	0.90 J	ND(1.0)	0.90 J	ND(1.0)	ND(100)	ND(200)	
	12/8/2010	467.17	23.18	ND	ND	443.99	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	1/12/2011	467.17	23.92	ND	ND	443.25	2.11	48.5	5.51	3.89	60.01	<1	**755**	**146**	
	5/9/2011	467.17	4.59	ND	ND	462.58	<1	<1	<1	<3	BRL	<1	**362**	<100	
	8/2/2011	467.17	19.95	ND	ND	447.22	<1	<1	<1	<3	BRL	<1	**412**	<100	
MW-47	4/12/2005	465.52	18.00	ND	ND	447.52	**96.5**	59.2	157	220	533	**321**	NA	NA	
	10/26/2005	465.52	26.52	ND	ND	439.00	1.8	ND(1.0)	3.6	1.1	6.5	**53.1**	NA	NA	
	4/20/2006	465.52	22.82	ND	ND	442.70	3.2	0.65 J	9.4	2.3	15.6 J	12.3	NA	NA	
	11/7/2006	465.52	25.45	ND	ND	440.07	4.1	3.8	4.3	11.2	23.4	0.64 J	NA	NA	
	12/14/2006	465.52	20.95	ND	ND	444.57	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	465.52	17.25	ND	ND	448.27	**38.7**	39.8	30.5	51.3	160.3	0.79 J	NA	NA	
	10/11/2007	465.52	26.45	ND	ND	439.07	0.82 J	ND(1.0)	ND(1.0)	ND(1.0)	0.82 J	ND(1.0)	NA	NA	
	1/21/2008	465.52	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	465.52	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well not located
	5/21/2008	465.52	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well not located
	10/28/2008	465.52	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well not located

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-48	4/12/2005	467.43	19.85	ND	ND	447.58	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	**41.8**	NA	NA	
	10/26/2005	467.43	27.04	ND	ND	440.39	1.0	54.5	0.67 J	2.9	59.1 J	19.1	NA	NA	
	4/20/2006	467.43	27.61	ND	ND	439.82	ND(1.0)	0.62 J	ND(1.0)	ND(1.0)	0.62 J	**38.6**	NA	NA	
	11/7/2006	467.43	26.55	ND	ND	440.88	**6.9**	8.8	3.4	17.4	36.5	5.9	NA	NA	
	12/14/2006	467.43	23.55	ND	ND	443.88	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	467.43	12.95	ND	ND	454.48	**43.0**	ND(1.0)	ND(1.0)	ND(1.0)	43.0	5.7	NA	NA	
	10/11/2007	467.43	26.78	ND	ND	440.65	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	4.0	NA	NA	
	2/19/2008	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	467.43	10.90	ND	ND	456.53	**12.1**	ND(1.0)	ND(1.0)	ND(1.0)	12.1	**20.3**	NA	NA	Gauged on 5/19/08
	10/28/2008	467.43	23.15	ND	ND	444.28	ND(1.0)	1.1	ND(1.0)	ND(1.0)	1.1	0.56 J	NA	NA	
	1/13/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	467.43	17.88	ND	ND	449.55	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	5.7	NA	NA	
	5/5/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/9/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	467.43	10.88	ND	ND	456.55	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	1.0	NA	ND(200)	
	4/30/2010	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	467.43	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
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		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-48	6/17/2010	467.43	20.12	ND	ND	447.31	ND(1.0)	0.60 J	ND(1.0)	ND(1.0)	0.60 J	0.55 J	**771**	ND(200)	
	9/2/2010	467.43	28.77	ND	ND	438.66	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	2.6	**4,800**	ND(200)	
	12/8/2010	467.43	23.41	ND	ND	444.02	ND(1.0)	12.2	ND(1.0)	ND(1.0)	12.2	0.90 J	**571**	ND(200)	
	1/11/2011	467.43	24.45	ND	ND	442.98	<1	<1	<1	<3	BRL	2.4	**972**	<100	
	5/9/2011	467.43	11.00	ND	ND	456.43	3.06	<1	<1	<3	3.06	<1	**481**	<100	
	8/2/2011	467.43	22.08	ND	ND	445.35	<1	1.32	<1	3.98	5.3	1.01	**214**	<100	

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		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-49	4/12/2005	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/3/2005	466.39	19.61	ND	ND	446.78	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	466.39	17.45	ND	ND	448.94	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/20/2005	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	466.39	17.38	ND	ND	449.01	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/25/2007	466.39	10.68	ND	ND	455.71	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	466.39	10.70	ND	ND	455.69	**8.2**	ND(1.0)	ND(1.0)	0.33 J	8.5 J	1.7	NA	NA	
	7/19/2007	466.39	16.90	ND	ND	449.49	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	466.39	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

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Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** **[5]**	**Toluene (µg/L)** **[1,000]**	**Ethyl-benzene (µg/L)** **[700]**	**Total Xylenes (µg/L)** **[10,000]**	**Total BTEX (µg/L)** **--**	**MTBE (µg/L)** **[20]**	**TPH-DRO (µg/L)** **[47]**	**TPH-GRO (µg/L)** **[47]**	**Comments**
MW-49	2/19/2008	466.39	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	466.39	8.16	ND	ND	458.23	**604**	37.2	63.8	44.1	749	**35.4**	NA	NA	
	10/28/2008	466.39	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/13/2009	466.39	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	466.39	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	466.39	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	466.39	14.73	ND	ND	451.66	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	

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Former Exxon Station #25553
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		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-50	4/12/2005	464.12	8.82	ND	ND	455.30	3.5	4.0	73.5	71.2	152.2	**53.9**	NA	NA	
	5/3/2005	464.12	9.30	ND	ND	454.82	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	464.12	10.29	ND	ND	453.83	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	464.12	11.78	ND	ND	452.34	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	464.12	13.16	ND	ND	450.96	**8.5**	5.1	24.4	16.3	54.3	6.5	NA	NA	
	12/20/2005	464.12	12.18	ND	ND	451.94	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	464.12	11.82	ND	ND	452.30	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	464.12	11.56	ND	ND	452.56	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	464.12	12.45	ND	ND	451.67	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	464.12	12.85	ND	ND	451.27	**114**	136	196	222	668	**75.6**	NA	NA	
	5/31/2006	464.12	12.91	ND	ND	451.21	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	464.12	14.69	ND	ND	449.43	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	464.12	13.40	ND	ND	450.72	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	464.12	12.65	ND	ND	451.47	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	464.12	12.59	ND	ND	451.53	**46.8**	49.4	147	157	400	**22.0**	NA	NA	
	11/22/2006	464.12	11.46	ND	ND	452.66	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	464.12	11.36	ND	ND	452.76	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	464.12	12.50	ND	ND	451.62	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	464.12	7.09	ND	ND	457.03	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	464.12	7.00	ND	ND	457.12	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	464.12	12.44	ND	ND	451.68	NS	NS	NS	NS	NS	NS	NS	NS	
	4/25/2007	464.12	7.53	ND	ND	456.59	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	464.12	7.55	ND	ND	456.57	**11.4**	15.0	26.2	179	232	**23.7**	NA	NA	
	7/19/2007	464.12	10.35	ND	ND	453.77	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	464.12	16.87	ND	ND	447.25	**16.8**	0.70 J	0.29 J	0.85 J	18.6 J	8.7	NA	NA	
	1/21/2008	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-50	5/21/2008	464.12	4.44	ND	ND	459.68	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	464.12	9.55	ND	ND	454.57	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/13/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	464.12	6.80	ND	ND	457.32	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	5/5/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well inaccessible
	12/9/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	464.12	5.96	ND	ND	458.16	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.29 J	NA	ND(200)	
	4/30/2010	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	464.12	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	464.12	9.98	ND	ND	454.14	0.32 J	0.48 J	7.0	7.3	15.1 J	1.4	ND(100)	**353**	
	9/2/2010	464.12	11.76	ND	ND	452.36	ND(1.0)	ND(1.0)	0.82 J	0.99 J	1.81 J	ND(1.0)	**444**	ND(200)	
	12/8/2010	464.12	9.43	ND	ND	454.69	ND(1.0)	0.45 J	6.3	6.3	13.1 J	0.62 J	**175**	**400**	
	1/12/2011	464.12	10.79	ND	ND	453.33	**7.9**	10.1	58.3	61.9	138.2	1.42	**1,060**	**3,600**	
	5/9/2011	464.12	5.75	ND	ND	458.37	<1	<1	4.08	<3	4.08	<1	<100	<100	sheen
	8/2/2011	464.12	11.18	ND	ND	452.94	2.3	1.75	8.94	9.23	22.22	<1	**170**	**1,210**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-51	4/12/2005	462.76	6.80	ND	ND	455.96	**114**	9.7	24.7	21.3	170	**67.2**	NA	NA	
	10/26/2005	462.76	14.46	ND	ND	448.30	**41.8**	3.7	20.6	17.8	83.9	**36.6**	NA	NA	
	4/20/2006	462.76	14.02	ND	ND	448.74	**17.7**	0.38 J	0.68 J	0.37 J	19.1 J	7.6	NA	NA	
	11/7/2006	462.76	13.10	ND	ND	449.66	**48.6**	20.8	9.2	45.0	123.6	8.8	NA	NA	
	12/14/2006	462.76	10.80	ND	ND	451.96	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	462.76	6.20	ND	ND	456.56	**274**	10.8	33.7	30.6	349	**23.5**	NA	NA	
	10/11/2007	462.76	17.98	ND	ND	444.78	**197**	6.2	4.1	8.8	216	**94.2**	NA	NA	
	1/21/2008	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	462.76	2.64	ND	ND	460.12	**38.2**	2.2	16.0	10.1	66.5	5.7	NA	NA	Gauged on 5/19/08
	10/28/2008	462.76	23.82	ND	ND	438.94	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.41 J	NA	NA	
	1/13/2009	462.76	7.39	ND	ND	455.37	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	462.76	5.07	ND	ND	457.69	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	462.76	4.96	ND	ND	457.80	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	462.76	4.46	ND	ND	458.30	**66.4**	6.5	21.7	13.1	107.7	0.74 J	NA	NA	
	5/5/2009	462.76	4.25	ND	ND	458.51	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	462.76	4.40	ND	ND	458.36	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	462.76	5.00	ND	ND	457.76	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	462.76	5.00	ND	ND	457.76	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	462.76	5.19	ND	ND	457.57	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	462.76	5.35	ND	ND	457.41	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	462.76	7.51	ND	ND	455.25	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/9/2009	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	462.76	4.00	ND	ND	458.76	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	462.76	3.56	ND	ND	459.20	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	462.76	5.00	ND	ND	457.76	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-51	3/31/2010	462.76	5.29	ND	ND	457.47	**156**	35.5	30.9	57.4	280	1.3	NA	**1,430**	
	4/30/2010	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	462.76	07.35	ND	ND	455.41	**359**	6.5	18.0	14.7	398.0	1.6 J	**264**	**1,860**	
	7/7/2010	462.76	07.18	ND	ND	455.58	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	462.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	462.76	5.91	ND	ND	456.85	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	462.76	6.43	ND	ND	456.33	**10.8**	ND(1.0)	0.47 J	ND(1.0)	11.3 J	ND(1.0)	ND(100)	ND(200)	
	9/16/2010	462.76	8.95	ND	ND	453.81	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	462.76	8.54	ND	ND	454.22	NS	NS	NS	NS	NS	NS	NS	NS	
	10/20/2010	462.76	8.15	ND	ND	454.61	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	462.76	8.71	ND	ND	454.05	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	462.76	8.76	ND	ND	454.00	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	462.76	3.86	ND	ND	458.90	3.3	0.53 J	0.70 J	3.4	7.9 J	ND(1.0)	**172**	**311**	
	12/21/2010	462.76	8.73	ND	ND	454.03	NS	NS	NS	NS	NS	NS	NS	NS	
	1/12/2011	462.76	5.60	ND	ND	457.16	**14.8**	3.56	3.87	8.26	30.49	<1	<96.2	**472**	
	5/9/2011	462.76	3.14	ND	ND	459.62	**16.1**	3.56	7.1	9.26	36.02	<1	**227**	**390**	
	8/2/2011	462.76	7.69	ND	ND	455.07	**8.7 [8.72]**	<1 [1.02]	<1 [1.03]	7.79 [8.06]	16.49 [18.83]	<1 [<1]	**200 [99]**	**553 [532]**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-53S	12/14/2006	469.08	22.20	ND	ND	446.88	**5,390**	**6,070**	685	4,110	16,255	**110**	NA	NA	
	1/3/2007	469.08	19.41	ND	ND	449.67	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	469.08	12.50	ND	ND	456.58	**16,100**	**23,100**	**1,960**	**12,100**	53,260	**238**	NA	NA	
	5/21/2007	469.08	12.55	ND	ND	456.53	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	469.08	12.46	ND	ND	456.62	NS	NS	NS	NS	NS	NS	NS	NS	
	8/3/2007	469.08	21.20	ND	ND	447.88	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	469.08	12.77	ND	ND	456.31	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	469.08	25.55	ND	ND	443.53	**5,610**	516	247	787	7,160	**585**	NA	NA	
	10/24/2007	469.08	26.98	ND	ND	442.10	NS	NS	NS	NS	NS	NS	NS	NS	
	12/4/2007	469.08	26.43	ND	ND	442.65	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	469.08	24.77	ND	ND	444.31	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	469.08	23.23	ND	ND	445.85	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	469.08	19.29	ND	ND	449.79	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	469.08	12.97	ND	ND	456.11	**14,800**	**12,200**	**1,390**	7,720	36,110	**272**	NA	NA	
	10/28/2008	469.08	21.39	ND	ND	447.69	**3,990**	**7,010**	603	3,460	15,063	**68.7**	NA	NA	
	1/13/2009	469.08	20.44	ND	ND	448.64	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	469.08	18.18	ND	ND	450.90	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	469.08	18.00	ND	ND	451.08	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	469.08	16.48	ND	ND	452.60	**8,870**	**11,200**	**752**	3,950	24,772	**122**	NA	NA	
	5/5/2009	469.08	16.23	ND	ND	452.85	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	469.08	16.37	ND	ND	452.71	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	469.08	16.49	ND	ND	452.59	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	469.08	16.78	ND	ND	452.30	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	469.08	17.05	ND	ND	452.03	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	469.08	17.18	ND	ND	451.90	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	469.08	20.10	ND	ND	448.98	**974**	**2,150**	148	970	4,242	12.7	NA	NA	
	12/9/2009	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-53S	1/22/2010	469.08	16.10	ND	ND	452.98	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	469.08	15.40	ND	ND	453.68	NS	NS	NS	NS	NS	NS	NS	NS	Well inaccessible
	3/23/2010	469.08	15.18	ND	ND	453.90	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	469.08	11.47	ND	ND	457.61	**11,100**	**11,700**	**991**	6,100	29,891	**389**	NA	**78,600**	
	4/30/2010	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	469.08	7.35	ND	ND	461.73	**20,000**	**23,700**	**1,770**	8,830	54,300	**645**	**3,440**	**133,000**	
	7/7/2010	469.08	7.16	ND	ND	461.92	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	469.08	22.75	ND	ND	446.33	**14,500**	**15,100**	672	4,980	35,252	**602**	**4,050**	**102,000**	
	9/16/2010	469.08	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	469.08	20.63	ND	ND	448.45	**14,200**	**15,600**	653	3,400	33,853	**844**	**1,180**	**72,200**	
	1/13/2011	469.08	20.55	NM	NM	448.53	**17,300**	**10,800**	**1,220**	5,690	35,010	**676**	**2,500**	**83,000**	
	5/9/2011	469.08	10.71	NM	NM	458.37	**19,200**	**26,000**	**1,440**	7,130	53,770	**761**	**2,060**	**110,000**	
	8/2/2011	469.08	19.36	NM	NM	449.72	**14,900**	**12,800**	**1,240**	5,310	34,250	**527**	**1,960**	**118,000**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
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Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-53I	12/14/2006	468.74	21.11	ND	ND	447.63	**2,090**	**2,110**	250	3,160	7,610	**186**	NA	NA	
	1/3/2007	468.74	19.07	ND	ND	449.67	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	468.74	12.65	ND	ND	456.09	**16,200**	**24,300**	**2,040**	**12,100**	54,640	**221**	NA	NA	
	5/21/2007	468.74	12.68	ND	ND	456.06	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	468.74	12.59	ND	ND	456.15	NS	NS	NS	NS	NS	NS	NS	NS	
	8/3/2007	468.74	21.05	ND	ND	447.69	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	468.74	12.84	ND	ND	455.90	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	469.09	25.70	ND	ND	443.39	2.0	0.41 J	ND(1.0)	0.59 J	3.0 J	0.87 J	NA	NA	
	10/24/2007	469.09	26.92	ND	ND	442.17	NS	NS	NS	NS	NS	NS	NS	NS	
	12/4/2007	468.74	26.01	ND	ND	442.73	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	468.74	24.50	ND	ND	444.24	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	468.74	23.21	ND	ND	445.53	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	468.74	19.23	ND	ND	449.51	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	468.74	12.78	ND	ND	455.96	**952**	64.2	22.2	41.3	1,080	**53.5**	NA	NA	Gauged on 5/19/08
	10/28/2008	468.74	21.31	ND	ND	447.43	**85.0**	17.9	7.3	18.0	128.2	9.9	NA	NA	
	1/13/2009	468.74	19.67	ND	ND	449.07	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	468.74	18.10	ND	ND	450.64	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	468.74	18.00	ND	ND	450.74	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	468.74	16.45	ND	ND	452.29	3.1	ND(1.0)	ND(1.0)	ND(1.0)	3.1	0.89 J	NA	NA	
	5/5/2009	468.74	16.25	ND	ND	452.49	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	468.74	16.37	ND	ND	452.37	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	468.74	16.52	ND	ND	452.22	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	468.74	16.72	ND	ND	452.02	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	468.74	17.00	ND	ND	451.74	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	468.74	17.18	ND	ND	451.56	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	468.74	19.80	ND	ND	448.94	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/9/2009	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-53I	1/22/2010	468.74	16.00	ND	ND	452.74	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	468.74	15.33	ND	ND	453.41	NS	NS	NS	NS	NS	NS	NS	NS	Well inaccessible
	3/23/2010	468.74	15.10	ND	ND	453.64	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	468.74	11.53	ND	ND	457.21	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	ND(200)	
	4/30/2010	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	468.74	18.58	ND	ND	450.16	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	205	ND(200)	
	7/7/2010	468.74	18.25	ND	ND	450.49	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	468.74	22.71	ND	ND	446.03	**222**	2.6	5.8	4.5	235	**20.5**	ND(100)	**1,110**	
	9/16/2010	468.74	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	468.74	20.50	ND	ND	448.24	**1,410**	56.5	24.4	24.3	1,515	**110**	ND(100)	**4,170**	
	1/13/2011	468.74	21.67	NM	NM	447.07	**3,060**	**1,410**	503	1,570	6,543	**161**	**861**	**18,100**	
	5/9/2011	468.74	10.40	NM	NM	458.34	**555**	159	49.6	205	969	**62.7**	**302**	**2,850**	
	8/2/2011	468.74	19.07	NM	NM	449.67	**1,800**	804	445	1,200	4,249	**69.5**	**835**	**18,000**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-53D	12/14/2006	469.09	21.47	ND	ND	447.62	**120**	99.9	23.9	117	361	12.5	NA	NA	
	1/3/2007	469.09	19.44	ND	ND	449.65	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	469.09	12.68	ND	ND	456.41	3.1	1.2	0.70 J	0.87 J	5.9 J	1.9	NA	NA	
	5/21/2007	469.09	12.71	ND	ND	456.38	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	469.09	12.60	ND	ND	456.49	NS	NS	NS	NS	NS	NS	NS	NS	
	8/3/2007	469.09	21.40	ND	ND	447.69	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	469.09	12.90	ND	ND	456.19	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2007	469.09	26.92	ND	ND	442.17	NS	NS	NS	NS	NS	NS	NS	NS	
	12/4/2007	469.09	26.50	ND	ND	442.59	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	469.09	24.78	ND	ND	444.31	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	469.09	23.57	ND	ND	445.52	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	469.09	19.58	ND	ND	449.51	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	469.09	10.12	ND	ND	458.97	0.62 J	ND(1.0)	0.87 J	ND(1.0)	1.49 J	ND(1.0)	NA	NA	Gauged on 5/19/08
	10/28/2008	469.09	21.68	ND	ND	447.41	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/13/2009	469.09	21.08	ND	ND	448.01	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	469.09	18.47	ND	ND	450.62	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	469.09	18.40	ND	ND	450.69	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	469.09	16.75	ND	ND	452.34	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	5/5/2009	469.09	16.48	ND	ND	452.61	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	469.09	16.54	ND	ND	452.55	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	469.09	16.67	ND	ND	452.42	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	469.09	17.07	ND	ND	452.02	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	469.09	17.39	ND	ND	451.70	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	469.09	17.50	ND	ND	451.59	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	469.09	19.75	ND	ND	449.34	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/9/2009	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	469.09	16.27	ND	ND	452.82	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-53D	2/4/2010	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	469.09	16.05	ND	ND	453.04	NS	NS	NS	NS	NS	NS	NS	NS	Well inaccessible
	3/23/2010	469.09	15.47	ND	ND	453.62	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	469.09	11.36	ND	ND	457.73	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.25 J	NA	ND(200)	
	4/30/2010	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	469.09	18.77	ND	ND	450.32	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	7/7/2010	469.09	18.35	ND	ND	450.74	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	469.09	23.04	ND	ND	446.05	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	9/16/2010	469.09	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	469.09	20.92	ND	ND	448.17	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	1/13/2011	469.09	21.68	NM	NM	447.41	**27.4**	14.4	3.25	14.7	59.75	2.52	<100	**188**	
	5/9/2011	469.09	10.75	NM	NM	458.34	<1	1.03	<1	<3	1.03	<1	<94.3	<100	sheen
	8/2/2011	469.09	19.43	NM	NM	449.66	**14**	2.39	1.26	5.64	23.29	1.83	<98	<100	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-54S	12/14/2006	468.11	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/3/2007	468.11	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	468.11	11.34	ND	ND	456.77	**40.9**	9.9	8.4	282	341	16.8	NA	NA	
	5/21/2007	468.11	11.70	ND	ND	456.41	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	468.11	11.30	ND	ND	456.81	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	468.11	11.40	ND	ND	456.71	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2007	468.11	25.60	ND	ND	442.51	NS	NS	NS	NS	NS	NS	NS	NS	
	12/4/2007	468.11	26.77	ND	ND	441.34	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	468.11	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	468.11	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	468.11	9.05	ND	ND	459.06	**48.1**	4.7	33.2	134	220	9.4	NA	NA	Gauged on 5/19/08
	10/28/2008	468.11	18.55	ND	ND	449.56	0.31 J	ND(1.0)	ND(1.0)	ND(1.0)	0.31 J	0.38 J	NA	NA	
	1/13/2009	468.11	18.15	ND	ND	449.96	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	468.11	17.15	ND	ND	450.96	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	468.11	17.06	ND	ND	451.05	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	468.11	18.19	ND	ND	449.92	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	5/5/2009	468.11	18.05	ND	ND	450.06	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	468.11	17.94	ND	ND	450.17	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	468.11	18.03	ND	ND	450.08	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	468.11	16.85	ND	ND	451.26	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	468.11	17.00	ND	ND	451.11	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	468.11	17.12	ND	ND	450.99	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	468.11	18.17	ND	ND	449.94	NS	NS	NS	NS	NS	NS	NS	NS	
	12/9/2009	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	468.11	16.08	ND	ND	452.03	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-54S	3/23/2010	468.11	14.44	ND	ND	453.67	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	468.11	9.12	ND	ND	458.99	**66.4**	3.5	16.3	212	298	4.9	NA	**3,670**	
	4/30/2010	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	468.11	11.02	ND	ND	457.09	**33**	2.9	80.8	6.2	122.9	3.1	**989**	**2,340**	
	7/7/2010	468.11	11.13	ND	ND	456.98	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	468.11	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	468.11	18.10	ND	ND	450.01	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	468.11	18.15	ND	ND	449.96	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2010	468.11	18.20	ND	ND	449.91	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	468.11	18.15	ND	ND	449.96	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	468.11	18.20	ND	ND	449.91	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	468.11	18.15	ND	ND	449.96	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	468.11	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/21/2010	468.11	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/11/2011	468.11	18.21	ND	ND	449.90	NS	NS	NS	NS	NS	NS	NS	NS	
	5/9/2011	468.11	9.21	ND	ND	458.90	**7.54**	<1	1.99	<3	9.53	<1	**1,050**	**248**	sheen
	8/2/2011	468.11	18.18	ND	ND	449.93	NS	NS	NS	NS	NS	NS	NS	NS	

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Former Exxon Station #25553
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		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-54D	12/14/2006	467.84	23.69	ND	ND	444.15	**1,880**	789	**1,990**	6,720	11,379	**119**	NA	NA	
	1/3/2007	467.84	24.43	ND	ND	443.41	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	467.84	13.08	ND	ND	454.76	**1,920**	**1,190**	**1,650**	9,260	14,020	**93.8**	NA	NA	
	5/21/2007	467.84	13.11	ND	ND	454.73	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	467.84	12.97	ND	ND	454.87	NS	NS	NS	NS	NS	NS	NS	NS	
	8/3/2007	467.84	23.22	ND	ND	444.62	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	467.84	13.17	ND	ND	454.67	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	467.84	28.76	ND	ND	439.08	**1,160**	**1,070**	**1,430**	7,740	11,400	**107**	NA	NA	
	10/24/2007	467.84	26.80	ND	ND	441.04	NS	NS	NS	NS	NS	NS	NS	NS	
	12/4/2007	467.84	25.40	ND	ND	442.44	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	467.84	27.84	ND	ND	440.00	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	467.84	26.04	ND	ND	441.80	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	467.84	21.55	ND	ND	446.29	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	467.84	10.14	ND	ND	457.70	**469**	122	299	3,940	4,830	**133**	NA	NA	Gauged on 5/19/08
	10/28/2008	467.84	24.10	ND	ND	443.74	**165**	84.0	12.6	1,270	1,532	**86.6**	NA	NA	
	1/13/2009	467.84	23.85	ND	ND	443.99	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	467.84	20.63	ND	ND	447.21	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	467.84	20.45	ND	ND	447.39	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	467.84	19.15	ND	ND	448.69	**157**	23.2	55.7	482	718	**51.4**	NA	NA	
	5/5/2009	467.84	18.96	ND	ND	448.88	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	467.84	18.98	ND	ND	448.86	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	467.84	19.07	ND	ND	448.77	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	467.84	18.69	ND	ND	449.15	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	467.84	18.84	ND	ND	449.00	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	467.84	19.00	ND	ND	448.84	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	467.84	22.40	ND	ND	445.44	**198**	79.1	194	923	1,394	**40.9**	NA	NA	
	12/9/2009	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-54D	1/22/2010	467.84	18.44	ND	ND	449.40	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	467.84	16.63	ND	ND	451.21	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	467.84	11.93	ND	ND	455.91	**130**	39.2	177	665	1,011	**29.7**	NA	**7,750**	
	4/30/2010	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	467.84	20.69	ND	ND	447.15	3.3	1.5	0.85 J	7.7	13.4 J	1.2	**201**	**275**	
	7/7/2010	467.84	20.55	ND	ND	447.29	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	467.84	25.86	ND	ND	441.98	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	467.84	25.90	ND	ND	441.94	0.30 J	0.57 J	ND(1.0)	ND(1.0)	0.87 J	ND(1.0)	ND(100)	ND(200)	
	9/16/2010	467.84	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	467.84	18.20	ND	ND	449.64	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	467.84	25.41	ND	ND	442.43	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	467.84	25.20	ND	ND	442.64	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	467.84	23.09	ND	ND	444.75	0.90 J	ND(1.0)	ND(1.0)	ND(1.0)	0.90 J	ND(1.0)	ND(100)	ND(200)	
	12/21/2010	467.84	23.92	ND	ND	443.92	NS	NS	NS	NS	NS	NS	NS	NS	
	1/11/2011	467.84	24.49	ND	ND	443.35	**1,340**	169	**1,630**	1,220	4,359	**63.3**	**3,010**	**22,400**	
	5/9/2011	467.84	11.31	ND	ND	456.53	**35.3**	7.41	27.6	26.3	96.6	1.44	**152**	**413**	sheen
	8/2/2011	467.84	21.91	ND	ND	445.93	**1,160**	144	**1,330**	1,320	3,954	**49.7**	**3,280**	**28,500**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-55S	12/14/2006	465.76	19.50	ND	ND	446.26	**1,050**	**4,390**	679	6,840	12,959	**26.4**	NA	NA	
	1/3/2007	465.76	14.91	ND	ND	450.85	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	465.76	9.30	ND	ND	456.46	**2,390**	**1,610**	**1,750**	7,310	13,060	**124**	NA	NA	
	5/21/2007	465.76	9.33	ND	ND	456.43	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	465.76	9.26	ND	ND	456.50	NS	NS	NS	NS	NS	NS	NS	NS	
	8/3/2007	465.76	15.11	ND	ND	450.65	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	465.76	9.39	ND	ND	456.37	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	465.76	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2007	465.76	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	465.76	19.81	ND	ND	445.95	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	465.76	19.83	ND	ND	445.93	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	465.76	19.83	ND	ND	445.93	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	465.76	6.90	ND	ND	458.86	**1,110**	620	**1,210**	1,720	4,660	**70.4**	NA	NA	
	10/28/2008	465.76	19.80	ND	ND	445.96	**521**	356	532	757	2,166	**75.8**	NA	NA	
	1/13/2009	465.76	19.26	ND	ND	446.50	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	465.76	13.58	ND	ND	452.18	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	465.76	13.48	ND	ND	452.28	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	465.76	13.30	ND	ND	452.46	**262**	43.9	547	273	1,126	**53.3**	NA	NA	
	5/5/2009	465.76	13.05	ND	ND	452.71	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	465.76	13.37	ND	ND	452.39	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	465.76	13.84	ND	ND	451.92	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	465.76	12.76	ND	ND	453.00	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	465.76	12.97	ND	ND	452.79	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	465.76	14.60	ND	ND	451.16	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	465.76	14.85	ND	ND	450.91	**203**	35.0	**1,550**	88.6	1,877	**27.2**	NA	NA	
	12/9/2009	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	465.76	11.20	ND	ND	454.56	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-55S	2/4/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	465.76	11.68	ND	ND	454.08	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	465.76	7.85	ND	ND	457.91	**187**	51.8	226	103	568	**47.4**	NA	**7,400**	
	4/30/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	465.76	13.56	ND	ND	452.20	**346**	85.8	**2,390**	590	3,412	17.0 J	**3,290**	**1,540**	
	7/7/2010	465.76	13.20	ND	ND	452.56	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	465.76	19.42	ND	ND	446.34	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	465.76	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/11/2011	465.76	15.98	NM	NM	449.78	NS	NS	NS	NS	NS	NS	NS	NS	
	5/9/2011	465.76	6.68	NM	NM	459.08	**588**	79.6	434	242	1,344	5.62	**4,230**	**5,430**	sheen
	8/2/2011	465.76	14.76	NM	NM	451.00	**470**	29.4	188	18.3	705.7	11.2	**2,280**	**4,360**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-55I	12/14/2006	465.67	21.40	ND	ND	444.27	**2,280**	**3,700**	**1,600**	6,870	14,450	**31.9**	NA	NA	
	1/3/2007	465.67	19.06	ND	ND	446.61	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	465.67	10.90	ND	ND	454.77	**10,700**	**12,300**	**820**	10,000	33,820	**32.6 J**	NA	NA	
	5/21/2007	465.67	10.96	ND	ND	454.71	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	465.67	10.86	ND	ND	454.81	NS	NS	NS	NS	NS	NS	NS	NS	
	8/3/2007	465.67	21.88	ND	ND	443.79	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	465.67	11.01	ND	ND	454.66	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	465.67	26.35	ND	ND	439.32	**13,700**	**15,400**	**1,610**	8,390	39,100	ND(100)	NA	NA	
	10/24/2007	465.67	26.41	ND	ND	439.26	NS	NS	NS	NS	NS	NS	NS	NS	
	12/4/2007	465.67	26.77	ND	ND	438.90	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	465.67	25.65	ND	ND	440.02	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	465.67	23.11	ND	ND	442.56	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	465.67	19.22	ND	ND	446.45	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	465.67	8.05	ND	ND	457.62	**7,840**	**14,300**	**3,790**	**17,100**	43,030	ND(100)	NA	NA	
	10/28/2008	465.67	22.08	ND	ND	443.59	**5,600**	**5,920**	**3,150**	**14,400**	29,070	**33.3**	NA	NA	
	1/13/2009	465.67	21.40	ND	ND	444.27	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	465.67	17.52	ND	ND	448.15	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	465.67	17.40	ND	ND	448.27	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	465.67	17.11	ND	ND	448.56	**6,580**	**6,870**	**1,050**	7,510	22,010	**24.3 J**	NA	NA	
	5/5/2009	465.67	16.89	ND	ND	448.78	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	465.67	16.88	ND	ND	448.79	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	465.67	17.01	ND	ND	448.66	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	465.67	16.45	ND	ND	449.22	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	465.67	16.70	ND	ND	448.97	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	465.67	19.13	ND	ND	446.54	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	465.67	20.12	ND	ND	445.55	**6,720**	**7,650**	**1,270**	7,430	23,070	15.9 J	NA	NA	
	12/9/2009	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-55I	1/22/2010	465.67	15.32	ND	ND	450.35	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	465.67	9.56	ND	ND	456.11	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	465.67	9.74	ND	ND	455.93	**948**	557	37.7	1,310	2,853	4.4	NA	**13,300**	
	4/30/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	465.67	18.41	ND	ND	447.26	**4,270**	**3,420**	172	3,450	11,312	**25.7**	**3,650**	**43,100**	
	7/7/2010	465.67	15.32	ND	ND	450.35	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	465.67	23.51	ND	ND	442.16	**50.2**	67.2	8.7	351	477	ND(5.0)	**17,300**	**7,850**	
	9/16/2010	465.67	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	465.67	23.61	ND	ND	444.06	**2,730**	742	27.1	153	3,652	19.3	**1,640**	**9,170**	
	1/12/2011	465.67	22.23	NM	NM	443.44	**3,990**	**2,980**	**1,740**	7,360	16,070	**25.4**	**42,300**	**46,900**	
	5/9/2011	465.67	16.83	NM	NM	448.84	**7,220**	**3,750**	**1,300**	6,080	18,350	**37**	**4,190**	**54,600**	sheen
	8/2/2011	465.67	19.63	NM	NM	446.04	**3,360**	**1,540**	**1,130**	3,780	9,810	19.2	**2,470**	**49,000**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MW-55D	12/14/2006	466.24	20.21	ND	ND	446.03	**181**	170	69.1	281	701	9.6	NA	NA	
	1/3/2007	466.24	18.23	ND	ND	448.01	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	466.24	10.50	ND	ND	455.74	**105**	82.6	52.5	123	363	10.5	NA	NA	
	5/21/2007	466.24	10.55	ND	ND	455.69	NS	NS	NS	NS	NS	NS	NS	NS	
	7/2/2007	466.24	10.45	ND	ND	455.79	NS	NS	NS	NS	NS	NS	NS	NS	
	8/3/2007	466.24	18.85	ND	ND	447.39	NS	NS	NS	NS	NS	NS	NS	NS	
	9/7/2007	466.24	10.62	ND	ND	455.62	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	466.24	23.80	ND	ND	442.44	**40.5**	31.1	17.8	53.7	143.1	7.6	NA	NA	
	10/24/2007	466.24	23.92	ND	ND	442.32	NS	NS	NS	NS	NS	NS	NS	NS	
	12/4/2007	466.24	23.78	ND	ND	442.46	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	466.24	23.30	ND	ND	442.94	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	466.24	21.10	ND	ND	445.14	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	466.24	18.46	ND	ND	447.78	NS	NS	NS	NS	NS	NS	NS	NS	
	5/21/2008	466.24	8.28	ND	ND	457.96	**50.0**	26.4	23.7	49.5	149.6	8.6	NA	NA	Gauged on 5/19/08
	10/28/2008	466.24	19.80	ND	ND	446.44	**88.9**	21.6	19.9	48.3	178.7	8.0	NA	NA	
	1/13/2009	466.24	19.20	ND	ND	447.04	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	466.24	16.00	ND	ND	450.24	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	466.24	15.84	ND	ND	450.40	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	466.24	15.62	ND	ND	450.62	**61.3**	9.5	6.7	32.9	110.4	7.1	NA	NA	
	5/5/2009	466.24	15.47	ND	ND	450.77	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	466.24	15.30	ND	ND	450.94	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	466.24	15.89	ND	ND	450.35	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	466.24	14.61	ND	ND	451.63	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	466.24	14.84	ND	ND	451.40	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	466.24	17.00	ND	ND	449.24	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	466.24	17.53	ND	ND	448.71	**42.7**	6.2	7.3	18.0	74.2	5.2	NA	NA	
	12/9/2009	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	

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Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MW-55D	1/22/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	466.24	14.93	ND	ND	451.31	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	466.24	8.61	ND	ND	457.63	**48.1**	7.8	21.5	19.9	97.3	7.3	NA	**812**	
	4/30/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	466.24	15.90	ND	ND	450.34	**78.2**	10.9	19.0	28.7	136.8	7.6	**290**	**1,260**	
	7/7/2010	466.24	18.01	ND	ND	448.23	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	466.24	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	466.24	21.00	ND	ND	445.24	**57.6**	7.8	13.2	21.9	100.5	5.5	**202**	**1,060**	
	9/16/2010	466.24	26.51	ND	ND	439.73	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	466.24	19.17	ND	ND	447.07	**67.4**	6.9	8.7	12.0	95.0	6.5	**121**	**535**	
	1/11/2011	466.24	19.24	ND	ND	447.00	**150**	9.29	16.3	9.45	185.04	6.6	**192**	**1,820**	
	5/9/2011	466.24	7.58	ND	ND	458.66	**92.4**	5.03	5.35	12.8	115.58	6.51	**168**	**589**	sheen
	8/2/2011	466.24	17.09	ND	ND	449.15	**142**	6.54	14.7	6.77	170.01	5.23	**134**	**1,520**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
RW-52	4/12/2005	459.78	18.52	ND	ND	441.26	**193**	19.6	61.9	71.8	346	**144**	NA	NA	
	7/7/2005	459.78	21.70	ND	ND	438.08	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	459.78	30.45	ND	ND	429.33	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	459.78	33.84	ND	ND	425.94	**127**	18.2	55.8	79.0	280	**37.8**	NA	NA	
	12/20/2005	459.78	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	460.36	35.95	ND	ND	424.41	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	460.36	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	460.36	NM	NM	NM	NM	**169**	31.8	80.2	116	397	**42.0**	NA	NA	
	5/31/2006	460.36	28.65	ND	ND	431.71	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	460.94	27.32	ND	ND	433.62	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	460.94	27.13	ND	ND	433.81	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	460.94	26.64	ND	ND	434.30	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	460.94	25.70	ND	ND	435.24	**183**	10.7	26.2	36.4	256	**35.3**	NA	NA	
	11/22/2006	460.94	21.84	ND	ND	439.10	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	460.94	25.63	ND	ND	435.31	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	460.94	25.40	ND	ND	435.54	**275**	13.0	10.4	18.8	317	**30.1**	NA	NA	
	12/19/2007	460.94	20.66	ND	ND	440.28	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	460.94	NM	NM	NM	NM	**254**	7.1	2.6	13.5	277	**28.7**	NA	NA	
	10/28/2008	460.94	NM	NM	NM	NM	**169**	7.7	3.7	17.8	198	**23.3**	NA	NA	
	1/13/2009	460.94	14.00	ND	ND	446.94	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	460.94	14.00	ND	ND	446.94	**131**	5.9	2.2	15.0	154	18.3	NA	NA	
	5/5/2009	460.94	14.00	ND	ND	446.94	NS	NS	NS	NS	NS	NS	NS	NS	
	6/30/2009	460.94	6.15	ND	ND	454.79	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	460.94	14.36	ND	ND	446.58	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
RW-52	8/10/2009	460.94	11.60	ND	ND	449.34	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	460.94	NM	NM	NM	NM	**21.9**	0.81 J	0.57 J	2.3	25.6 J	0.44 J	NA	NA	
	12/9/2009	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	460.94	12.10	ND	ND	448.84	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	460.94	4.63	ND	ND	456.31	**85.9**	5.1	0.42 J	1.4	92.8 J	8.1	NA	**1,120**	
	4/30/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	460.94	14.08	ND	ND	446.86	ND(1.0)	51.4	ND(1.0)	ND(1.0)	51.4	ND(1.0)	**34,800**	**254**	
	7/7/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	460.94	19.61	ND	ND	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	460.94	19.74	ND	ND	441.20	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	9/16/2010	460.94	20.21	ND	ND	440.73	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	460.94	19.85	ND	ND	441.09	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	460.94	19.51	ND	ND	441.43	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	460.94	18.95	ND	ND	441.99	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	460.94	18.45	ND	ND	442.49	0.83 J	12.2	0.47 J	ND(1.0)	13.5 J	ND(1.0)	**22,100**	**266**	
	12/21/2010	460.94	19.20	ND	ND	441.74	NS	NS	NS	NS	NS	NS	NS	NS	
	1/11/2011	460.94	18.12	ND	ND	442.82	**163 [173]**	6.56 [7.61]	3.21 [3.75]	12.8 [14.5]	185.57 [198.86]	12.2 [13.9]	**361 [308]**	**3,270 [3,240]**	
	5/9/2011	460.94	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/2/2011	460.94	15.14	ND	ND	445.80	**175**	9.1	5.07	9.67	198.84	8.03	**270**	**2,820**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
VW-1	4/12/2005	465.07	7.80	ND	ND	457.27	**637**	**1,040**	157	718	2,552	13.0	NA	NA	
	5/3/2005	465.07	4.08	ND	ND	460.99	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	465.07	9.55	ND	ND	455.52	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	465.07	10.03	ND	ND	455.04	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	465.07	11.41	ND	ND	453.66	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	465.07	10.02	ND	ND	455.05	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/20/2005	465.07	11.16	ND	ND	453.91	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	465.07	10.53	ND	ND	454.54	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	465.07	10.34	ND	ND	454.73	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	465.07	10.65	ND	ND	454.42	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	465.07	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	465.07	11.31	ND	ND	453.76	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	465.07	9.66	ND	ND	455.41	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	465.07	11.26	ND	ND	453.81	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	465.07	11.52	ND	ND	453.55	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	465.07	10.07	ND	ND	455.00	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	465.07	10.75	ND	ND	454.32	**43.3**	1.8	3.3	3.5	51.9	0.46 J	NA	NA	
	11/22/2006	465.07	10.53	ND	ND	454.54	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	465.07	10.20	ND	ND	454.87	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	465.07	10.71	ND	ND	454.36	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	465.07	7.28	ND	ND	457.79	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	465.07	7.25	ND	ND	457.82	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	465.07	10.63	ND	ND	454.44	NS	NS	NS	NS	NS	NS	NS	NS	
	4/25/2007	465.07	7.30	ND	ND	457.77	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	465.07	7.30	ND	ND	457.77	**95.2**	173	35.1	225	528	12.8	NA	NA	
	7/19/2007	465.07	9.21	ND	ND	455.86	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	465.07	11.90	ND	ND	453.17	**465**	97.3	94.2	189	846	33.6	NA	NA	
	10/24/2007	465.07	12.60	ND	ND	452.47	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
VW-1	12/19/2007	465.07	10.13	ND	ND	454.94	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	465.07	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	465.07	8.41	ND	ND	456.66	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	465.07	4.00	ND	ND	461.07	**637**	49.9	330	573	1,590	**28.3**	NA	NA	
	10/28/2008	465.07	6.00	ND	ND	459.07	0.71 J	0.40 J	0.28 J	0.69 J	2.08 J	0.38 J	NA	NA	
	1/13/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	NSVD	6.16	ND	ND	NSVD	ND(1.0)	ND(1.0)	ND(1.0)	1.6	1.6	ND(1.0)	NA	NA	
	5/5/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	NSVD	9.76	ND	ND	NSVD	**224**	140	109	757	1,230	ND(1.0)	NA	NA	
	12/9/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	NSVD	5.82	ND	ND	NSVD	**52.9**	10.6	7.9	73.2	144.6	0.65 J	NA	**631**	
	4/30/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	NSVD	9.00	ND	ND	NSVD	**1,150**	397	**834**	2,440	4,821	**88.0**	**2,010**	**20,600**	
	9/2/2010	NSVD	9.86	ND	ND	NSVD	**599**	51	254	389	1,293	**47.7**	**1,040**	**6,380**	
	12/8/2010	NSVD	6.80	ND	ND	NSVD	**567**	29	44	56	696	**51.6**	**241**	**2,040**	
	1/13/2011	NSVD	7.74	ND	ND	NSVD	**1,030**	375	**956**	1,570	3,931	**40.8**	**2,310**	**12,600**	
	5/9/2011	NSVD	4.65	ND	ND	NSVD	**623**	80.1	610	718	2,031	16.5	**1,630**	**5,160**	sheen
	8/2/2011	NSVD	9.31	ND	ND	NSVD	**2,020**	199	**1,490**	2,620	6,329	**86**	**1,930**	**30,600**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
VW-2	4/12/2005	465.88	3.05	ND	ND	462.83	ND(5.0)	ND(5.0)	ND(5.0)	2.5 J	2.5 J	**5,760**	NA	NA	
	5/3/2005	465.88	3.10	ND	ND	462.78	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	465.88	3.42	ND	ND	462.46	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	465.88	3.67	ND	ND	462.21	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	465.88	6.00	ND	ND	459.88	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	465.88	3.01	ND	ND	462.87	ND(1.0)	3.7	0.59 J	3.1	7.4 J	4.1	NA	NA	
	12/20/2005	465.88	3.12	ND	ND	462.76	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	465.88	3.10	ND	ND	462.78	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	465.88	3.10	ND	ND	462.78	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	465.88	3.35	ND	ND	462.53	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	465.88	3.57	ND	ND	462.31	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	6.5	NA	NA	
	5/31/2006	465.88	3.53	ND	ND	462.35	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	465.88	3.30	ND	ND	462.58	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	465.88	11.40	ND	ND	454.48	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	465.88	3.48	ND	ND	462.40	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	465.88	3.29	ND	ND	462.59	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	465.88	3.35	ND	ND	462.53	ND(1.0)	ND(1.0)	ND(1.0)	0.35 J	0.35 J	4.8	NA	NA	
	11/22/2006	465.88	3.08	ND	ND	462.80	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	465.88	3.05	ND	ND	462.83	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	465.88	2.65	ND	ND	463.23	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	7/19/2007	465.88	3.30	ND	ND	462.58	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2007	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	465.88	3.12	ND	ND	462.76	NS	NS	NS	NS	NS	NS	NS	NS	
	1/21/2008	465.88	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	465.88	2.92	ND	ND	462.96	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
VW-2	5/19/2008	465.88	3.27	ND	ND	462.61	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	1.4	NA	NA	
	10/28/2008	465.88	6.80	ND	ND	459.08	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	3.4	NA	NA	
	1/13/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	NSVD	2.87	ND	ND	NSVD	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.96 J	NA	NA	
	5/5/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	NSVD	3.67	ND	ND	NSVD	2.1	3.8	ND(1.0)	1.7	7.6	2.1	NA	NA	
	12/9/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	NSVD	2.69	ND	ND	NSVD	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	4.0	NA	ND(200)	
	4/30/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	NSVD	3.25	ND	ND	NSVD	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.55 J	**272**	**244**	
	9/2/2010	NSVD	6.78	ND	ND	NSVD	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	2.3	ND(100)	ND(200)	
	12/8/2010	NSVD	3.02	ND	ND	NSVD	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.89 J	**147**	ND(200)	
	1/13/2011	NSVD	3.50	ND	ND	NSVD	**10.7**	29.4	7.79	32.7	80.59	1.0	**128**	**226**	
	5/9/2011	NSVD	2.85	ND	ND	NSVD	<1	<1	<1	<3	BRL	<1	**146**	**114**	
	8/2/2011	NSVD	9.07	ND	ND	NSVD	1.05	2.29	<1	<3	3.34	1.5	**669**	**227**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
VW-3	4/12/2005	464.65	7.10	ND	ND	457.55	2.7	5.9	1.5	6.2	16.3	**30.3**	NA	NA	
	5/3/2005	464.65	5.89	ND	ND	458.76	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	464.65	8.55	ND	ND	456.10	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	464.65	10.85	ND	ND	453.80	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	464.65	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	464.65	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/20/2005	464.65	10.40	ND	ND	454.25	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	464.65	10.49	ND	ND	454.16	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	464.65	10.31	ND	ND	454.34	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	464.65	10.63	ND	ND	454.02	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	464.65	11.03	ND	ND	453.62	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	0.64 J	NA	NA	
	5/31/2006	464.65	11.30	ND	ND	453.35	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	464.65	10.35	ND	ND	454.30	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	464.65	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	464.65	11.00	ND	ND	453.65	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	464.65	10.77	ND	ND	453.88	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	464.65	10.77	ND	ND	453.88	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	11/22/2006	464.65	10.23	ND	ND	454.42	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	464.65	10.35	ND	ND	454.30	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	464.65	7.15	ND	ND	457.50	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	464.65	7.04	ND	ND	457.61	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	464.65	10.82	ND	ND	453.83	NS	NS	NS	NS	NS	NS	NS	NS	
	4/25/2007	464.65	6.70	ND	ND	457.95	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	464.65	6.70	ND	ND	457.95	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	7/19/2007	464.65	8.58	ND	ND	456.07	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	464.65	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2007	464.65	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	464.65	10.57	ND	ND	454.08	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
VW-3	1/21/2008	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/19/2008	464.65	7.98	ND	ND	456.67	NS	NS	NS	NS	NS	NS	NS	NS	
	5/19/2008	464.65	2.52	ND	ND	462.13	**88.3**	4.4	2.5	1.1	96.3	**20.9**	NA	NA	
	10/28/2008	464.65	6.83	ND	ND	457.82	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	1/13/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/6/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/16/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	4/23/2009	464.65	5.12	ND	ND	459.53	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	5/5/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/1/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	7/9/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	8/10/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	9/3/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/1/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	10/12/2009	464.65	8.89	ND	ND	455.76	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	12/9/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	464.65	5.23	ND	ND	459.42	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	ND(200)	
	4/30/2010	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	464.65	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	464.65	8.08	ND	ND	456.57	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	**269**	ND(200)	
	9/2/2010	464.65	9.05	ND	ND	455.60	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	**144**	ND(200)	
	12/8/2010	464.65	5.33	ND	ND	459.32	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	ND(100)	ND(200)	
	1/12/2011	464.65	6.71	ND	ND	457.94	<1	2.31	2.03	9.37	13.71	<1	**129**	<100	
	5/9/2011	464.65	4.31	ND	ND	460.34	<1	<1	<1	<3	BRL	<1	**1,050**	<100	
	8/2/2011	464.65	8.40	ND	ND	456.25	<1	<1	<1	<3	BRL	<1	<98	<100	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
VW-4	4/12/2005	465.88	8.06	ND	ND	457.82	2.3	ND(1.0)	ND(1.0)	ND(1.0)	2.3	3.9	NA	NA	
	5/3/2005	465.88	7.96	ND	ND	457.92	NS	NS	NS	NS	NS	NS	NS	NS	
	6/21/2005	465.88	9.69	ND	ND	456.19	NS	NS	NS	NS	NS	NS	NS	NS	
	7/7/2005	465.88	11.85	ND	ND	454.03	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2005	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/20/2005	465.88	11.62	ND	ND	454.26	NS	NS	NS	NS	NS	NS	NS	NS	
	1/9/2006	465.88	12.11	ND	ND	453.77	NS	NS	NS	NS	NS	NS	NS	NS	
	2/10/2006	465.88	11.96	ND	ND	453.92	NS	NS	NS	NS	NS	NS	NS	NS	
	3/24/2006	465.88	12.53	ND	ND	453.35	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	465.88	12.90	ND	ND	452.98	NS	NS	NS	NS	NS	NS	NS	NS	
	6/27/2006	465.88	12.15	ND	ND	453.73	NS	NS	NS	NS	NS	NS	NS	NS	
	8/24/2006	465.88	12.70	ND	ND	453.18	NS	NS	NS	NS	NS	NS	NS	NS	
	9/28/2006	465.88	12.43	ND	ND	453.45	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2006	465.88	12.05	ND	ND	453.83	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2006	465.88	11.91	ND	ND	453.97	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	465.88	12.05	ND	ND	453.83	NS	NS	NS	NS	NS	NS	NS	NS	
	12/22/2006	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/19/2007	465.88	8.01	ND	ND	457.87	NS	NS	NS	NS	NS	NS	NS	NS	
	2/23/2007	465.88	7.97	ND	ND	457.91	NS	NS	NS	NS	NS	NS	NS	NS	
	3/22/2007	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	7/19/2007	465.88	9.51	ND	ND	456.37	NS	NS	NS	NS	NS	NS	NS	NS	
	10/11/2007	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2007	465.88	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/19/2007	465.88	10.57	ND	ND	455.31	NS	NS	NS	NS	NS	NS	NS	NS	Well destroyed

Table 1
Groundwater Monitoring and Analytical Data
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		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
VW-5	4/12/2005	466.31	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well destroyed
VW-6	4/12/2005	466.53	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/26/2005	466.53	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	466.53	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	466.53	9.27	ND	ND	457.26	**22.3**	8.2	3.2	17.2	50.9	ND(1.0)	NA	NA	
	12/14/2006	466.53	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	466.53	7.50	ND	ND	459.03	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/11/2007	466.53	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	10/24/2007	466.53	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well destroyed
VW-7	4/12/2005	466.53	9.39	ND	ND	457.17	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/26/2005	466.56	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/20/2006	466.56	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	5/31/2006	466.56	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	11/7/2006	466.56	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	12/14/2006	466.56	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	4/26/2007	466.56	8.35	ND	ND	458.21	ND(1.0)	ND(1.0)	ND(1.0)	ND(1.0)	BRL	ND(1.0)	NA	NA	
	10/11/2007	466.56	9.76	ND	ND	456.80	**165**	10.1	54.4	134	364	ND(1.0)	NA	NA	
	12/19/2007	466.56	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well destroyed

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
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Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
DPE-1	11/11/2009	NSVD	13.43	ND	ND	NSVD	NS	NS	NS	NS	NS	NS	NS	NS	
	11/30/2009	NSVD	14.29	ND	ND	NSVD	**4,190**	**15,500**	**1,170**	8,150	29,010	**241**	NA	NA	
	12/9/2009	468.37	13.15	ND	ND	455.22	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	468.37	10.10	ND	ND	458.27	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	468.37	14	ND	ND	454.37	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	468.37	8.6	ND	ND	459.77	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	468.37	8.52	ND	ND	459.85	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	468.37	6.80	ND	ND	461.57	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	468.37	8.04	ND	ND	460.33	**33.4**	340	307	699	1,379	3.3	NA	**5,970**	
	4/30/2010	468.37	14.00	ND	ND	454.37	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	468.37	13.7	ND	ND	454.67	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	468.37	17.70	ND	ND	450.67	**2,360**	**14,300**	**5,920**	**26,900**	49,480	**490**	**31,500**	**135,000**	
	7/7/2010	468.37	17.22	ND	ND	451.15	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	468.37	22.00	ND	ND	446.37	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	468.37	22.51	22.5	0.01	445.87	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	468.37	21.79	ND	ND	446.58	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	468.37	22.11	ND	ND	446.26	**6,950**	**16,200**	**1,940**	9,730	34,820	**405**	**7,950**	**104,000**	
	9/16/2010	468.37	22.70	ND	ND	445.67	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	468.37	22.52	ND	ND	445.85	NS	NS	NS	NS	NS	NS	NS	NS	
	10/20/2010	468.37	19.61	ND	ND	448.76	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	468.37	22.78	ND	ND	445.59	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	468.37	22.02	ND	ND	446.35	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	468.37	18.13	ND	ND	450.24	**6,890**	**16,500**	**1,120**	5,290	26,800	**478**	**3,320**	**70,000**	
	12/21/2010	468.37	18.23	ND	ND	450.14	NS	NS	NS	NS	NS	NS	NS	NS	
	1/13/2011	468.37	18.61	ND	ND	449.76	**6,050**	**15,600**	**2,060**	**11,000**	34,710	**372**	**4,770**	**78,000**	
	5/9/2011	468.37	9.41	ND	ND	458.96	**4,240**	**11,000**	**1,970**	8,720	25,930	**216**	**3,780**	**67,200**	
	8/2/2011	468.37	18.14	ND	ND	450.23	**5,630**	**14,300**	**1,830**	9,960	31,720	**263**	**2,880**	**115,000**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L) [5]**	**Toluene (µg/L) [1,000]**	**Ethyl-benzene (µg/L) [700]**	**Total Xylenes (µg/L) [10,000]**	**Total BTEX (µg/L) --**	**MTBE (µg/L) [20]**	**TPH-DRO (µg/L) [47]**	**TPH-GRO (µg/L) [47]**	**Comments**
MP-1	11/30/2009	NSVD	6.50	ND	ND	NSVD	**35.7**	210	140	452	838	2.8	NA	NA	
	12/9/2009	468.51	7.22	ND	ND	461.29	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	468.51	7.74	ND	ND	460.77	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	468.51	6.27	ND	ND	462.24	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	468.51	9.35	ND	ND	459.16	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	468.51	8.87	ND	ND	459.64	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	468.51	8.40	ND	ND	460.11	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	468.51	9.65	ND	ND	458.86	**1,050**	**1,960**	436	4,050	7,496	**82.6**	NA	**23,700**	
	4/30/2010	468.51	16.29	ND	ND	452.22	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	468.51	14.30	ND	ND	454.21	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	468.51	15.58	ND	ND	452.93	**2,950**	**5,190**	628	3,470	12,238	**327**	**2,380**	**50,000**	
	7/7/2010	468.51	15.40	ND	ND	453.11	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	468.51	22.85	ND	ND	445.66	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	468.51	15.22	ND	ND	453.29	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	468.51	20.05	ND	ND	448.46	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	468.51	20.75	ND	ND	447.76	**684**	231	168	576	1,659	**228**	**3,080**	**15,600**	
	9/16/2010	468.51	21.65	ND	ND	446.86	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	468.51	21.47	ND	ND	447.04	NS	NS	NS	NS	NS	NS	NS	NS	
	10/20/2010	468.51	23.01	ND	ND	445.50	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	468.51	21.41	ND	ND	447.10	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	468.51	18.44	ND	ND	450.07	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	468.51	14.93	ND	ND	453.58	**3,780**	**7,830**	491	3,400	15,501	**405**	**4,270**	**42,600**	
	12/21/2010	468.51	17.31	ND	ND	451.20	NS	NS	NS	NS	NS	NS	NS	NS	
	1/13/2011	468.51	13.66	ND	ND	454.85	**5,940**	**11,200**	**929**	5,790	23,859	**443**	**3,080**	**57,300**	
	5/9/2011	468.51	7.21	ND	ND	461.30	**4,240**	**7,550**	**1,120**	4,960	17,870	**188**	**2,190**	**39,100**	
	8/2/2011	468.51	15.10	ND	ND	453.41	**7,580**	**13,400**	**1,490**	7,390	29,860	**329**	**1,710**	**101,000**	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MP-2	11/11/2009	NSVD	19.50	18.30	1.20	NSVD	NS	NS	NS	NS	NS	NS	NS	NS	
	11/19/2009	NSVD	18.10	16.80	1.30	NSVD	NS	NS	NS	NS	NS	NS	NS	NS	
	11/30/2009	NSVD	16.74	ND	ND	NSVD	**4,040**	**18,000**	**3,010**	**15,200**	40,250	**535**	NA	NA	
	12/9/2009	468.20	14.20	ND	ND	454.00	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	468.20	12.08	ND	ND	456.12	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	468.20	16.00	ND	ND	452.20	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	468.20	12.15	ND	ND	456.05	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	468.20	11.85	ND	ND	456.35	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	468.20	9.50	ND	ND	458.70	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	468.20	11.25	ND	ND	456.95	**434**	**2,200**	**2,140**	9,150	13,924	**116**	**NA**	**45,300**	
	4/30/2010	468.20	13.15	ND	ND	455.05	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	468.20	16.90	ND	ND	451.30	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	468.20	18.23	ND	ND	449.97	**4,810**	**16,700**	**2,220**	10,000	33,730	**232**	**2,850**	**116,000**	
	7/7/2010	468.20	17.93	ND	ND	450.27	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	468.20	23.00	21.85	1.15	446.07	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	468.20	23.72	22.60	1.12	445.33	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	468.20	23.10	21.98	1.12	445.95	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	468.20	22.46	22.30	0.16	445.86	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2010	468.20	22.99	22.80	0.19	445.35	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	468.20	22.04	22.01	0.03	446.18	NS	NS	NS	NS	NS	NS	NS	NS	
	10/20/2010	468.20	21.73	21.70	0.03	446.49	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	468.20	22.07	21.97	0.10	446.21	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	468.20	21.43	21.41	0.02	446.79	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	468.20	20.41	20.36	0.05	447.83	NS	NS	NS	NS	NS	NS	NS	NS	
	12/21/2010	468.20	20.57	20.54	0.03	447.65	NS	NS	NS	NS	NS	NS	NS	NS	
	1/13/2011	468.20	20.89	ND	ND	447.31	**2,520**	**9,240**	**2,200**	**17,000**	30,960	**425**	**20,200**	**69,200**	
	5/9/2011	468.20	10.03	ND	ND	458.17	**887**	**3,650**	**1,830**	8,900	15,267	**159**	**9,210**	**46,800**	
	8/2/2011	468.20	18.76	ND	ND	449.44	NS	NS	NS	NS	NS	NS	NS	NS	

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

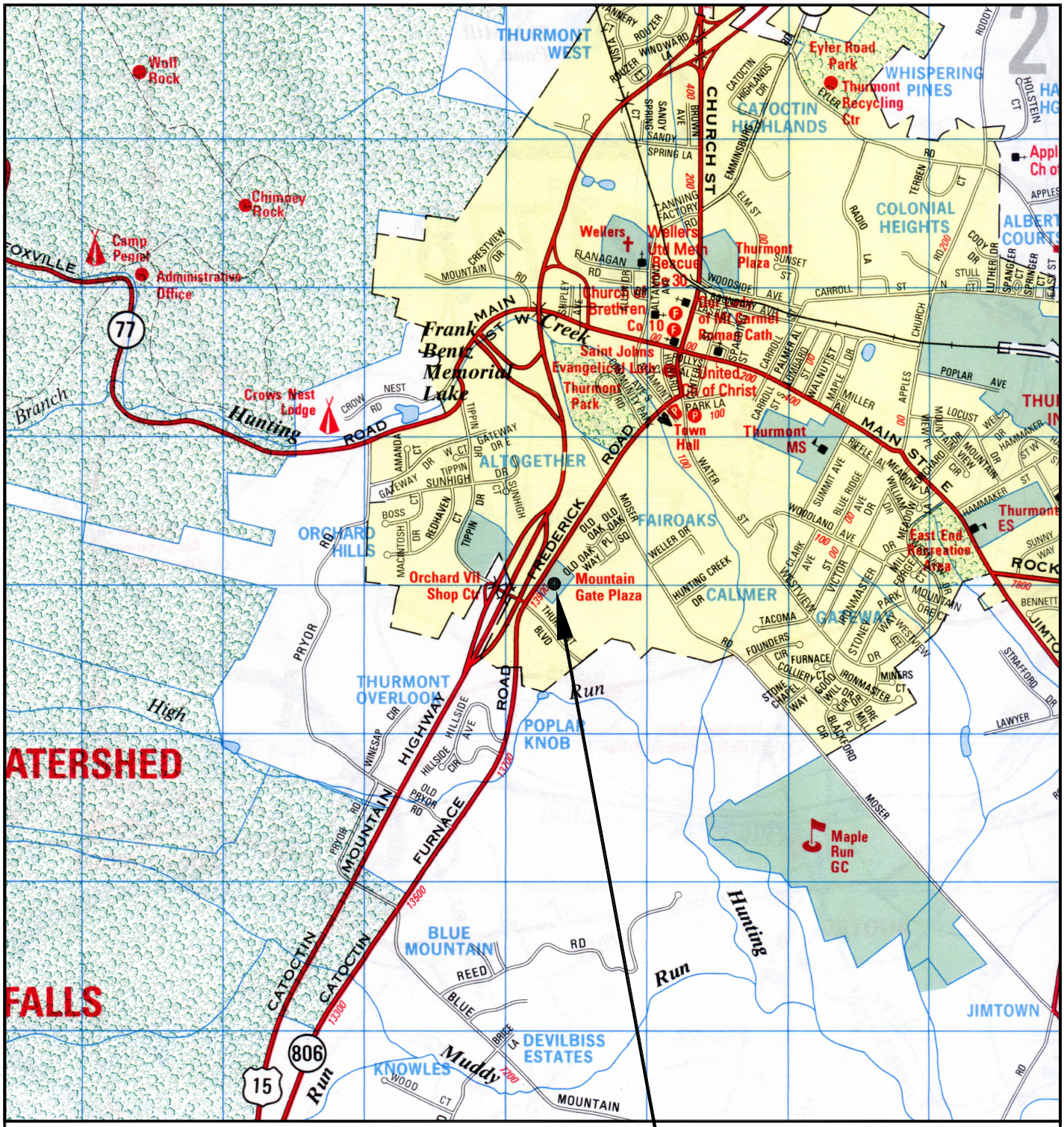
		Gauging Data					Analytical Data								
Well ID	**Date**	**Top of Casing Elevation (feet)**	**Depth to Water (feet)**	**Depth to Hydro-carbon (feet)**	**Hydro-carbon Thickness (feet)**	**Corrected GW Elevation (feet)**	**Benzene (µg/L)** [5]	**Toluene (µg/L)** [1,000]	**Ethyl-benzene (µg/L)** [700]	**Total Xylenes (µg/L)** [10,000]	**Total BTEX (µg/L)** --	**MTBE (µg/L)** [20]	**TPH-DRO (µg/L)** [47]	**TPH-GRO (µg/L)** [47]	**Comments**
MP-3	11/30/2009	NSVD	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	Well inaccessible
	12/9/2009	468.62	16.75	ND	ND	451.87	NS	NS	NS	NS	NS	NS	NS	NS	
	12/30/2009	468.62	DRY	DRY	DRY	DRY	NS	NS	NS	NS	NS	NS	NS	NS	
	1/22/2010	468.62	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	2/4/2010	468.62	13.00	ND	ND	455.62	NS	NS	NS	NS	NS	NS	NS	NS	
	2/16/2010	468.62	12.00	ND	ND	456.62	NS	NS	NS	NS	NS	NS	NS	NS	
	3/23/2010	468.62	10.00	ND	ND	458.62	NS	NS	NS	NS	NS	NS	NS	NS	
	3/31/2010	468.62	11.35	ND	ND	457.27	**435**	223	350	975	1,983	**64.5**	NA	**8,580**	
	4/30/2010	468.62	14.36	ND	ND	454.26	NS	NS	NS	NS	NS	NS	NS	NS	
	5/16/2010	468.62	14.80	ND	ND	453.82	NS	NS	NS	NS	NS	NS	NS	NS	
	6/17/2010	468.62	6.00	ND	ND	462.62	**489**	221	472	549	1,731	**69.6**	**1,110**	**11,400**	
	7/7/2010	468.62	6.89	ND	ND	461.73	NS	NS	NS	NS	NS	NS	NS	NS	
	7/30/2010	468.62	22.10	ND	ND	446.52	NS	NS	NS	NS	NS	NS	NS	NS	
	8/13/2010	468.62	22.80	ND	ND	445.82	NS	NS	NS	NS	NS	NS	NS	NS	
	8/27/2010	468.62	22.80	ND	ND	445.82	NS	NS	NS	NS	NS	NS	NS	NS	
	9/2/2010	468.62	22.95	ND	ND	445.67	NS	NS	NS	NS	NS	NS	NS	NS	
	9/16/2010	468.62	22.98	ND	ND	445.64	NS	NS	NS	NS	NS	NS	NS	NS	
	10/8/2010	468.62	22.23	ND	ND	446.39	NS	NS	NS	NS	NS	NS	NS	NS	
	10/20/2010	468.62	21.23	ND	ND	447.39	NS	NS	NS	NS	NS	NS	NS	NS	
	11/10/2010	468.62	19.20	ND	ND	449.42	NS	NS	NS	NS	NS	NS	NS	NS	
	11/22/2010	468.62	23.10	ND	ND	445.52	NS	NS	NS	NS	NS	NS	NS	NS	
	12/8/2010	468.62	23.81	ND	ND	444.81	**454**	323	**749**	1,800	3,326	**72.7**	NA	**26,000**	
	12/21/2010	468.62	23.75	ND	ND	444.87	NS	NS	NS	NS	NS	NS	NS	NS	
	1/13/2011	468.62	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	
	5/9/2011	468.62	10.46	NM	NM	458.16	**543**	427	**1,200**	4,710	6,880	**47.9**	**3,170**	**22,400**	
	8/2/2011	468.62	NM	NM	NM	NM	NS	NS	NS	NS	NS	NS	NS	NS	obstruction

Table 1
Groundwater Monitoring and Analytical Data
Former Exxon Station #25553
143 Frederick Road
Thurmont, Maryland

Notes:

µg/L - micrograms/liter
BRL - Below laboratory reporting limits
BTEX - Benzene, toluene, ethyl benzene, and total xylenes
DRO - Diesel Range Organics
GRO - Gasoli ne Range Organics
GW - Groundwater
J - Indi cates an estimated value
mg/L - milligram per liter
M TB E - Methyl Tert B utyl Ether
NA - Not analyzed
N D - Not detected
N D(5.0) - Not detected at or above the laboratory reporting l i mit, laboratory reporting limit included.
NM - Not monitored
NS - Not sampled
NSV D - Not surveyed to vertical datum
TPH - Total Petroleum Hydrocarbons

Figures

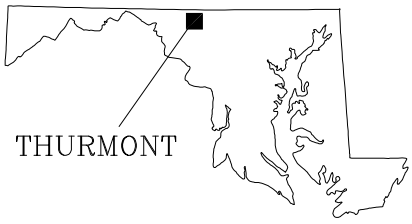


MAP SOURCE: ADC Frederick Co. MD. Street Map Book

SITE

0 1,000 2,000

SCALE: FEET



N

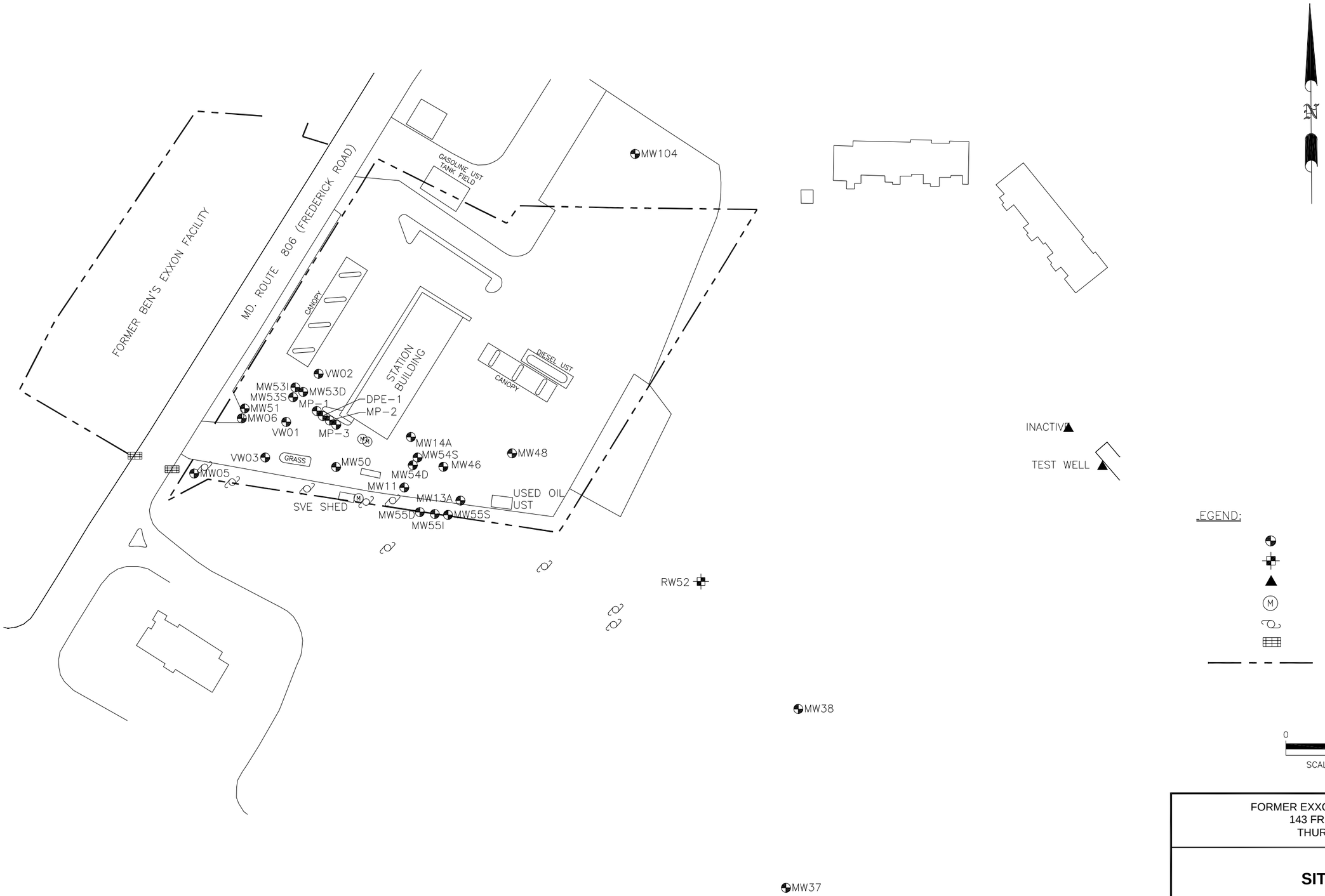
FORMER EXXON FACILITY # 25553
143 FREDERICK RD
THURMONT, MD.

LOCAL AREA MAP

ARCADIS

FIGURE
1

CITY:(Reqd) DIV/GROUP:(Reqd) DB:(Reqd) LD:(Opt) PIC:(Opt) PM:(Reqd) TM:(Opt) LYR:(Opt)ON=*;OFF=*REF*
G:\ENVCAD\BALTIMORE\ACT\B0085851\0014\00002\143 Frederick RD-Site-Loc.dwg LAYOUT: 1 SAVED: 4/15/2011 10:11 AM ACADVER: 18.0S (LMS TECH) PAGESETUP: ---- PLOTSTYLETABLE: ACAD.CTB PLOTTED: 4/29/2011 12:39 PM BY: GOFORTH, JOHN



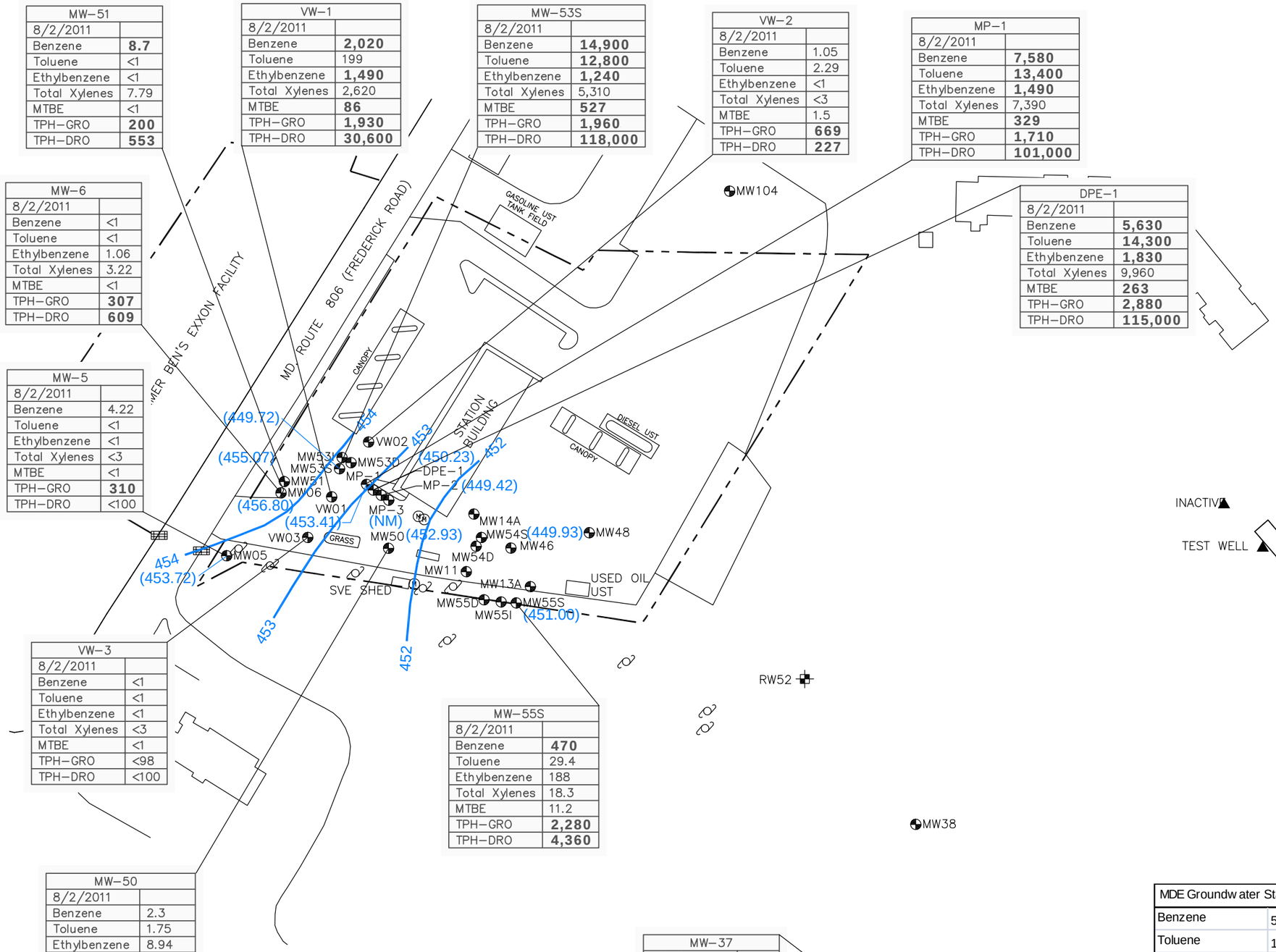
FORMER EXXON FACILITY # 25553
143 FREDERICK RD
THURMONT, MD.

SITE MAP

ARCADIS

FIGURE
2

CITY:(Reqd) DIV/GROUP:(Reqd) DB:(Reqd) LD:(Opt) PIC:(Opt) PM:(Reqd) TM:(Opt) LYR:(Opt)ON=*;OFF=*REF*
G:\ENVCAD\BALTIMORE\ACT\B0085510\0140\00021\43 Frederick Site Base.dwg LAYOUT: 2 SAVED: 4/29/2011 12:50 PM ACADVER: 18.0S (LMS TECH) PAGESETUP: ---- PLOTSTYLETABLE: ACAD.CTB PLOTTED: 4/29/2011 12:50 PM BY: GOFORTH, JOHN



MW-51	
8/2/2011	
Benzene	**8.7**
Toluene	<1
Ethylbenzene	<1
Total Xylenes	7.79
MTBE	<1
TPH-GRO	**200**
TPH-DRO	**553**

VW-1	
8/2/2011	
Benzene	**2,020**
Toluene	199
Ethylbenzene	**1,490**
Total Xylenes	2,620
MTBE	**86**
TPH-GRO	**1,930**
TPH-DRO	**30,600**

MW-53S	
8/2/2011	
Benzene	**14,900**
Toluene	**12,800**
Ethylbenzene	**1,240**
Total Xylenes	5,310
MTBE	**527**
TPH-GRO	**1,960**
TPH-DRO	**118,000**

VW-2	
8/2/2011	
Benzene	1.05
Toluene	2.29
Ethylbenzene	<1
Total Xylenes	<3
MTBE	1.5
TPH-GRO	**669**
TPH-DRO	**227**

MP-1	
8/2/2011	
Benzene	**7,580**
Toluene	**13,400**
Ethylbenzene	**1,490**
Total Xylenes	7,390
MTBE	**329**
TPH-GRO	**1,710**
TPH-DRO	**101,000**

DPE-1	
8/2/2011	
Benzene	**5,630**
Toluene	**14,300**
Ethylbenzene	**1,830**
Total Xylenes	9,960
MTBE	**263**
TPH-GRO	**2,880**
TPH-DRO	**115,000**

MW-6	
8/2/2011	
Benzene	<1
Toluene	<1
Ethylbenzene	1.06
Total Xylenes	3.22
MTBE	<1
TPH-GRO	**307**
TPH-DRO	**609**

MW-5	
8/2/2011	
Benzene	4.22
Toluene	<1
Ethylbenzene	<1
Total Xylenes	<3
MTBE	<1
TPH-GRO	**310**
TPH-DRO	<100

VW-3	
8/2/2011	
Benzene	<1
Toluene	<1
Ethylbenzene	<1
Total Xylenes	<3
MTBE	<1
TPH-GRO	<98
TPH-DRO	<100

MW-55S	
8/2/2011	
Benzene	**470**
Toluene	29.4
Ethylbenzene	188
Total Xylenes	18.3
MTBE	11.2
TPH-GRO	**2,280**
TPH-DRO	**4,360**

MW-50	
8/2/2011	
Benzene	2.3
Toluene	1.75
Ethylbenzene	8.94
Total Xylenes	9.23
MTBE	<1
TPH-GRO	**170**
TPH-DRO	**1,210**

MW-37	
8/2/2011	
Benzene	<1
Toluene	<1
Ethylbenzene	<1
Total Xylenes	<3
MTBE	<1
TPH-GRO	<111
TPH-DRO	<100

N

LEGEND:

- MONITORING WELL
- RECOVERY WELL
- MUNICIPAL SUPPLY WELL
- UTILITY MANHOLE
- UTILITY POLE
- CATCH BASIN
- PROPERTY LINE
- 444 — GROUNDWATER ELEVATION CONTOUR (FEET ABOVE MEAN SEA LEVEL)
- (440.50) GROUND WATER ELEVATION (FEET ABOVE MEAN SEA LEVEL)
- NM NOT MEASURED
- INACTIVE
- TEST WELL

NOTES:

[254] – MDE Groundwater Clean-Up Standard

MDE – Maryland Department of the Environment

Boldface – results exceed MDE groundwater clean-up standards

MTBE – methyl tertiary butyl ether

TPH-GRO – total petroleum hydrocarbons – gasoline range organics

TPH-DRO – total petroleum hydrocarbons – diesle range organics

All results listed in µg/L

0 50 100

SCALE: FEET

MDE Groundwater Standards	
Benzene	5
Toluene	1000
Ethylbenzene	700
Total Xylenes	10000
MTBE	20
Naphthalene	0.65
TPH-GRO	47
TPH-DRO	47

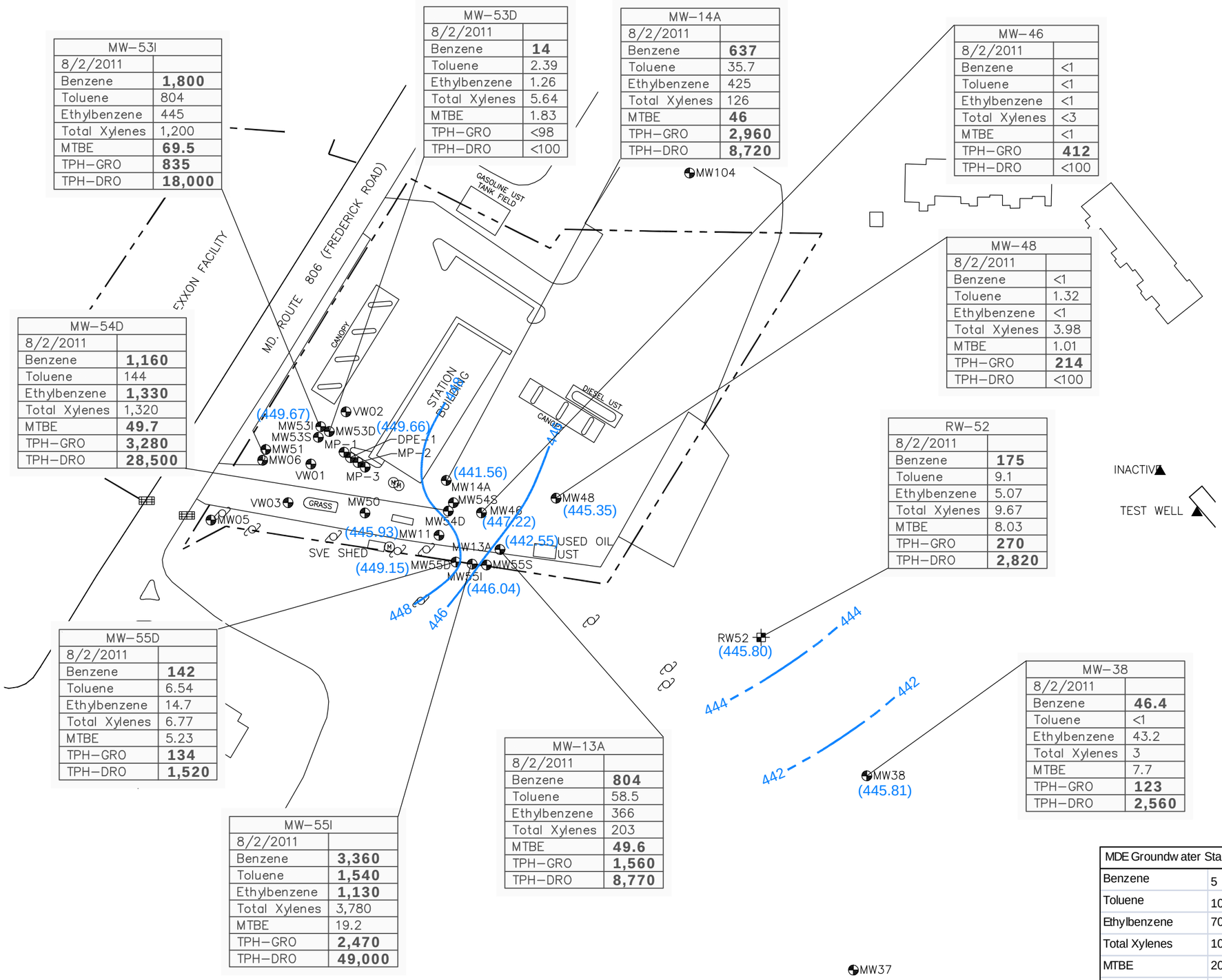
FORMER EXXON FACILITY # 25553
143 FREDERICK RD
THURMONT, MD.

SHALLOW GROUNDWATER ELEVATION CONTOURS AND ANALYTICAL DATA 2 AUGUST 2011

ARCADIS

FIGURE 3

CITY:(Req'd) DIV/GROUP:(Req'd) DB:(Req'd) LD:(Opt) PIC:(Opt) PM:(Req'd) TM:(Opt) LYR:(On)ON=*;OFF=*REF*
G:\ENVCAD\BALTIMORE\ACT\B0085951\00140\0002\Frederick Site-August-2011.dwg LAYOUT: 3 SAVED: 10/5/2011 3:46 PM ACADVER: 18.0S (LMS TECH) PAGESETUP: ---- PLOTSTYLETABLE: ACAD.CTB PLOTTED: 10/5/2011 3:47 PM BY: GOFORTH, JOHN



MW-53I	
8/2/2011	
Benzene	**1,800**
Toluene	804
Ethylbenzene	445
Total Xylenes	1,200
MTBE	**69.5**
TPH-GRO	**835**
TPH-DRO	**18,000**

MW-53D	
8/2/2011	
Benzene	**14**
Toluene	2.39
Ethylbenzene	1.26
Total Xylenes	5.64
MTBE	1.83
TPH-GRO	<98
TPH-DRO	<100

MW-14A	
8/2/2011	
Benzene	**637**
Toluene	35.7
Ethylbenzene	425
Total Xylenes	126
MTBE	**46**
TPH-GRO	**2,960**
TPH-DRO	**8,720**

MW-46	
8/2/2011	
Benzene	<1
Toluene	<1
Ethylbenzene	<1
Total Xylenes	<3
MTBE	<1
TPH-GRO	**412**
TPH-DRO	<100

MW-48	
8/2/2011	
Benzene	<1
Toluene	1.32
Ethylbenzene	<1
Total Xylenes	3.98
MTBE	1.01
TPH-GRO	**214**
TPH-DRO	<100

MW-54D	
8/2/2011	
Benzene	**1,160**
Toluene	144
Ethylbenzene	**1,330**
Total Xylenes	1,320
MTBE	**49.7**
TPH-GRO	**3,280**
TPH-DRO	**28,500**

RW-52	
8/2/2011	
Benzene	**175**
Toluene	9.1
Ethylbenzene	5.07
Total Xylenes	9.67
MTBE	8.03
TPH-GRO	**270**
TPH-DRO	**2,820**

MW-55D	
8/2/2011	
Benzene	**142**
Toluene	6.54
Ethylbenzene	14.7
Total Xylenes	6.77
MTBE	5.23
TPH-GRO	**134**
TPH-DRO	**1,520**

MW-13A	
8/2/2011	
Benzene	**804**
Toluene	58.5
Ethylbenzene	366
Total Xylenes	203
MTBE	**49.6**
TPH-GRO	**1,560**
TPH-DRO	**8,770**

MW-38	
8/2/2011	
Benzene	**46.4**
Toluene	<1
Ethylbenzene	43.2
Total Xylenes	3
MTBE	7.7
TPH-GRO	**123**
TPH-DRO	**2,560**

MW-55I	
8/2/2011	
Benzene	**3,360**
Toluene	**1,540**
Ethylbenzene	**1,130**
Total Xylenes	3,780
MTBE	19.2
TPH-GRO	**2,470**
TPH-DRO	**49,000**

LEGEND:

- MONITORING WELL
- RECOVERY WELL
- MUNICIPAL SUPPLY WELL
- UTILITY MANHOLE
- UTILITY POLE
- CATCH BASIN
- PROPERTY LINE
- 444 — GROUNDWATER ELEVATION CONTOUR (FEET ABOVE MEAN SEA LEVEL)
- (440.50) — GROUND WATER ELEVATION (FEET ABOVE MEAN SEA LEVEL)
- NM — NOT MEASURED

NOTES:

[254] – MDE Groundwater Clean-Up Standard

MDE – Maryland Department of the Environment

Boldface – results exceed MDE groundwater clean-up standards

MTBE – methyl tertiary butyl ether

TPH-GRO – total petroleum hydrocarbons – gasoline range organics

TPH-DRO – total petroleum hydrocarbons – diesle range organics

All results listed in µg/L

0 50 100
SCALE: FEET

MDE Groundwater Standards	
Benzene	5
Toluene	1000
Ethylbenzene	700
Total Xylenes	10000
MTBE	20
Naphthalene	0.65
TPH-GRO	47
TPH-DRO	47

FORMER EXXON FACILITY # 25553
143 FREDERICK RD
THURMONT, MD.

DEEP GROUNDWATER ELEVATION CONTOURS AND ANALYTICAL DATA 2 AUGUST 2011

ARCADIS

FIGURE **4**

CITY:(Reqd) DIV/GROUP:(Reqd) DB:(Reqd) LD:(Opt) PIC:(Opt) PM:(Reqd) TM:(Opt) LYR:(Opt)ON=*;OFF=*REF* PAGESETUP: ---- PLOTSTYLETABLE: ACAD.CTB PLOTTED: 10/5/2011 3:49 PM BY: GOFORTH, JOHN
G:\ENVCAD\BALTIMORE\ACT\B0085851\0014\00002\Frederick Site-August-2011.dwg LAYOUT: 4 SAVED: 10/5/2011 3:46 PM ACADVER: 18.0S (LMS TECH)

Appendix A

TestAmerica Laboratory Analytical Report (16 August 2011)

TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

ANALYTICAL REPORT

TestAmerica Laboratories, Inc.
TestAmerica Nashville
2960 Foster Creighton Road
Nashville, TN 37204
Tel: 800-765-0980

TestAmerica Job ID: NUH0938
Client Project/Site: 25553 - Thurmont, MD
Client Project Description: Exxon 25553

For:
ARCADIS Lexington - Exxon
1114 Benfield Boulevard, Suite A
Millersville, MD 21108

Attn: Rusty Kahl

Authorized for release by:
08/16/2011 01:55:37 PM

Heather Wagner
Project Manager
heather.wagner@testamericainc.com

LINKS

Review your project results through TotalAccess

Have a Question? Ask The Expert

Visit us at: www.testamericainc.com

This report has been electronically signed and authorized by the signatory. Electronic signature is intended to be the legally binding equivalent of a traditionally handwritten signature.

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Sample Summary

Client: ARCADIS Lexington - Exxon
Project/Site: 25553 - Thurmont, MD

TestAmerica Job ID: NUH0938

Lab Sample ID	Client Sample ID	Matrix	Collected	Received
NUH0938-01	MW-5 (08032011)	Ground Water	08/03/11 12:40	08/06/11 08:25
NUH0938-02	MW-6 (08032011)	Ground Water	08/03/11 15:05	08/06/11 08:25
NUH0938-03	MW-13A (08022011)	Ground Water	08/02/11 20:15	08/06/11 08:25
NUH0938-04	MW-14A (08022011)	Ground Water	08/02/11 20:05	08/06/11 08:25
NUH0938-05	MW-37 (08022011)	Ground Water	08/02/11 15:20	08/06/11 08:25
NUH0938-06	MW-38 (08022011)	Ground Water	08/02/11 16:35	08/06/11 08:25
NUH0938-07	MW-46 (08022011)	Ground Water	08/02/11 12:45	08/06/11 08:25
NUH0938-08	MW-48 (08022011)	Ground Water	08/02/11 18:00	08/06/11 08:25
NUH0938-09	MW-50 (08032011)	Ground Water	08/03/11 09:30	08/06/11 08:25
NUH0938-10	MW-51 (08032011)	Ground Water	08/03/11 18:00	08/06/11 08:25
NUH0938-11	MW-53D (08032011)	Ground Water	08/03/11 15:10	08/06/11 08:25
NUH0938-12	MW-53I (08032011)	Ground Water	08/03/11 12:10	08/06/11 08:25
NUH0938-13	MW-53S (08032011)	Ground Water	08/03/11 11:15	08/06/11 08:25
NUH0938-14	MW-54D (08032011)	Ground Water	08/03/11 10:15	08/06/11 08:25
NUH0938-15	MW-55D (08022011)	Ground Water	08/02/11 12:20	08/06/11 08:25
NUH0938-16	MW-55S (08022011)	Ground Water	08/02/11 13:05	08/06/11 08:25
NUH0938-17	RW-52 (08022011)	Ground Water	08/02/11 18:05	08/06/11 08:25
NUH0938-18	VW-1 (08032011)	Ground Water	08/03/11 15:45	08/06/11 08:25
NUH0938-19	VW-2 (08032011)	Ground Water	08/03/11 16:05	08/06/11 08:25
NUH0938-20	VW-3 (08032011)	Ground Water	08/03/11 10:40	08/06/11 08:25
NUH0938-21	MP-1 (08032011)	Ground Water	08/03/11 16:50	08/06/11 08:25
NUH0938-22	DPE-1 (08032011)	Ground Water	08/03/11 17:45	08/06/11 08:25
NUH0938-23	DUP-1 (08022011)	Ground Water	08/02/11 12:00	08/06/11 08:25
NUH0938-24	DUP-2 (08032011)	Ground Water	08/03/11 12:00	08/06/11 08:25
NUH0938-25	RB-1 (08022011)	Ground Water	08/02/11 15:20	08/06/11 08:25
NUH0938-26	RB-2 (08032011)	Ground Water	08/03/11 17:55	08/06/11 08:25
NUH0938-27	TB-1 (08032011)	Water	08/03/11 00:01	08/06/11 08:25
NUH0938-28	TB-2 (08032011)	Water	08/03/11 00:01	08/06/11 08:25
NUH0938-29	MW-55I (08022011)	Ground Water	08/02/11 15:40	08/06/11 08:25

Qualifiers

GCMS Volatiles

Qualifier	Qualifier Description
M8	The MS and/or MSD were below the acceptance limits. See Blank Spike (LCS).
R2	The RPD exceeded the acceptance limit.
L1	Laboratory Control Sample and/or Laboratory Control Sample Duplicate recovery was above acceptance limits.
L	Laboratory Control Sample and/or Laboratory Control Sample Duplicate recovery was above the acceptance limits. Analyte not detected, data not impacted.

GC Semivolatiles

Qualifier	Qualifier Description
MNR1	There was no MS/MSD analyzed with this batch due to insufficient sample volume. See Blank Spike.
ZX	Due to sample matrix effects, the surrogate recovery was outside the acceptance limits.
S10	Insufficient sample available for reanalysis.
Z6	Surrogate recovery was below acceptance limits.

Glossary

Abbreviation	These commonly used abbreviations may or may not be present in this report.
☼	Listed under the "D" column to designate that the result is reported on a dry weight basis
%R	Percent Recovery
DL, RA, RE, IN	Indicates a Dilution, Reanalysis, Re-extraction, or additional Initial metals/anion analysis of the sample
EDL	Estimated Detection Limit (Dioxin)
EPA	United States Environmental Protection Agency
MDL	Method Detection Limit
ML	Minimum Level (Dioxin)
ND	Not detected at the reporting limit (or method detection limit if shown)
PQL	Practical Quantitation Limit
RL	Reporting Limit
RPD	Relative Percent Difference, a measure of the relative difference between two points
TEF	Toxicity Equivalent Factor (Dioxin)
TEQ	Toxicity Equivalent Quotient (Dioxin)

Client Sample ID: MW-5 (08032011)

Date Collected: 08/03/11 12:40
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-01
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Benzene	**4.22**		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Bromoform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Chloroform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 17:45	08/09/11 21:49	1.00

Client Sample ID: MW-5 (08032011)

Date Collected: 08/03/11 12:40
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-01
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1 (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Styrene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Toluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00
Xylenes, total	ND		3.00		ug/L		08/09/11 17:45	08/09/11 21:49	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*98*		*63 - 140*	*08/09/11 17:45*	*08/09/11 21:49*	*1.00*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 17:45*	*08/09/11 21:49*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 17:45*	*08/09/11 21:49*	*1.00*
4-Bromofluorobenzene	*90*		*79 - 125*	*08/09/11 17:45*	*08/09/11 21:49*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/03/11 12:40	08/10/11 13:54	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*98*		*60 - 150*	*08/03/11 12:40*	*08/10/11 13:54*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**310**		99.0		ug/L		08/08/11 12:05	08/09/11 07:36	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*84*		*53 - 140*	*08/08/11 12:05*	*08/09/11 07:36*	*1.00*

Client Sample ID: MW-6 (08032011)

Date Collected: 08/03/11 15:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-02
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Benzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Bromoform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
n-Butylbenzene	**1.04**		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Chloroform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Ethylbenzene	**1.06**		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Isopropylbenzene	**1.34**		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 17:45	08/09/11 22:17	1.00

Client Sample ID: MW-6 (08032011)

Date Collected: 08/03/11 15:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-02

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1 (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
n-Propylbenzene	**2.18**		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Styrene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Toluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00
Xylenes, total	**3.22**		3.00		ug/L		08/09/11 17:45	08/09/11 22:17	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/09/11 17:45*	*08/09/11 22:17*	*1.00*
Dibromofluoromethane	*104*		*73 - 131*	*08/09/11 17:45*	*08/09/11 22:17*	*1.00*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/09/11 22:17*	*1.00*
4-Bromofluorobenzene	*91*		*79 - 125*	*08/09/11 17:45*	*08/09/11 22:17*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**609**		100		ug/L		08/03/11 15:05	08/10/11 14:24	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*93*		*60 - 150*	*08/03/11 15:05*	*08/10/11 14:24*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**307**		94.3		ug/L		08/08/11 12:05	08/09/11 08:00	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*81*		*53 - 140*	*08/08/11 12:05*	*08/09/11 08:00*	*1.00*

Client Sample ID: MW-13A (08022011)

Lab Sample ID: NUH0938-03

Date Collected: 08/02/11 20:15

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Tert-Amyl Methyl Ether	**1.95**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Diisopropyl Ether	**9.63**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Methyl tert-Butyl Ether	**49.6**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Tertiary Butyl Alcohol	**99.7**		20.0		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
sec-Butylbenzene	**4.97**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
n-Butylbenzene	**8.78**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
cis-1,2-Dichloroethene	**3.03**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Isopropylbenzene	**50.0**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
p-Isopropyltoluene	**1.83**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Naphthalene	**100**		5.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
n-Propylbenzene	**98.6**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00

Client Sample ID: MW-13A (08022011)

Date Collected: 08/02/11 20:15
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-03
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Toluene	**58.5**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,3,5-Trimethylbenzene	**22.0**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
1,2,4-Trimethylbenzene	**41.6**		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00
Xylenes, total	**203**		3.00		ug/L		08/08/11 17:45	08/09/11 04:02	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*103*		*63 - 140*	*08/08/11 17:45*	*08/09/11 04:02*	*1.00*
Dibromofluoromethane	*101*		*73 - 131*	*08/08/11 17:45*	*08/09/11 04:02*	*1.00*
Toluene-d8	*97*		*80 - 120*	*08/08/11 17:45*	*08/09/11 04:02*	*1.00*
4-Bromofluorobenzene	*92*		*79 - 125*	*08/08/11 17:45*	*08/09/11 04:02*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**804**		10.0		ug/L		08/09/11 17:45	08/10/11 00:09	10.0
Ethylbenzene	**366**		10.0		ug/L		08/09/11 17:45	08/10/11 00:09	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*97*		*63 - 140*	*08/09/11 17:45*	*08/10/11 00:09*	*10.0*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 17:45*	*08/10/11 00:09*	*10.0*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/10/11 00:09*	*10.0*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/09/11 17:45*	*08/10/11 00:09*	*10.0*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**8770**		1000		ug/L		08/02/11 20:15	08/11/11 19:01	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*98*		*60 - 150*	*08/02/11 20:15*	*08/11/11 19:01*	*10.0*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**1560**		94.3		ug/L		08/08/11 12:05	08/09/11 21:18	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*43*	*ZX*	*53 - 140*	*08/08/11 12:05*	*08/09/11 21:18*	*1.00*

Client Sample ID: MW-14A (08022011)

Date Collected: 08/02/11 20:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-04
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Diisopropyl Ether	**10.1**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Methyl tert-Butyl Ether	**46.0**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Tertiary Butyl Alcohol	**113**		20.0		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
sec-Butylbenzene	**6.51**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
n-Butylbenzene	**14.4**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Isopropylbenzene	**42.6**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
p-Isopropyltoluene	**3.11**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Naphthalene	**132**		5.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
n-Propylbenzene	**89.3**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00

Client Sample ID: MW-14A (08022011)

Date Collected: 08/02/11 20:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-04

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Toluene	**35.7**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,3,5-Trimethylbenzene	**17.1**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
1,2,4-Trimethylbenzene	**28.8**		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00
Xylenes, total	**126**		3.00		ug/L		08/08/11 17:45	08/09/11 04:30	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*101*		*63 - 140*	*08/08/11 17:45*	*08/09/11 04:30*	*1.00*
Dibromofluoromethane	*103*		*73 - 131*	*08/08/11 17:45*	*08/09/11 04:30*	*1.00*
Toluene-d8	*98*		*80 - 120*	*08/08/11 17:45*	*08/09/11 04:30*	*1.00*
4-Bromofluorobenzene	*91*		*79 - 125*	*08/08/11 17:45*	*08/09/11 04:30*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**637**		10.0		ug/L		08/09/11 17:45	08/10/11 00:37	10.0
Ethylbenzene	**425**		10.0		ug/L		08/09/11 17:45	08/10/11 00:37	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*97*		*63 - 140*	*08/09/11 17:45*	*08/10/11 00:37*	*10.0*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 17:45*	*08/10/11 00:37*	*10.0*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/10/11 00:37*	*10.0*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/09/11 17:45*	*08/10/11 00:37*	*10.0*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**8720**		1000		ug/L		08/02/11 20:05	08/11/11 19:31	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*95*		*60 - 150*	*08/02/11 20:05*	*08/11/11 19:31*	*10.0*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**2960**		94.3		ug/L		08/08/11 12:05	08/09/11 08:24	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*57*		*53 - 140*	*08/08/11 12:05*	*08/09/11 08:24*	*1.00*

Client Sample ID: MW-37 (08022011)

Date Collected: 08/02/11 15:20
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-05
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Benzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
n-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Ethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Isopropylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 04:58	1.00

Client Sample ID: MW-37 (08022011)

Date Collected: 08/02/11 15:20
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-05

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
n-Propylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Toluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00
Xylenes, total	ND		3.00		ug/L		08/08/11 17:45	08/09/11 04:58	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*96*		*63 - 140*	*08/08/11 17:45*	*08/09/11 04:58*	*1.00*
Dibromofluoromethane	*101*		*73 - 131*	*08/08/11 17:45*	*08/09/11 04:58*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/08/11 17:45*	*08/09/11 04:58*	*1.00*
4-Bromofluorobenzene	*93*		*79 - 125*	*08/08/11 17:45*	*08/09/11 04:58*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/02/11 15:20	08/10/11 15:54	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*92*		*60 - 150*	*08/02/11 15:20*	*08/10/11 15:54*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	ND		111		ug/L		08/08/11 12:05	08/09/11 18:43	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*85*		*53 - 140*	*08/08/11 12:05*	*08/09/11 18:43*	*1.00*

Client Sample ID: MW-38 (08022011)

Date Collected: 08/02/11 16:35
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-06

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Benzene	**46.4**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Diisopropyl Ether	**2.31**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Methyl tert-Butyl Ether	**7.70**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
sec-Butylbenzene	**4.29**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
n-Butylbenzene	**3.36**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Ethylbenzene	**43.2**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Isopropylbenzene	**22.7**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 05:26	1.00

Client Sample ID: MW-38 (08022011)

Date Collected: 08/02/11 16:35
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-06

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
n-Propylbenzene	**16.6**		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Toluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00
Xylenes, total	**3.00**		3.00		ug/L		08/08/11 17:45	08/09/11 05:26	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/08/11 17:45*	*08/09/11 05:26*	*1.00*
Dibromofluoromethane	*106*		*73 - 131*	*08/08/11 17:45*	*08/09/11 05:26*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/08/11 17:45*	*08/09/11 05:26*	*1.00*
4-Bromofluorobenzene	*92*		*79 - 125*	*08/08/11 17:45*	*08/09/11 05:26*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**2560**		100		ug/L		08/02/11 16:35	08/10/11 16:24	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*90*		*60 - 150*	*08/02/11 16:35*	*08/10/11 16:24*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**123**		97.1		ug/L		08/08/11 12:05	08/09/11 19:02	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*84*		*53 - 140*	*08/08/11 12:05*	*08/09/11 19:02*	*1.00*

Client Sample ID: MW-46 (08022011)

Date Collected: 08/02/11 12:45
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-07
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Benzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 09:10	1.00

Client Sample ID: MW-46 (08022011)

Lab Sample ID: NUH0938-07

Date Collected: 08/02/11 12:45

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Toluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00
Xylenes, total	ND		3.00		ug/L		08/09/11 05:54	08/09/11 09:10	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*100*		*63 - 140*	*08/09/11 05:54*	*08/09/11 09:10*	*1.00*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 05:54*	*08/09/11 09:10*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/09/11 05:54*	*08/09/11 09:10*	*1.00*
4-Bromofluorobenzene	*92*		*79 - 125*	*08/09/11 05:54*	*08/09/11 09:10*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/02/11 12:45	08/10/11 16:54	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*94*		*60 - 150*	*08/02/11 12:45*	*08/10/11 16:54*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**412**		95.2		ug/L		08/08/11 12:05	08/09/11 21:38	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*24*	*ZX*	*53 - 140*	*08/08/11 12:05*	*08/09/11 21:38*	*1.00*

Client Sample ID: MW-48 (08022011)

Date Collected: 08/02/11 18:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-08
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Benzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Methyl tert-Butyl Ether	**1.01**		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 09:38	1.00

Client Sample ID: MW-48 (08022011)

Date Collected: 08/02/11 18:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-08

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Toluene	**1.32**		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
1,2,4-Trimethylbenzene	**1.07**		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00
Xylenes, total	**3.98**		3.00		ug/L		08/09/11 05:54	08/09/11 09:38	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/09/11 05:54*	*08/09/11 09:38*	*1.00*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 05:54*	*08/09/11 09:38*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 05:54*	*08/09/11 09:38*	*1.00*
4-Bromofluorobenzene	*90*		*79 - 125*	*08/09/11 05:54*	*08/09/11 09:38*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/02/11 18:00	08/10/11 17:24	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*96*		*60 - 150*	*08/02/11 18:00*	*08/10/11 17:24*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**214**		99.0		ug/L		08/08/11 12:05	08/09/11 21:57	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*75*		*53 - 140*	*08/08/11 12:05*	*08/09/11 21:57*	*1.00*

Client Sample ID: MW-50 (08032011)

Date Collected: 08/03/11 09:30
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-09
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Benzene	**2.30**		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
sec-Butylbenzene	**2.73**		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
n-Butylbenzene	**1.88**		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Ethylbenzene	**8.94**		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Isopropylbenzene	**8.25**		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 10:06	1.00

Client Sample ID: MW-50 (08032011)

Date Collected: 08/03/11 09:30
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-09

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
n-Propylbenzene	**11.6**		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Toluene	**1.75**		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00
Xylenes, total	**9.23**		3.00		ug/L		08/09/11 05:54	08/09/11 10:06	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*105*		*63 - 140*	*08/09/11 05:54*	*08/09/11 10:06*	*1.00*
Dibromofluoromethane	*108*		*73 - 131*	*08/09/11 05:54*	*08/09/11 10:06*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 05:54*	*08/09/11 10:06*	*1.00*
4-Bromofluorobenzene	*90*		*79 - 125*	*08/09/11 05:54*	*08/09/11 10:06*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**1210**		100		ug/L		08/03/11 09:30	08/10/11 17:53	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*90*		*60 - 150*	*08/03/11 09:30*	*08/10/11 17:53*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**170**		100		ug/L		08/08/11 11:05	08/10/11 14:48	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*107*		*53 - 140*	*08/08/11 11:05*	*08/10/11 14:48*	*1.00*

Client Sample ID: MW-51 (08032011)

Date Collected: 08/03/11 18:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-10

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Benzene	**8.70**		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 10:34	1.00

Client Sample ID: MW-51 (08032011)

Date Collected: 08/03/11 18:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-10

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Toluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,3,5-Trimethylbenzene	**1.02**		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
1,2,4-Trimethylbenzene	**3.49**		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00
Xylenes, total	**7.79**		3.00		ug/L		08/09/11 05:54	08/09/11 10:34	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*105*		*63 - 140*	*08/09/11 05:54*	*08/09/11 10:34*	*1.00*
Dibromofluoromethane	*106*		*73 - 131*	*08/09/11 05:54*	*08/09/11 10:34*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 05:54*	*08/09/11 10:34*	*1.00*
4-Bromofluorobenzene	*91*		*79 - 125*	*08/09/11 05:54*	*08/09/11 10:34*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**553**		100		ug/L		08/03/11 18:00	08/10/11 18:23	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*92*		*60 - 150*	*08/03/11 18:00*	*08/10/11 18:23*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**200**		98.0		ug/L		08/08/11 11:05	08/10/11 15:12	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*92*		*53 - 140*	*08/08/11 11:05*	*08/10/11 15:12*	*1.00*

Client Sample ID: MW-53D (08032011)

Date Collected: 08/03/11 15:10
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-11
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Benzene	**14.0**		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Methyl tert-Butyl Ether	**1.83**		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Ethylbenzene	**1.26**		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 11:02	1.00

Client Sample ID: MW-53D (08032011)

Date Collected: 08/03/11 15:10
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-11
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Toluene	**2.39**		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
1,2,4-Trimethylbenzene	**1.34**		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00
Xylenes, total	**5.64**		3.00		ug/L		08/09/11 05:54	08/09/11 11:02	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*103*		*63 - 140*	*08/09/11 05:54*	*08/09/11 11:02*	*1.00*
Dibromofluoromethane	*104*		*73 - 131*	*08/09/11 05:54*	*08/09/11 11:02*	*1.00*
Toluene-d8	*94*		*80 - 120*	*08/09/11 05:54*	*08/09/11 11:02*	*1.00*
4-Bromofluorobenzene	*92*		*79 - 125*	*08/09/11 05:54*	*08/09/11 11:02*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/03/11 15:10	08/11/11 13:32	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*94*		*60 - 150*	*08/03/11 15:10*	*08/11/11 13:32*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	ND		98.0		ug/L		08/08/11 11:05	08/10/11 15:36	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*102*		*53 - 140*	*08/08/11 11:05*	*08/10/11 15:36*	*1.00*

Client Sample ID: MW-53I (08032011)

Date Collected: 08/03/11 12:10
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-12
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Tert-Amyl Methyl Ether	**3.24**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Diisopropyl Ether	**6.24**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Methyl tert-Butyl Ether	**69.5**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Tertiary Butyl Alcohol	**98.7**		20.0		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
sec-Butylbenzene	**3.50**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
n-Butylbenzene	**5.42**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Isopropylbenzene	**28.4**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
p-Isopropyltoluene	**2.14**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Naphthalene	**141**		5.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
n-Propylbenzene	**53.8**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00

Client Sample ID: MW-53I (08032011)

Date Collected: 08/03/11 12:10
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-12

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
1,3,5-Trimethylbenzene	**79.0**		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:30	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/09/11 05:54*	*08/09/11 11:30*	*1.00*
Dibromofluoromethane	*100*		*73 - 131*	*08/09/11 05:54*	*08/09/11 11:30*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 05:54*	*08/09/11 11:30*	*1.00*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/09/11 05:54*	*08/09/11 11:30*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**1800**		10.0		ug/L		08/09/11 17:45	08/10/11 01:05	10.0
Ethylbenzene	**445**		10.0		ug/L		08/09/11 17:45	08/10/11 01:05	10.0
Toluene	**804**		10.0		ug/L		08/09/11 17:45	08/10/11 01:05	10.0
1,2,4-Trimethylbenzene	**312**		10.0		ug/L		08/09/11 17:45	08/10/11 01:05	10.0
Xylenes, total	**1200**		30.0		ug/L		08/09/11 17:45	08/10/11 01:05	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*101*		*63 - 140*	*08/09/11 17:45*	*08/10/11 01:05*	*10.0*
Dibromofluoromethane	*108*		*73 - 131*	*08/09/11 17:45*	*08/10/11 01:05*	*10.0*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/10/11 01:05*	*10.0*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/09/11 17:45*	*08/10/11 01:05*	*10.0*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**18000**		2000		ug/L		08/03/11 12:10	08/11/11 18:01	20.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*99*		*60 - 150*	*08/03/11 12:10*	*08/11/11 18:01*	*20.0*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**835**		97.1		ug/L		08/08/11 11:05	08/10/11 16:00	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*116*		*53 - 140*	*08/08/11 11:05*	*08/10/11 16:00*	*1.00*

Client Sample ID: MW-53S (08032011)

Date Collected: 08/03/11 11:15
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-13
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Diisopropyl Ether	**27.1**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Tertiary Butyl Alcohol	**1700**		20.0		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
sec-Butylbenzene	**3.80**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
n-Butylbenzene	**13.2**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2-Dibromoethane (EDB)	**14.8**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Isopropylbenzene	**39.4**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
p-Isopropyltoluene	**2.86**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
n-Propylbenzene	**79.2**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00

Client Sample ID: MW-53S (08032011)

Date Collected: 08/03/11 11:15
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-13
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
1,3,5-Trimethylbenzene	**166**		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 11:58	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*91*		*63 - 140*	*08/09/11 05:54*	*08/09/11 11:58*	*1.00*
Dibromofluoromethane	*125*		*73 - 131*	*08/09/11 05:54*	*08/09/11 11:58*	*1.00*
Toluene-d8	*83*		*80 - 120*	*08/09/11 05:54*	*08/09/11 11:58*	*1.00*
4-Bromofluorobenzene	*94*		*79 - 125*	*08/09/11 05:54*	*08/09/11 11:58*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methyl tert-Butyl Ether	**527**		10.0		ug/L		08/10/11 06:14	08/10/11 09:30	10.0
Ethylbenzene	**1240**		10.0		ug/L		08/10/11 06:14	08/10/11 09:30	10.0
Naphthalene	**247**		50.0		ug/L		08/10/11 06:14	08/10/11 09:30	10.0
1,2,4-Trimethylbenzene	**724**		10.0		ug/L		08/10/11 06:14	08/10/11 09:30	10.0
Xylenes, total	**5310**		30.0		ug/L		08/10/11 06:14	08/10/11 09:30	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*101*		*63 - 140*	*08/10/11 06:14*	*08/10/11 09:30*	*10.0*
Dibromofluoromethane	*101*		*73 - 131*	*08/10/11 06:14*	*08/10/11 09:30*	*10.0*
Toluene-d8	*93*		*80 - 120*	*08/10/11 06:14*	*08/10/11 09:30*	*10.0*
4-Bromofluorobenzene	*88*		*79 - 125*	*08/10/11 06:14*	*08/10/11 09:30*	*10.0*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE2

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**14900**		100		ug/L		08/10/11 06:14	08/10/11 09:59	100
Toluene	**12800**		100		ug/L		08/10/11 06:14	08/10/11 09:59	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*102*		*63 - 140*	*08/10/11 06:14*	*08/10/11 09:59*	*100*
Dibromofluoromethane	*108*		*73 - 131*	*08/10/11 06:14*	*08/10/11 09:59*	*100*
Toluene-d8	*94*		*80 - 120*	*08/10/11 06:14*	*08/10/11 09:59*	*100*
4-Bromofluorobenzene	*88*		*79 - 125*	*08/10/11 06:14*	*08/10/11 09:59*	*100*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**118000**		10000		ug/L		08/03/11 11:15	08/11/11 15:32	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*98*		*60 - 150*	*08/03/11 11:15*	*08/11/11 15:32*	*100*

Client Sample Results

Client: ARCADIS Lexington - Exxon
Project/Site: 25553 - Thurmont, MD

TestAmerica Job ID: NUH0938

Client Sample ID: MW-53S (08032011)

Date Collected: 08/03/11 11:15
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-13

Matrix: Ground Water

Method: SW846 8015B - Extractable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	1960		196		ug/L		08/08/11 11:05	08/11/11 06:23	2.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*76*		*53 - 140*	*08/08/11 11:05*	*08/11/11 06:23*	*2.00*

Client Sample ID: MW-54D (08032011)

Lab Sample ID: NUH0938-14

Date Collected: 08/03/11 10:15

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Diisopropyl Ether	**13.7**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Methyl tert-Butyl Ether	**49.7**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Tertiary Butyl Alcohol	**384**		20.0		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
sec-Butylbenzene	**11.3**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
n-Butylbenzene	**33.0**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Isopropylbenzene	**99.5**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
p-Isopropyltoluene	**9.36**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
n-Propylbenzene	**182**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00

Client Sample ID: MW-54D (08032011)

Date Collected: 08/03/11 10:15
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-14
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Toluene	**144**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
1,3,5-Trimethylbenzene	**180**		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 12:26	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*106*		*63 - 140*	*08/09/11 05:54*	*08/09/11 12:26*	*1.00*
Dibromofluoromethane	*106*		*73 - 131*	*08/09/11 05:54*	*08/09/11 12:26*	*1.00*
Toluene-d8	*94*		*80 - 120*	*08/09/11 05:54*	*08/09/11 12:26*	*1.00*
4-Bromofluorobenzene	*101*		*79 - 125*	*08/09/11 05:54*	*08/09/11 12:26*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**1160**		10.0		ug/L		08/09/11 17:45	08/10/11 01:33	10.0
Ethylbenzene	**1330**		10.0		ug/L		08/09/11 17:45	08/10/11 01:33	10.0
Naphthalene	**476**		50.0		ug/L		08/09/11 17:45	08/10/11 01:33	10.0
1,2,4-Trimethylbenzene	**964**		10.0		ug/L		08/09/11 17:45	08/10/11 01:33	10.0
Xylenes, total	**1320**		30.0		ug/L		08/09/11 17:45	08/10/11 01:33	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/09/11 17:45*	*08/10/11 01:33*	*10.0*
Dibromofluoromethane	*106*		*73 - 131*	*08/09/11 17:45*	*08/10/11 01:33*	*10.0*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/10/11 01:33*	*10.0*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/09/11 17:45*	*08/10/11 01:33*	*10.0*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**28500**		2500		ug/L		08/03/11 10:15	08/11/11 17:31	25.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*92*		*60 - 150*	*08/03/11 10:15*	*08/11/11 17:31*	*25.0*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**3280**		211		ug/L		08/08/11 11:05	08/11/11 06:47	2.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*71*		*53 - 140*	*08/08/11 11:05*	*08/11/11 06:47*	*2.00*

Client Sample ID: MW-55D (08022011)

Date Collected: 08/02/11 12:20
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-15
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Benzene	**142**		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Diisopropyl Ether	**14.4**		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Methyl tert-Butyl Ether	**5.23**		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Bromoform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Chloroform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Ethylbenzene	**14.7**		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Isopropylbenzene	**4.33**		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 17:45	08/09/11 22:45	1.00

Client Sample ID: MW-55D (08022011)

Date Collected: 08/02/11 12:20
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-15
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1 (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
n-Propylbenzene	**3.47**		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Styrene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Toluene	**6.54**		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00
Xylenes, total	**6.77**		3.00		ug/L		08/09/11 17:45	08/09/11 22:45	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/09/11 17:45*	*08/09/11 22:45*	*1.00*
Dibromofluoromethane	*106*		*73 - 131*	*08/09/11 17:45*	*08/09/11 22:45*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 17:45*	*08/09/11 22:45*	*1.00*
4-Bromofluorobenzene	*90*		*79 - 125*	*08/09/11 17:45*	*08/09/11 22:45*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**1520**		100		ug/L		08/02/11 12:20	08/11/11 15:02	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*94*		*60 - 150*	*08/02/11 12:20*	*08/11/11 15:02*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**134**		100		ug/L		08/08/11 11:05	08/10/11 17:12	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*108*		*53 - 140*	*08/08/11 11:05*	*08/10/11 17:12*	*1.00*

Client Sample ID: MW-55S (08022011)

Lab Sample ID: NUH0938-16

Date Collected: 08/02/11 13:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Diisopropyl Ether	**1.05**		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Methyl tert-Butyl Ether	**11.2**		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Tertiary Butyl Alcohol	**37.4**		20.0		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Ethylbenzene	**188**		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Isopropylbenzene	**11.2**		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Naphthalene	**33.4**		5.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00

Client Sample ID: MW-55S (08022011)

Date Collected: 08/02/11 13:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-16

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
n-Propylbenzene	**11.8**		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Toluene	**29.4**		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
1,2,4-Trimethylbenzene	**3.09**		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00
Xylenes, total	**18.3**		3.00		ug/L		08/09/11 05:54	08/09/11 13:23	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*98*		*63 - 140*	*08/09/11 05:54*	*08/09/11 13:23*	*1.00*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 05:54*	*08/09/11 13:23*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/09/11 05:54*	*08/09/11 13:23*	*1.00*
4-Bromofluorobenzene	*92*		*79 - 125*	*08/09/11 05:54*	*08/09/11 13:23*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**470**		10.0		ug/L		08/09/11 17:45	08/10/11 02:01	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*97*		*63 - 140*	*08/09/11 17:45*	*08/10/11 02:01*	*10.0*
Dibromofluoromethane	*104*		*73 - 131*	*08/09/11 17:45*	*08/10/11 02:01*	*10.0*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/10/11 02:01*	*10.0*
4-Bromofluorobenzene	*90*		*79 - 125*	*08/09/11 17:45*	*08/10/11 02:01*	*10.0*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**4360**		100		ug/L		08/02/11 13:05	08/10/11 21:23	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*99*		*60 - 150*	*08/02/11 13:05*	*08/10/11 21:23*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**2280**		97.1		ug/L		08/08/11 11:05	08/10/11 17:36	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*58*		*53 - 140*	*08/08/11 11:05*	*08/10/11 17:36*	*1.00*

Client Sample ID: RW-52 (08022011)

Lab Sample ID: NUH0938-17

Date Collected: 08/02/11 18:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Benzene	**175**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Diisopropyl Ether	**4.27**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Methyl tert-Butyl Ether	**8.03**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
sec-Butylbenzene	**3.54**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
n-Butylbenzene	**6.24**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Ethylbenzene	**5.07**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Isopropylbenzene	**13.7**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 13:51	1.00

Client Sample ID: RW-52 (08022011)

Date Collected: 08/02/11 18:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-17

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
n-Propylbenzene	**17.7**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Toluene	**9.10**		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00
Xylenes, total	**9.67**		3.00		ug/L		08/09/11 05:54	08/09/11 13:51	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*98*		*63 - 140*	*08/09/11 05:54*	*08/09/11 13:51*	*1.00*
Dibromofluoromethane	*106*		*73 - 131*	*08/09/11 05:54*	*08/09/11 13:51*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/09/11 05:54*	*08/09/11 13:51*	*1.00*
4-Bromofluorobenzene	*91*		*79 - 125*	*08/09/11 05:54*	*08/09/11 13:51*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**2820**		100		ug/L		08/02/11 18:05	08/10/11 21:52	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*94*		*60 - 150*	*08/02/11 18:05*	*08/10/11 21:52*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**270**		95.2		ug/L		08/08/11 11:05	08/10/11 18:00	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*103*		*53 - 140*	*08/08/11 11:05*	*08/10/11 18:00*	*1.00*

Client Sample ID: VW-1 (08032011)

Lab Sample ID: NUH0938-18

Date Collected: 08/03/11 15:45

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Diisopropyl Ether	**7.06**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Methyl tert-Butyl Ether	**86.0**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Tertiary Butyl Alcohol	**262**		20.0		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
sec-Butylbenzene	**6.75**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
n-Butylbenzene	**17.5**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Isopropylbenzene	**69.9**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
p-Isopropyltoluene	**7.16**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
n-Propylbenzene	**121**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00

Client Sample ID: VW-1 (08032011)

Date Collected: 08/03/11 15:45
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-18

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Toluene	**199**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
1,3,5-Trimethylbenzene	**169**		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 14:19	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*97*		*63 - 140*	*08/09/11 05:54*	*08/09/11 14:19*	*1.00*
Dibromofluoromethane	*100*		*73 - 131*	*08/09/11 05:54*	*08/09/11 14:19*	*1.00*
Toluene-d8	*94*		*80 - 120*	*08/09/11 05:54*	*08/09/11 14:19*	*1.00*
4-Bromofluorobenzene	*96*		*79 - 125*	*08/09/11 05:54*	*08/09/11 14:19*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**2020**		20.0		ug/L		08/09/11 17:45	08/10/11 02:57	20.0
Ethylbenzene	**1490**		20.0		ug/L		08/09/11 17:45	08/10/11 02:57	20.0
Naphthalene	**387**		100		ug/L		08/09/11 17:45	08/10/11 02:57	20.0
1,2,4-Trimethylbenzene	**1140**		20.0		ug/L		08/09/11 17:45	08/10/11 02:57	20.0
Xylenes, total	**2620**		60.0		ug/L		08/09/11 17:45	08/10/11 02:57	20.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*97*		*63 - 140*	*08/09/11 17:45*	*08/10/11 02:57*	*20.0*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 17:45*	*08/10/11 02:57*	*20.0*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/10/11 02:57*	*20.0*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/09/11 17:45*	*08/10/11 02:57*	*20.0*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**30600**		2000		ug/L		08/03/11 15:45	08/11/11 18:31	20.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*92*		*60 - 150*	*08/03/11 15:45*	*08/11/11 18:31*	*20.0*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**1930**		190		ug/L		08/08/11 11:05	08/11/11 07:11	2.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*50*	*ZX*	*53 - 140*	*08/08/11 11:05*	*08/11/11 07:11*	*2.00*

Client Sample ID: VW-2 (08032011)

Date Collected: 08/03/11 16:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-19

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Benzene	**1.05**		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Diisopropyl Ether	**1.23**		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Methyl tert-Butyl Ether	**1.50**		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Bromoform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Chloroform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Isopropylbenzene	**1.07**		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 17:45	08/09/11 23:13	1.00

Client Sample ID: VW-2 (08032011)

Date Collected: 08/03/11 16:05
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-19

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1 (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Styrene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Toluene	**2.29**		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00
Xylenes, total	ND		3.00		ug/L		08/09/11 17:45	08/09/11 23:13	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*96*		*63 - 140*	*08/09/11 17:45*	*08/09/11 23:13*	*1.00*
Dibromofluoromethane	*103*		*73 - 131*	*08/09/11 17:45*	*08/09/11 23:13*	*1.00*
Toluene-d8	*94*		*80 - 120*	*08/09/11 17:45*	*08/09/11 23:13*	*1.00*
4-Bromofluorobenzene	*88*		*79 - 125*	*08/09/11 17:45*	*08/09/11 23:13*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**227**		100		ug/L		08/03/11 16:05	08/10/11 22:52	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*91*		*60 - 150*	*08/03/11 16:05*	*08/10/11 22:52*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**669**		98.0		ug/L		08/08/11 11:05	08/10/11 18:48	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*52*	*S10 Z6*	*53 - 140*	*08/08/11 11:05*	*08/10/11 18:48*	*1.00*

Client Sample ID: VW-3 (08032011)

Lab Sample ID: NUH0938-20

Date Collected: 08/03/11 10:40

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Benzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 15:15	1.00

Client Sample ID: VW-3 (08032011)

Date Collected: 08/03/11 10:40
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-20

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Toluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00
Xylenes, total	ND		3.00		ug/L		08/09/11 05:54	08/09/11 15:15	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*96*		*63 - 140*	*08/09/11 05:54*	*08/09/11 15:15*	*1.00*
Dibromofluoromethane	*102*		*73 - 131*	*08/09/11 05:54*	*08/09/11 15:15*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 05:54*	*08/09/11 15:15*	*1.00*
4-Bromofluorobenzene	*91*		*79 - 125*	*08/09/11 05:54*	*08/09/11 15:15*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/03/11 10:40	08/11/11 14:02	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*94*		*60 - 150*	*08/03/11 10:40*	*08/11/11 14:02*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	ND		98.0		ug/L		08/08/11 11:05	08/10/11 20:00	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*102*		*53 - 140*	*08/08/11 11:05*	*08/10/11 20:00*	*1.00*

Client Sample ID: MP-1 (08032011)

Date Collected: 08/03/11 16:50
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-21
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Tert-Amyl Methyl Ether	**19.2**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Diisopropyl Ether	**16.4**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Tertiary Butyl Alcohol	**424**		20.0		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
sec-Butylbenzene	**5.35**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
n-Butylbenzene	**20.4**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Isopropylbenzene	**50.6**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
p-Isopropyltoluene	**4.10**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
n-Propylbenzene	**111**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00

Client Sample ID: MP-1 (08032011)

Date Collected: 08/03/11 16:50
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-21

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
1,3,5-Trimethylbenzene	**198**		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 15:43	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*112*		*63 - 140*	*08/09/11 05:54*	*08/09/11 15:43*	*1.00*
Dibromofluoromethane	*110*		*73 - 131*	*08/09/11 05:54*	*08/09/11 15:43*	*1.00*
Toluene-d8	*83*		*80 - 120*	*08/09/11 05:54*	*08/09/11 15:43*	*1.00*
4-Bromofluorobenzene	*95*		*79 - 125*	*08/09/11 05:54*	*08/09/11 15:43*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methyl tert-Butyl Ether	**329**		10.0		ug/L		08/10/11 06:14	08/10/11 10:27	10.0
Ethylbenzene	**1490**		10.0		ug/L		08/10/11 06:14	08/10/11 10:27	10.0
Naphthalene	**360**		50.0		ug/L		08/10/11 06:14	08/10/11 10:27	10.0
1,2,4-Trimethylbenzene	**1060**		10.0		ug/L		08/10/11 06:14	08/10/11 10:27	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*106*		*63 - 140*	*08/10/11 06:14*	*08/10/11 10:27*	*10.0*
Dibromofluoromethane	*109*		*73 - 131*	*08/10/11 06:14*	*08/10/11 10:27*	*10.0*
Toluene-d8	*94*		*80 - 120*	*08/10/11 06:14*	*08/10/11 10:27*	*10.0*
4-Bromofluorobenzene	*88*		*79 - 125*	*08/10/11 06:14*	*08/10/11 10:27*	*10.0*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE2

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**7580**		100		ug/L		08/10/11 06:14	08/10/11 10:55	100
Toluene	**13400**		100		ug/L		08/10/11 06:14	08/10/11 10:55	100
Xylenes, total	**7390**		300		ug/L		08/10/11 06:14	08/10/11 10:55	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*105*		*63 - 140*	*08/10/11 06:14*	*08/10/11 10:55*	*100*
Dibromofluoromethane	*108*		*73 - 131*	*08/10/11 06:14*	*08/10/11 10:55*	*100*
Toluene-d8	*92*		*80 - 120*	*08/10/11 06:14*	*08/10/11 10:55*	*100*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/10/11 06:14*	*08/10/11 10:55*	*100*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**101000**		10000		ug/L		08/03/11 16:50	08/11/11 16:02	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*99*		*60 - 150*	*08/03/11 16:50*	*08/11/11 16:02*	*100*

Client Sample Results

Client: ARCADIS Lexington - Exxon
Project/Site: 25553 - Thurmont, MD

TestAmerica Job ID: NUH0938

Client Sample ID: MP-1 (08032011)

Date Collected: 08/03/11 16:50
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-21

Matrix: Ground Water

Method: SW846 8015B - Extractable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	1710		198		ug/L		08/08/11 11:05	08/11/11 07:35	2.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*47*	*ZX*	*53 - 140*	*08/08/11 11:05*	*08/11/11 07:35*	*2.00*

Client Sample ID: DPE-1 (08032011)

Date Collected: 08/03/11 17:45
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-22

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Tert-Amyl Methyl Ether	**24.2**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Diisopropyl Ether	**19.8**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Tertiary Butyl Alcohol	**595**		20.0		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
sec-Butylbenzene	**9.41**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
n-Butylbenzene	**37.8**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2-Dibromoethane (EDB)	**1.56**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Isopropylbenzene	**70.5**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
p-Isopropyltoluene	**6.91**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
n-Propylbenzene	**157**		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00

Client Sample ID: DPE-1 (08032011)

Date Collected: 08/03/11 17:45
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-22

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:11	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*116*		*63 - 140*	*08/09/11 05:54*	*08/09/11 16:11*	*1.00*
Dibromofluoromethane	*116*		*73 - 131*	*08/09/11 05:54*	*08/09/11 16:11*	*1.00*
Toluene-d8	*82*		*80 - 120*	*08/09/11 05:54*	*08/09/11 16:11*	*1.00*
4-Bromofluorobenzene	*95*		*79 - 125*	*08/09/11 05:54*	*08/09/11 16:11*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Methyl tert-Butyl Ether	**263**		10.0		ug/L		08/10/11 06:14	08/10/11 11:23	10.0
Ethylbenzene	**1830**		10.0		ug/L		08/10/11 06:14	08/10/11 11:23	10.0
Naphthalene	**486**		50.0		ug/L		08/10/11 06:14	08/10/11 11:23	10.0
1,3,5-Trimethylbenzene	**404**		10.0		ug/L		08/10/11 06:14	08/10/11 11:23	10.0
1,2,4-Trimethylbenzene	**1490**		10.0		ug/L		08/10/11 06:14	08/10/11 11:23	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*107*		*63 - 140*	*08/10/11 06:14*	*08/10/11 11:23*	*10.0*
Dibromofluoromethane	*110*		*73 - 131*	*08/10/11 06:14*	*08/10/11 11:23*	*10.0*
Toluene-d8	*95*		*80 - 120*	*08/10/11 06:14*	*08/10/11 11:23*	*10.0*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/10/11 06:14*	*08/10/11 11:23*	*10.0*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE2

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**5630**		100		ug/L		08/10/11 06:14	08/10/11 11:51	100
Toluene	**14300**		100		ug/L		08/10/11 06:14	08/10/11 11:51	100
Xylenes, total	**9960**		300		ug/L		08/10/11 06:14	08/10/11 11:51	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*102*		*63 - 140*	*08/10/11 06:14*	*08/10/11 11:51*	*100*
Dibromofluoromethane	*107*		*73 - 131*	*08/10/11 06:14*	*08/10/11 11:51*	*100*
Toluene-d8	*94*		*80 - 120*	*08/10/11 06:14*	*08/10/11 11:51*	*100*
4-Bromofluorobenzene	*90*		*79 - 125*	*08/10/11 06:14*	*08/10/11 11:51*	*100*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**115000**		10000		ug/L		08/03/11 17:45	08/11/11 16:32	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*100*		*60 - 150*	*08/03/11 17:45*	*08/11/11 16:32*	*100*

Client Sample Results

Client: ARCADIS Lexington - Exxon
Project/Site: 25553 - Thurmont, MD

TestAmerica Job ID: NUH0938

Client Sample ID: DPE-1 (08032011)

Date Collected: 08/03/11 17:45
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-22

Matrix: Ground Water

Method: SW846 8015B - Extractable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	2880		196		ug/L		08/08/11 11:05	08/11/11 07:59	2.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*59*		*53 - 140*	*08/08/11 11:05*	*08/11/11 07:59*	*2.00*

Client Sample ID: DUP-1 (08022011)

Date Collected: 08/02/11 12:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-23

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Tert-Amyl Methyl Ether	**2.20**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Diisopropyl Ether	**10.4**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Methyl tert-Butyl Ether	**50.0**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Tertiary Butyl Alcohol	**214**		20.0		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
sec-Butylbenzene	**6.68**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
n-Butylbenzene	**15.1**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Isopropylbenzene	**44.4**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
p-Isopropyltoluene	**3.23**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Naphthalene	**142**		5.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
n-Propylbenzene	**88.4**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00

Client Sample ID: DUP-1 (08022011)

Date Collected: 08/02/11 12:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-23

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Toluene	**42.7**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,3,5-Trimethylbenzene	**19.5**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
1,2,4-Trimethylbenzene	**34.7**		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00
Xylenes, total	**133**		3.00		ug/L		08/09/11 05:54	08/09/11 16:39	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*103*		*63 - 140*	*08/09/11 05:54*	*08/09/11 16:39*	*1.00*
Dibromofluoromethane	*102*		*73 - 131*	*08/09/11 05:54*	*08/09/11 16:39*	*1.00*
Toluene-d8	*97*		*80 - 120*	*08/09/11 05:54*	*08/09/11 16:39*	*1.00*
4-Bromofluorobenzene	*94*		*79 - 125*	*08/09/11 05:54*	*08/09/11 16:39*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**693**		10.0		ug/L		08/09/11 17:45	08/10/11 02:29	10.0
Ethylbenzene	**448**		10.0		ug/L		08/09/11 17:45	08/10/11 02:29	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*96*		*63 - 140*	*08/09/11 17:45*	*08/10/11 02:29*	*10.0*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 17:45*	*08/10/11 02:29*	*10.0*
Toluene-d8	*95*		*80 - 120*	*08/09/11 17:45*	*08/10/11 02:29*	*10.0*
4-Bromofluorobenzene	*88*		*79 - 125*	*08/09/11 17:45*	*08/10/11 02:29*	*10.0*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**42600**		2500		ug/L		08/02/11 12:00	08/11/11 20:01	25.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*96*		*60 - 150*	*08/02/11 12:00*	*08/11/11 20:01*	*25.0*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**2430**		100		ug/L		08/08/11 11:05	08/10/11 21:12	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*55*		*53 - 140*	*08/08/11 11:05*	*08/10/11 21:12*	*1.00*

Client Sample ID: DUP-2 (08032011)

Date Collected: 08/03/11 12:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-24

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Benzene	**8.72**		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Bromoform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Chloroform	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Ethylbenzene	**1.03**		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 17:45	08/09/11 23:41	1.00

Client Sample ID: DUP-2 (08032011)

Date Collected: 08/03/11 12:00
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-24

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1 (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Styrene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Toluene	**1.02**		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,3,5-Trimethylbenzene	**1.09**		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
1,2,4-Trimethylbenzene	**3.65**		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00
Xylenes, total	**8.06**		3.00		ug/L		08/09/11 17:45	08/09/11 23:41	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*98*		*63 - 140*	*08/09/11 17:45*	*08/09/11 23:41*	*1.00*
Dibromofluoromethane	*105*		*73 - 131*	*08/09/11 17:45*	*08/09/11 23:41*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 17:45*	*08/09/11 23:41*	*1.00*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/09/11 17:45*	*08/09/11 23:41*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**532**		100		ug/L		08/03/11 12:00	08/11/11 14:32	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*93*		*60 - 150*	*08/03/11 12:00*	*08/11/11 14:32*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	**99.0**		98.0		ug/L		08/08/11 11:05	08/10/11 21:36	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*86*		*53 - 140*	*08/08/11 11:05*	*08/10/11 21:36*	*1.00*

Client Sample ID: RB-1 (08022011)

Date Collected: 08/02/11 15:20
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-25
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Benzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
n-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Ethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Isopropylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 01:42	1.00

Client Sample ID: RB-1 (08022011)

Date Collected: 08/02/11 15:20
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-25

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
n-Propylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Toluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00
Xylenes, total	ND		3.00		ug/L		08/08/11 17:45	08/09/11 01:42	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*98*		*63 - 140*	*08/08/11 17:45*	*08/09/11 01:42*	*1.00*
Dibromofluoromethane	*103*		*73 - 131*	*08/08/11 17:45*	*08/09/11 01:42*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/08/11 17:45*	*08/09/11 01:42*	*1.00*
4-Bromofluorobenzene	*93*		*79 - 125*	*08/08/11 17:45*	*08/09/11 01:42*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/02/11 15:20	08/11/11 02:21	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*91*		*60 - 150*	*08/02/11 15:20*	*08/11/11 02:21*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	ND		98.0		ug/L		08/08/11 11:05	08/10/11 22:00	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*116*		*53 - 140*	*08/08/11 11:05*	*08/10/11 22:00*	*1.00*

Client Sample ID: RB-2 (08032011)

Date Collected: 08/03/11 17:55
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-26
Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Benzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
n-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Ethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Isopropylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 02:10	1.00

Client Sample ID: RB-2 (08032011)

Lab Sample ID: NUH0938-26

Date Collected: 08/03/11 17:55

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
n-Propylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Toluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00
Xylenes, total	ND		3.00		ug/L		08/08/11 17:45	08/09/11 02:10	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*97*		*63 - 140*	*08/08/11 17:45*	*08/09/11 02:10*	*1.00*
Dibromofluoromethane	*103*		*73 - 131*	*08/08/11 17:45*	*08/09/11 02:10*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/08/11 17:45*	*08/09/11 02:10*	*1.00*
4-Bromofluorobenzene	*92*		*79 - 125*	*08/08/11 17:45*	*08/09/11 02:10*	*1.00*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/03/11 17:55	08/11/11 02:51	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*93*		*60 - 150*	*08/03/11 17:55*	*08/11/11 02:51*	*1.00*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	ND		94.3		ug/L		08/08/11 11:05	08/10/11 22:24	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*98*		*53 - 140*	*08/08/11 11:05*	*08/10/11 22:24*	*1.00*

Client Sample ID: TB-1 (08032011)

Lab Sample ID: NUH0938-27

Date Collected: 08/03/11 00:01

Matrix: Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Benzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
n-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Ethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Isopropylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 00:46	1.00

Client Sample ID: TB-1 (08032011)

Lab Sample ID: NUH0938-27

Date Collected: 08/03/11 00:01

Matrix: Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
n-Propylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Toluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00
Xylenes, total	ND		3.00		ug/L		08/08/11 17:45	08/09/11 00:46	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/08/11 17:45*	*08/09/11 00:46*	*1.00*
Dibromofluoromethane	*102*		*73 - 131*	*08/08/11 17:45*	*08/09/11 00:46*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/08/11 17:45*	*08/09/11 00:46*	*1.00*
4-Bromofluorobenzene	*92*		*79 - 125*	*08/08/11 17:45*	*08/09/11 00:46*	*1.00*

Client Sample ID: TB-2 (08032011)

Lab Sample ID: NUH0938-28

Date Collected: 08/03/11 00:01

Matrix: Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND	L	50.0		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Benzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
n-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Ethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Isopropylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/09/11 01:14	1.00

Client Sample ID: TB-2 (08032011)

Lab Sample ID: NUH0938-28

Date Collected: 08/03/11 00:01

Matrix: Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Naphthalene	ND		5.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
n-Propylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Toluene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00
Xylenes, total	ND		3.00		ug/L		08/08/11 17:45	08/09/11 01:14	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*97*		*63 - 140*	*08/08/11 17:45*	*08/09/11 01:14*	*1.00*
Dibromofluoromethane	*103*		*73 - 131*	*08/08/11 17:45*	*08/09/11 01:14*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/08/11 17:45*	*08/09/11 01:14*	*1.00*
4-Bromofluorobenzene	*93*		*79 - 125*	*08/08/11 17:45*	*08/09/11 01:14*	*1.00*

Client Sample ID: MW-55I (08022011)

Lab Sample ID: NUH0938-29

Date Collected: 08/02/11 15:40

Matrix: Ground Water

Date Received: 08/06/11 08:25

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Diisopropyl Ether	**66.4**		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Methyl tert-Butyl Ether	**19.2**		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Tertiary Butyl Alcohol	**95.4**		20.0		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
sec-Butylbenzene	**10.5**		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
n-Butylbenzene	**33.0**		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Isopropylbenzene	**101**		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
p-Isopropyltoluene	**7.59**		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
4-Methyl-2-pentanone	**70.3**		10.0		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
n-Propylbenzene	**181**		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00

Client Sample ID: MW-55I (08022011)

Date Collected: 08/02/11 15:40
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-29

Matrix: Ground Water

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 17:36	1.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*101*		*63 - 140*	*08/09/11 05:54*	*08/09/11 17:36*	*1.00*
Dibromofluoromethane	*102*		*73 - 131*	*08/09/11 05:54*	*08/09/11 17:36*	*1.00*
Toluene-d8	*95*		*80 - 120*	*08/09/11 05:54*	*08/09/11 17:36*	*1.00*
4-Bromofluorobenzene	*97*		*79 - 125*	*08/09/11 05:54*	*08/09/11 17:36*	*1.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Ethylbenzene	**1130**		10.0		ug/L		08/10/11 06:14	08/10/11 12:19	10.0
Naphthalene	**464**		50.0		ug/L		08/10/11 06:14	08/10/11 12:19	10.0
Toluene	**1540**		10.0		ug/L		08/10/11 06:14	08/10/11 12:19	10.0
1,3,5-Trimethylbenzene	**351**		10.0		ug/L		08/10/11 06:14	08/10/11 12:19	10.0
1,2,4-Trimethylbenzene	**1160**		10.0		ug/L		08/10/11 06:14	08/10/11 12:19	10.0
Xylenes, total	**3780**		30.0		ug/L		08/10/11 06:14	08/10/11 12:19	10.0

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*106*		*63 - 140*	*08/10/11 06:14*	*08/10/11 12:19*	*10.0*
Dibromofluoromethane	*110*		*73 - 131*	*08/10/11 06:14*	*08/10/11 12:19*	*10.0*
Toluene-d8	*93*		*80 - 120*	*08/10/11 06:14*	*08/10/11 12:19*	*10.0*
4-Bromofluorobenzene	*86*		*79 - 125*	*08/10/11 06:14*	*08/10/11 12:19*	*10.0*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B - RE2

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Benzene	**3360**		100		ug/L		08/10/11 06:14	08/10/11 12:47	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*103*		*63 - 140*	*08/10/11 06:14*	*08/10/11 12:47*	*100*
Dibromofluoromethane	*107*		*73 - 131*	*08/10/11 06:14*	*08/10/11 12:47*	*100*
Toluene-d8	*93*		*80 - 120*	*08/10/11 06:14*	*08/10/11 12:47*	*100*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/10/11 06:14*	*08/10/11 12:47*	*100*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	**49000**		10000		ug/L		08/02/11 15:40	08/11/11 17:01	100

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*97*		*60 - 150*	*08/02/11 15:40*	*08/11/11 17:01*	*100*

Client Sample Results

Client: ARCADIS Lexington - Exxon
Project/Site: 25553 - Thurmont, MD

TestAmerica Job ID: NUH0938

Client Sample ID: MW-55I (08022011)

Date Collected: 08/02/11 15:40
Date Received: 08/06/11 08:25

Lab Sample ID: NUH0938-29

Matrix: Ground Water

Method: SW846 8015B - Extractable Petroleum Hydrocarbons - RE1

Analyte	Result	Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	2470		198		ug/L		08/08/11 11:05	08/11/11 08:23	2.00

Surrogate	*% Recovery*	*Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*69*		*53 - 140*	*08/08/11 11:05*	*08/11/11 08:23*	*2.00*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B

Lab Sample ID: 11H1622-BLK1
Matrix: Water
Analysis Batch: U014132

Client Sample ID: Method Blank
Prep Type: Total
Prep Batch: 11H1622_P

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Benzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Bromoform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Chloroform	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1622-BLK1
Matrix: Water
Analysis Batch: U014132

Client Sample ID: Method Blank
Prep Type: Total
Prep Batch: 11H1622_P

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Naphthalene	ND		5.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Styrene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Toluene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00
Xylenes, total	ND		3.00		ug/L		08/09/11 05:54	08/09/11 08:42	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/09/11 05:54*	*08/09/11 08:42*	*1.00*
Dibromofluoromethane	*103*		*73 - 131*	*08/09/11 05:54*	*08/09/11 08:42*	*1.00*
Toluene-d8	*96*		*80 - 120*	*08/09/11 05:54*	*08/09/11 08:42*	*1.00*
4-Bromofluorobenzene	*91*		*79 - 125*	*08/09/11 05:54*	*08/09/11 08:42*	*1.00*

Lab Sample ID: 11H1622-BS1
Matrix: Water
Analysis Batch: U014132

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H1622_P

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	250	336		ug/L		134	56 - 150
Tert-Amyl Methyl Ether	50.0	43.5		ug/L		87	70 - 133
Benzene	50.0	47.7		ug/L		95	80 - 121
Bromobenzene	50.0	43.2		ug/L		86	72 - 130
Bromochloromethane	50.0	51.5		ug/L		103	73 - 137
Ethyl tert-Butyl Ether	50.0	42.0		ug/L		84	68 - 138
Bromodichloromethane	50.0	49.6		ug/L		99	75 - 131
Diisopropyl Ether	50.0	40.7		ug/L		81	63 - 136
Methyl tert-Butyl Ether	50.0	45.7		ug/L		91	76 - 120
Bromoform	50.0	47.7		ug/L		95	65 - 140
Tertiary Butyl Alcohol	500	513		ug/L		103	60 - 140
Bromomethane	50.0	41.4		ug/L		83	50 - 150
2-Butanone	250	285		ug/L		114	70 - 144
sec-Butylbenzene	50.0	46.9		ug/L		94	72 - 140
n-Butylbenzene	50.0	44.7		ug/L		89	68 - 140
tert-Butylbenzene	50.0	46.9		ug/L		94	76 - 135
Carbon disulfide	50.0	43.5		ug/L		87	74 - 137
Carbon Tetrachloride	50.0	54.4		ug/L		109	71 - 137

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1622-BS1
Matrix: Water
Analysis Batch: U014132

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H1622_P

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Chlorobenzene	50.0	48.8		ug/L		98	80 - 121
Chlorodibromomethane	50.0	55.1		ug/L		110	68 - 137
Chloroethane	50.0	40.3		ug/L		81	50 - 146
Chloroform	50.0	49.7		ug/L		99	73 - 131
Chloromethane	50.0	29.8		ug/L		60	30 - 132
2-Chlorotoluene	50.0	43.6		ug/L		87	74 - 135
4-Chlorotoluene	50.0	45.3		ug/L		91	74 - 132
1,2-Dibromo-3-chloropropane	50.0	47.5		ug/L		95	56 - 145
1,2-Dibromoethane (EDB)	50.0	49.2		ug/L		98	80 - 135
Dibromomethane	50.0	50.6		ug/L		101	78 - 133
1,4-Dichlorobenzene	50.0	49.2		ug/L		98	80 - 120
1,3-Dichlorobenzene	50.0	48.7		ug/L		97	80 - 128
1,2-Dichlorobenzene	50.0	49.1		ug/L		98	80 - 125
Dichlorodifluoromethane	50.0	27.8		ug/L		56	30 - 132
1,1-Dichloroethane	50.0	47.5		ug/L		95	75 - 125
1,2-Dichloroethane	50.0	48.3		ug/L		97	70 - 134
cis-1,2-Dichloroethene	50.0	45.6		ug/L		91	71 - 132
1,1-Dichloroethene	50.0	49.2		ug/L		98	73 - 125
trans-1,2-Dichloroethene	50.0	47.3		ug/L		95	77 - 125
1,3-Dichloropropane	50.0	46.0		ug/L		92	76 - 125
1,2-Dichloropropane	50.0	43.3		ug/L		87	72 - 120
2,2-Dichloropropane	50.0	39.0		ug/L		78	50 - 150
cis-1,3-Dichloropropene	50.0	47.9		ug/L		96	70 - 140
trans-1,3-Dichloropropene	50.0	46.0		ug/L		92	62 - 139
1,1-Dichloropropene	50.0	48.0		ug/L		96	78 - 126
Ethylbenzene	50.0	46.7		ug/L		93	78 - 133
Hexachlorobutadiene	50.0	47.2		ug/L		94	70 - 150
2-Hexanone	250	258		ug/L		103	60 - 150
Isopropylbenzene	50.0	52.6		ug/L		105	69 - 120
p-Isopropyltoluene	50.0	46.3		ug/L		93	72 - 134
Methylene Chloride	50.0	49.3		ug/L		99	80 - 133
4-Methyl-2-pentanone	250	225		ug/L		90	62 - 146
Naphthalene	50.0	48.9		ug/L		98	71 - 139
n-Propylbenzene	50.0	44.7		ug/L		89	70 - 143
Styrene	50.0	49.0		ug/L		98	80 - 136
1,1,1,2-Tetrachloroethane	50.0	52.0		ug/L		104	80 - 130
1,1,2,2-Tetrachloroethane	50.0	43.2		ug/L		86	73 - 131
Tetrachloroethene	50.0	51.8		ug/L		104	77 - 131
Toluene	50.0	47.3		ug/L		95	78 - 125
1,2,3-Trichlorobenzene	50.0	51.4		ug/L		103	71 - 138
1,2,4-Trichlorobenzene	50.0	51.6		ug/L		103	74 - 136
1,1,2-Trichloroethane	50.0	49.1		ug/L		98	80 - 123
1,1,1-Trichloroethane	50.0	51.5		ug/L		103	75 - 137
Trichloroethene	50.0	54.8		ug/L		110	74 - 139
Trichlorofluoromethane	50.0	42.4		ug/L		85	60 - 133
1,2,3-Trichloropropane	50.0	44.6		ug/L		89	64 - 127
1,3,5-Trimethylbenzene	50.0	46.1		ug/L		92	75 - 134
1,2,4-Trimethylbenzene	50.0	45.1		ug/L		90	77 - 134
Vinyl chloride	50.0	36.3		ug/L		73	60 - 122
Xylenes, total	150	143		ug/L		95	78 - 134

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1622-BS1
Matrix: Water
Analysis Batch: U014132

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H1622_P

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
1,2-Dichloroethane-d4	*93*		*63 - 140*
Dibromofluoromethane	*105*		*73 - 131*
Toluene-d8	*96*		*80 - 120*
4-Bromofluorobenzene	*91*		*79 - 125*

Lab Sample ID: 11H1622-BSD1
Matrix: Water
Analysis Batch: U014132

Client Sample ID: Lab Control Sample Dup
Prep Type: Total
Prep Batch: 11H1622_P

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	250	315		ug/L		126	56 - 150	7	31
Tert-Amyl Methyl Ether	50.0	42.5		ug/L		85	70 - 133	2	16
Benzene	50.0	46.7		ug/L		93	80 - 121	2	12
Bromobenzene	50.0	42.2		ug/L		84	72 - 130	2	23
Bromochloromethane	50.0	51.0		ug/L		102	73 - 137	1	32
Ethyl tert-Butyl Ether	50.0	41.2		ug/L		82	68 - 138	2	16
Bromodichloromethane	50.0	48.5		ug/L		97	75 - 131	2	13
Diisopropyl Ether	50.0	39.8		ug/L		80	63 - 136	2	32
Methyl tert-Butyl Ether	50.0	44.4		ug/L		89	76 - 120	3	32
Bromoform	50.0	47.1		ug/L		94	65 - 140	1	18
Tertiary Butyl Alcohol	500	511		ug/L		102	60 - 140	0.3	30
Bromomethane	50.0	40.8		ug/L		82	50 - 150	1	50
2-Butanone	250	273		ug/L		109	70 - 144	4	37
sec-Butylbenzene	50.0	46.2		ug/L		92	72 - 140	1	21
n-Butylbenzene	50.0	43.8		ug/L		88	68 - 140	2	11
tert-Butylbenzene	50.0	47.1		ug/L		94	76 - 135	0.3	20
Carbon disulfide	50.0	42.1		ug/L		84	74 - 137	3	28
Carbon Tetrachloride	50.0	53.3		ug/L		107	71 - 137	2	26
Chlorobenzene	50.0	48.2		ug/L		96	80 - 121	1	11
Chlorodibromomethane	50.0	54.0		ug/L		108	68 - 137	2	16
Chloroethane	50.0	39.3		ug/L		79	50 - 146	3	35
Chloroform	50.0	48.7		ug/L		97	73 - 131	2	32
Chloromethane	50.0	29.2		ug/L		58	30 - 132	2	34
2-Chlorotoluene	50.0	43.2		ug/L		86	74 - 135	0.9	22
4-Chlorotoluene	50.0	45.0		ug/L		90	74 - 132	0.5	22
1,2-Dibromo-3-chloropropane	50.0	44.4		ug/L		89	56 - 145	7	21
1,2-Dibromoethane (EDB)	50.0	49.1		ug/L		98	80 - 135	0.3	10
Dibromomethane	50.0	49.1		ug/L		98	78 - 133	3	11
1,4-Dichlorobenzene	50.0	48.5		ug/L		97	80 - 120	1	10
1,3-Dichlorobenzene	50.0	47.9		ug/L		96	80 - 128	2	18
1,2-Dichlorobenzene	50.0	48.5		ug/L		97	80 - 125	1	11
Dichlorodifluoromethane	50.0	26.8		ug/L		54	30 - 132	4	32
1,1-Dichloroethane	50.0	46.2		ug/L		92	75 - 125	3	34
1,2-Dichloroethane	50.0	46.9		ug/L		94	70 - 134	3	25
cis-1,2-Dichloroethene	50.0	44.6		ug/L		89	71 - 132	2	32
1,1-Dichloroethene	50.0	48.4		ug/L		97	73 - 125	2	31
trans-1,2-Dichloroethene	50.0	46.9		ug/L		94	77 - 125	0.8	32
1,3-Dichloropropane	50.0	45.3		ug/L		91	76 - 125	2	20
1,2-Dichloropropane	50.0	42.3		ug/L		85	72 - 120	2	11

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1622-BSD1 | **Client Sample ID: Lab Control Sample Dup**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014132 | **Prep Batch: 11H1622_P**

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
2,2-Dichloropropane	50.0	37.0		ug/L		74	50 - 150	5	11
cis-1,3-Dichloropropene	50.0	46.8		ug/L		94	70 - 140	2	35
trans-1,3-Dichloropropene	50.0	44.9		ug/L		90	62 - 139	2	26
1,1-Dichloropropene	50.0	47.3		ug/L		95	78 - 126	1	18
Ethylbenzene	50.0	46.3		ug/L		93	78 - 133	0.8	12
Hexachlorobutadiene	50.0	47.6		ug/L		95	70 - 150	0.7	21
2-Hexanone	250	241		ug/L		96	60 - 150	7	20
Isopropylbenzene	50.0	52.2		ug/L		104	69 - 120	0.8	15
p-Isopropyltoluene	50.0	45.8		ug/L		92	72 - 134	1	18
Methylene Chloride	50.0	48.1		ug/L		96	80 - 133	3	36
4-Methyl-2-pentanone	250	219		ug/L		88	62 - 146	3	35
Naphthalene	50.0	48.2		ug/L		96	71 - 139	2	30
n-Propylbenzene	50.0	44.3		ug/L		89	70 - 143	1	23
Styrene	50.0	48.5		ug/L		97	80 - 136	1	29
1,1,1,2-Tetrachloroethane	50.0	51.6		ug/L		103	80 - 130	0.8	11
1,1,2,2-Tetrachloroethane	50.0	42.8		ug/L		86	73 - 131	0.9	28
Tetrachloroethene	50.0	51.3		ug/L		103	77 - 131	0.9	16
Toluene	50.0	46.8		ug/L		94	78 - 125	1	35
1,2,3-Trichlorobenzene	50.0	51.4		ug/L		103	71 - 138	0.08	28
1,2,4-Trichlorobenzene	50.0	50.7		ug/L		101	74 - 136	2	23
1,1,2-Trichloroethane	50.0	48.4		ug/L		97	80 - 123	1	21
1,1,1-Trichloroethane	50.0	50.8		ug/L		102	75 - 137	1	29
Trichloroethene	50.0	53.4		ug/L		107	74 - 139	3	11
Trichlorofluoromethane	50.0	43.3		ug/L		87	60 - 133	2	33
1,2,3-Trichloropropane	50.0	43.6		ug/L		87	64 - 127	2	25
1,3,5-Trimethylbenzene	50.0	45.6		ug/L		91	75 - 134	1	21
1,2,4-Trimethylbenzene	50.0	44.5		ug/L		89	77 - 134	1	20
Vinyl chloride	50.0	36.4		ug/L		73	60 - 122	0.1	32
Xylenes, total	150	142		ug/L		94	78 - 134	1	18

Surrogate	*LCS Dup % Recovery*	*LCS Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*94*		*63 - 140*
Dibromofluoromethane	*104*		*73 - 131*
Toluene-d8	*97*		*80 - 120*
4-Bromofluorobenzene	*91*		*79 - 125*

Lab Sample ID: 11H1622-MS1 | **Client Sample ID: MW-46 (08022011)**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014132 | **Prep Batch: 11H1622_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	ND		250	191		ug/L		77	56 - 150
Tert-Amyl Methyl Ether	ND		50.0	39.5		ug/L		79	69 - 139
Benzene	ND		50.0	44.5		ug/L		89	65 - 151
Bromobenzene	ND		50.0	36.6		ug/L		73	69 - 142
Bromochloromethane	ND		50.0	50.6		ug/L		101	64 - 154
Ethyl tert-Butyl Ether	ND		50.0	37.6		ug/L		75	68 - 139
Bromodichloromethane	ND		50.0	46.1		ug/L		92	75 - 138
Diisopropyl Ether	ND		50.0	35.6		ug/L		71	59 - 145

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1622-MS1 | **Client Sample ID: MW-46 (08022011)**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014132 | **Prep Batch: 11H1622_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Methyl tert-Butyl Ether	ND		50.0	41.8		ug/L		84	56 - 152
Bromoform	ND		50.0	42.7		ug/L		85	55 - 153
Tertiary Butyl Alcohol	ND		500	503		ug/L		101	14 - 200
Bromomethane	ND		50.0	34.1		ug/L		68	13 - 176
2-Butanone	ND		250	209		ug/L		84	45 - 164
sec-Butylbenzene	ND		50.0	38.7		ug/L		77	68 - 159
n-Butylbenzene	ND		50.0	36.2		ug/L		72	67 - 151
tert-Butylbenzene	ND		50.0	40.2		ug/L		80	73 - 153
Carbon disulfide	ND		50.0	41.1		ug/L		82	33 - 187
Carbon Tetrachloride	ND		50.0	53.5		ug/L		107	64 - 157
Chlorobenzene	ND		50.0	44.5		ug/L		89	78 - 136
Chlorodibromomethane	ND		50.0	49.6		ug/L		99	64 - 145
Chloroethane	ND		50.0	38.1		ug/L		76	48 - 159
Chloroform	ND		50.0	47.3		ug/L		95	72 - 145
Chloromethane	ND		50.0	29.4		ug/L		59	10 - 194
2-Chlorotoluene	ND		50.0	37.0		ug/L		74	66 - 155
4-Chlorotoluene	ND		50.0	38.8		ug/L		78	69 - 149
1,2-Dibromo-3-chloropropane	ND		50.0	40.7		ug/L		81	49 - 162
1,2-Dibromoethane (EDB)	ND		50.0	44.0		ug/L		88	70 - 152
Dibromomethane	ND		50.0	45.8		ug/L		92	75 - 141
1,4-Dichlorobenzene	ND		50.0	42.1		ug/L		84	75 - 135
1,3-Dichlorobenzene	ND		50.0	41.8		ug/L		84	72 - 146
1,2-Dichlorobenzene	ND		50.0	42.5		ug/L		85	80 - 136
Dichlorodifluoromethane	ND		50.0	31.2		ug/L		62	23 - 159
1,1-Dichloroethane	ND		50.0	44.1		ug/L		88	64 - 154
1,2-Dichloroethane	ND		50.0	44.3		ug/L		89	72 - 137
cis-1,2-Dichloroethene	ND		50.0	42.8		ug/L		86	57 - 154
1,1-Dichloroethene	ND		50.0	47.5		ug/L		95	34 - 151
trans-1,2-Dichloroethene	ND		50.0	44.8		ug/L		90	57 - 157
1,3-Dichloropropane	ND		50.0	40.5		ug/L		81	71 - 137
1,2-Dichloropropane	ND		50.0	38.8		ug/L		78	71 - 139
2,2-Dichloropropane	ND		50.0	42.2		ug/L		84	10 - 198
cis-1,3-Dichloropropene	ND		50.0	42.7		ug/L		85	56 - 156
trans-1,3-Dichloropropene	ND		50.0	41.1		ug/L		82	47 - 157
1,1-Dichloropropene	ND		50.0	45.5		ug/L		91	70 - 155
Ethylbenzene	ND		50.0	41.6		ug/L		83	68 - 157
Hexachlorobutadiene	ND		50.0	42.8		ug/L		86	47 - 173
2-Hexanone	ND		250	195		ug/L		78	57 - 154
Isopropylbenzene	ND		50.0	46.4		ug/L		93	69 - 139
p-Isopropyltoluene	ND		50.0	38.5		ug/L		77	69 - 151
Methylene Chloride	ND		50.0	45.3		ug/L		91	71 - 136
4-Methyl-2-pentanone	ND		250	193		ug/L		77	62 - 159
Naphthalene	ND		50.0	41.4		ug/L		83	56 - 161
n-Propylbenzene	ND		50.0	37.6		ug/L		75	61 - 167
Styrene	ND		50.0	42.6		ug/L		85	69 - 150
1,1,1,2-Tetrachloroethane	ND		50.0	48.4		ug/L		97	80 - 140
1,1,2,2-Tetrachloroethane	ND		50.0	37.2	M8	ug/L		74	76 - 141
Tetrachloroethene	ND		50.0	48.0		ug/L		96	63 - 155
Toluene	ND		50.0	43.2		ug/L		86	61 - 153
1,2,3-Trichlorobenzene	ND		50.0	44.5		ug/L		89	57 - 155

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1622-MS1 — **Client Sample ID: MW-46 (08022011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014132 — **Prep Batch: 11H1622_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
1,2,4-Trichlorobenzene	ND		50.0	44.6		ug/L		89	64 - 147
1,1,2-Trichloroethane	ND		50.0	43.6		ug/L		87	74 - 138
1,1,1-Trichloroethane	ND		50.0	50.7		ug/L		101	78 - 153
Trichloroethene	ND		50.0	51.6		ug/L		103	74 - 139
Trichlorofluoromethane	ND		50.0	42.7		ug/L		85	53 - 149
1,2,3-Trichloropropane	ND		50.0	36.9		ug/L		74	49 - 148
1,3,5-Trimethylbenzene	ND		50.0	39.2		ug/L		78	67 - 151
1,2,4-Trimethylbenzene	ND		50.0	38.2		ug/L		76	69 - 150
Vinyl chloride	ND		50.0	37.0		ug/L		74	53 - 137
Xylenes, total	ND		150	128		ug/L		85	68 - 158

Surrogate	*Matrix Spike % Recovery*	*Matrix Spike Qualifier*	*Limits*
1,2-Dichloroethane-d4	*94*		*63 - 140*
Dibromofluoromethane	*107*		*73 - 131*
Toluene-d8	*94*		*80 - 120*
4-Bromofluorobenzene	*89*		*79 - 125*

Lab Sample ID: 11H1622-MSD1 — **Client Sample ID: MW-46 (08022011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014132 — **Prep Batch: 11H1622_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	ND		250	208		ug/L		83	56 - 150	8	31
Tert-Amyl Methyl Ether	ND		50.0	44.1		ug/L		88	69 - 139	11	16
Benzene	ND		50.0	48.4		ug/L		97	65 - 151	8	12
Bromobenzene	ND		50.0	40.7		ug/L		81	69 - 142	11	23
Bromochloromethane	ND		50.0	56.4		ug/L		113	64 - 154	11	32
Ethyl tert-Butyl Ether	ND		50.0	42.1		ug/L		84	68 - 139	11	16
Bromodichloromethane	ND		50.0	51.1		ug/L		102	75 - 138	10	13
Diisopropyl Ether	ND		50.0	39.2		ug/L		78	59 - 145	10	32
Methyl tert-Butyl Ether	ND		50.0	47.0		ug/L		94	56 - 152	12	32
Bromoform	ND		50.0	49.4		ug/L		99	55 - 153	15	18
Tertiary Butyl Alcohol	ND		500	568		ug/L		114	14 - 200	12	30
Bromomethane	ND		50.0	40.0		ug/L		80	13 - 176	16	50
2-Butanone	ND		250	243		ug/L		97	45 - 164	15	37
sec-Butylbenzene	ND		50.0	43.5		ug/L		87	68 - 159	12	21
n-Butylbenzene	ND		50.0	41.2	R2	ug/L		82	67 - 151	13	11
tert-Butylbenzene	ND		50.0	44.4		ug/L		89	73 - 153	10	20
Carbon disulfide	ND		50.0	43.7		ug/L		87	33 - 187	6	28
Carbon Tetrachloride	ND		50.0	58.2		ug/L		116	64 - 157	8	26
Chlorobenzene	ND		50.0	49.0		ug/L		98	78 - 136	10	11
Chlorodibromomethane	ND		50.0	55.8		ug/L		112	64 - 145	12	16
Chloroethane	ND		50.0	40.5		ug/L		81	48 - 159	6	35
Chloroform	ND		50.0	51.6		ug/L		103	72 - 145	9	32
Chloromethane	ND		50.0	31.7		ug/L		63	10 - 194	7	34
2-Chlorotoluene	ND		50.0	40.8		ug/L		82	66 - 155	10	22
4-Chlorotoluene	ND		50.0	42.6		ug/L		85	69 - 149	9	22
1,2-Dibromo-3-chloropropane	ND		50.0	47.6		ug/L		95	49 - 162	16	21
1,2-Dibromoethane (EDB)	ND		50.0	50.1	R2	ug/L		100	70 - 152	13	10

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1622-MSD1 | **Client Sample ID: MW-46 (08022011)**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014132 | **Prep Batch: 11H1622_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Dibromomethane	ND		50.0	51.0		ug/L		102	75 - 141	11	11
1,4-Dichlorobenzene	ND		50.0	46.9	R2	ug/L		94	75 - 135	11	10
1,3-Dichlorobenzene	ND		50.0	46.3		ug/L		93	72 - 146	10	18
1,2-Dichlorobenzene	ND		50.0	47.3		ug/L		95	80 - 136	11	11
Dichlorodifluoromethane	ND		50.0	35.2		ug/L		70	23 - 159	12	32
1,1-Dichloroethane	ND		50.0	47.5		ug/L		95	64 - 154	7	34
1,2-Dichloroethane	ND		50.0	49.2		ug/L		98	72 - 137	10	25
cis-1,2-Dichloroethene	ND		50.0	46.2		ug/L		92	57 - 154	8	32
1,1-Dichloroethene	ND		50.0	51.2		ug/L		102	34 - 151	7	31
trans-1,2-Dichloroethene	ND		50.0	48.0		ug/L		96	57 - 157	7	32
1,3-Dichloropropane	ND		50.0	45.3		ug/L		91	71 - 137	11	20
1,2-Dichloropropane	ND		50.0	43.1		ug/L		86	71 - 139	11	11
2,2-Dichloropropane	ND		50.0	44.8		ug/L		90	10 - 198	6	11
cis-1,3-Dichloropropene	ND		50.0	47.6		ug/L		95	56 - 156	11	35
trans-1,3-Dichloropropene	ND		50.0	46.0		ug/L		92	47 - 157	11	26
1,1-Dichloropropene	ND		50.0	49.0		ug/L		98	70 - 155	7	18
Ethylbenzene	ND		50.0	46.1		ug/L		92	68 - 157	10	12
Hexachlorobutadiene	ND		50.0	47.8		ug/L		96	47 - 173	11	21
2-Hexanone	ND		250	229		ug/L		92	57 - 154	16	20
Isopropylbenzene	ND		50.0	51.8		ug/L		104	69 - 139	11	15
p-Isopropyltoluene	ND		50.0	43.2		ug/L		86	69 - 151	11	18
Methylene Chloride	ND		50.0	49.9		ug/L		100	71 - 136	9	36
4-Methyl-2-pentanone	ND		250	224		ug/L		90	62 - 159	15	35
Naphthalene	ND		50.0	47.4		ug/L		95	56 - 161	13	30
n-Propylbenzene	ND		50.0	41.9		ug/L		84	61 - 167	11	23
Styrene	ND		50.0	47.2		ug/L		94	69 - 150	10	29
1,1,1,2-Tetrachloroethane	ND		50.0	53.6		ug/L		107	80 - 140	10	11
1,1,2,2-Tetrachloroethane	ND		50.0	42.4		ug/L		85	76 - 141	13	28
Tetrachloroethene	ND		50.0	53.1		ug/L		106	63 - 155	10	16
Toluene	ND		50.0	47.0		ug/L		94	61 - 153	8	35
1,2,3-Trichlorobenzene	ND		50.0	50.3		ug/L		101	57 - 155	12	28
1,2,4-Trichlorobenzene	ND		50.0	49.5		ug/L		99	64 - 147	10	23
1,1,2-Trichloroethane	ND		50.0	48.9		ug/L		98	74 - 138	12	21
1,1,1-Trichloroethane	ND		50.0	55.3		ug/L		111	78 - 153	9	29
Trichloroethene	ND		50.0	55.9		ug/L		112	74 - 139	8	11
Trichlorofluoromethane	ND		50.0	47.8		ug/L		96	53 - 149	11	33
1,2,3-Trichloropropane	ND		50.0	42.2		ug/L		84	49 - 148	13	25
1,3,5-Trimethylbenzene	ND		50.0	43.4		ug/L		87	67 - 151	10	21
1,2,4-Trimethylbenzene	ND		50.0	42.4		ug/L		85	69 - 150	10	20
Vinyl chloride	ND		50.0	39.8		ug/L		80	53 - 137	7	32
Xylenes, total	ND		150	140		ug/L		94	68 - 158	10	18

Surrogate	*Matrix Spike Dup % Recovery*	*Matrix Spike Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*94*		*63 - 140*
Dibromofluoromethane	*107*		*73 - 131*
Toluene-d8	*95*		*80 - 120*
4-Bromofluorobenzene	*89*		*79 - 125*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-BLK1
Matrix: Water
Analysis Batch: U014048

Client Sample ID: Method Blank
Prep Type: Total
Prep Batch: 11H1881_P

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Benzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Bromobenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Bromochloromethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Bromodichloromethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Bromoform	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Bromomethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
2-Butanone	ND		50.0		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
n-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Carbon disulfide	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Chlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Chloroethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Chloroform	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Chloromethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Dibromomethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Ethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
2-Hexanone	ND		50.0		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Isopropylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Methylene Chloride	ND		5.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-BLK1
Matrix: Water
Analysis Batch: U014048

Client Sample ID: Method Blank
Prep Type: Total
Prep Batch: 11H1881_P

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
4-Methyl-2-pentanone	ND		10.0		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Naphthalene	ND		5.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
n-Propylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Styrene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Tetrachloroethene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Toluene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Trichloroethene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Vinyl chloride	ND		1.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00
Xylenes, total	ND		3.00		ug/L		08/08/11 17:45	08/08/11 20:33	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*110*		*63 - 140*	*08/08/11 17:45*	*08/08/11 20:33*	*1.00*
Dibromofluoromethane	*107*		*73 - 131*	*08/08/11 17:45*	*08/08/11 20:33*	*1.00*
Toluene-d8	*94*		*80 - 120*	*08/08/11 17:45*	*08/08/11 20:33*	*1.00*
4-Bromofluorobenzene	*93*		*79 - 125*	*08/08/11 17:45*	*08/08/11 20:33*	*1.00*

Lab Sample ID: 11H1881-BS1
Matrix: Water
Analysis Batch: U014048

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H1881_P

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	250	407	L1	ug/L		163	56 - 150
Tert-Amyl Methyl Ether	50.0	44.8		ug/L		90	70 - 133
Benzene	50.0	48.3		ug/L		97	80 - 121
Bromobenzene	50.0	44.0		ug/L		88	72 - 130
Bromochloromethane	50.0	51.9		ug/L		104	73 - 137
Ethyl tert-Butyl Ether	50.0	43.7		ug/L		87	68 - 138
Bromodichloromethane	50.0	52.3		ug/L		105	75 - 131
Diisopropyl Ether	50.0	43.1		ug/L		86	63 - 136
Methyl tert-Butyl Ether	50.0	48.2		ug/L		96	76 - 120
Bromoform	50.0	50.5		ug/L		101	65 - 140
Tertiary Butyl Alcohol	500	567		ug/L		113	60 - 140
Bromomethane	50.0	44.7		ug/L		89	50 - 150
2-Butanone	250	305		ug/L		122	70 - 144
sec-Butylbenzene	50.0	47.8		ug/L		96	72 - 140
n-Butylbenzene	50.0	50.4		ug/L		101	68 - 140
tert-Butylbenzene	50.0	48.0		ug/L		96	76 - 135
Carbon disulfide	50.0	45.2		ug/L		90	74 - 137
Carbon Tetrachloride	50.0	58.7		ug/L		117	71 - 137

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-BS1
Matrix: Water
Analysis Batch: U014048

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H1881_P

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Chlorobenzene	50.0	49.3		ug/L		99	80 - 121
Chlorodibromomethane	50.0	57.1		ug/L		114	68 - 137
Chloroethane	50.0	44.3		ug/L		89	50 - 146
Chloroform	50.0	52.0		ug/L		104	73 - 131
Chloromethane	50.0	33.6		ug/L		67	30 - 132
2-Chlorotoluene	50.0	44.2		ug/L		88	74 - 135
4-Chlorotoluene	50.0	46.8		ug/L		94	74 - 132
1,2-Dibromo-3-chloropropane	50.0	51.6		ug/L		103	56 - 145
1,2-Dibromoethane (EDB)	50.0	51.0		ug/L		102	80 - 135
Dibromomethane	50.0	52.2		ug/L		104	78 - 133
1,4-Dichlorobenzene	50.0	50.0		ug/L		100	80 - 120
1,3-Dichlorobenzene	50.0	49.3		ug/L		99	80 - 128
1,2-Dichlorobenzene	50.0	53.4		ug/L		107	80 - 125
Dichlorodifluoromethane	50.0	31.2		ug/L		62	30 - 132
1,1-Dichloroethane	50.0	50.0		ug/L		100	75 - 125
1,2-Dichloroethane	50.0	52.7		ug/L		105	70 - 134
cis-1,2-Dichloroethene	50.0	49.0		ug/L		98	71 - 132
1,1-Dichloroethene	50.0	51.0		ug/L		102	73 - 125
trans-1,2-Dichloroethene	50.0	50.4		ug/L		101	77 - 125
1,3-Dichloropropane	50.0	48.0		ug/L		96	76 - 125
1,2-Dichloropropane	50.0	44.7		ug/L		89	72 - 120
2,2-Dichloropropane	50.0	53.3		ug/L		107	50 - 150
cis-1,3-Dichloropropene	50.0	51.0		ug/L		102	70 - 140
trans-1,3-Dichloropropene	50.0	50.9		ug/L		102	62 - 139
1,1-Dichloropropene	50.0	50.6		ug/L		101	78 - 126
Ethylbenzene	50.0	47.8		ug/L		96	78 - 133
Hexachlorobutadiene	50.0	50.4		ug/L		101	70 - 150
2-Hexanone	250	281		ug/L		112	60 - 150
Isopropylbenzene	50.0	53.9		ug/L		108	69 - 120
p-Isopropyltoluene	50.0	47.9		ug/L		96	72 - 134
Methylene Chloride	50.0	50.7		ug/L		101	80 - 133
4-Methyl-2-pentanone	250	240		ug/L		96	62 - 146
Naphthalene	50.0	49.8		ug/L		100	71 - 139
n-Propylbenzene	50.0	45.3		ug/L		91	70 - 143
Styrene	50.0	49.8		ug/L		100	80 - 136
1,1,1,2-Tetrachloroethane	50.0	53.7		ug/L		107	80 - 130
1,1,2,2-Tetrachloroethane	50.0	46.4		ug/L		93	73 - 131
Tetrachloroethene	50.0	53.1		ug/L		106	77 - 131
Toluene	50.0	47.2		ug/L		94	78 - 125
1,2,3-Trichlorobenzene	50.0	53.3		ug/L		107	71 - 138
1,2,4-Trichlorobenzene	50.0	52.6		ug/L		105	74 - 136
1,1,2-Trichloroethane	50.0	50.3		ug/L		101	80 - 123
1,1,1-Trichloroethane	50.0	54.9		ug/L		110	75 - 137
Trichloroethene	50.0	53.2		ug/L		106	74 - 139
Trichlorofluoromethane	50.0	49.7		ug/L		99	60 - 133
1,2,3-Trichloropropane	50.0	48.2		ug/L		96	64 - 127
1,3,5-Trimethylbenzene	50.0	46.7		ug/L		93	75 - 134
1,2,4-Trimethylbenzene	50.0	45.6		ug/L		91	77 - 134
Vinyl chloride	50.0	40.6		ug/L		81	60 - 122
Xylenes, total	150	146		ug/L		98	78 - 134

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-BS1
Matrix: Water
Analysis Batch: U014048

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H1881_P

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
1,2-Dichloroethane-d4	*104*		*63 - 140*
Dibromofluoromethane	*107*		*73 - 131*
Toluene-d8	*95*		*80 - 120*
4-Bromofluorobenzene	*90*		*79 - 125*

Lab Sample ID: 11H1881-BSD1
Matrix: Water
Analysis Batch: U014048

Client Sample ID: Lab Control Sample Dup
Prep Type: Total
Prep Batch: 11H1881_P

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	250	394	L1	ug/L		157	56 - 150	3	31
Tert-Amyl Methyl Ether	50.0	44.0		ug/L		88	70 - 133	2	16
Benzene	50.0	47.5		ug/L		95	80 - 121	2	12
Bromobenzene	50.0	43.2		ug/L		86	72 - 130	2	23
Bromochloromethane	50.0	51.8		ug/L		104	73 - 137	0.3	32
Ethyl tert-Butyl Ether	50.0	42.8		ug/L		86	68 - 138	2	16
Bromodichloromethane	50.0	51.6		ug/L		103	75 - 131	1	13
Diisopropyl Ether	50.0	42.7		ug/L		85	63 - 136	1	32
Methyl tert-Butyl Ether	50.0	47.3		ug/L		95	76 - 120	2	32
Bromoform	50.0	50.2		ug/L		100	65 - 140	0.6	18
Tertiary Butyl Alcohol	500	566		ug/L		113	60 - 140	0.2	30
Bromomethane	50.0	43.5		ug/L		87	50 - 150	3	50
2-Butanone	250	305		ug/L		122	70 - 144	0.1	37
sec-Butylbenzene	50.0	47.0		ug/L		94	72 - 140	2	21
n-Butylbenzene	50.0	49.9		ug/L		100	68 - 140	1	11
tert-Butylbenzene	50.0	46.9		ug/L		94	76 - 135	2	20
Carbon disulfide	50.0	44.3		ug/L		89	74 - 137	2	28
Carbon Tetrachloride	50.0	57.1		ug/L		114	71 - 137	3	26
Chlorobenzene	50.0	48.4		ug/L		97	80 - 121	2	11
Chlorodibromomethane	50.0	56.0		ug/L		112	68 - 137	2	16
Chloroethane	50.0	42.7		ug/L		85	50 - 146	4	35
Chloroform	50.0	50.7		ug/L		101	73 - 131	3	32
Chloromethane	50.0	32.6		ug/L		65	30 - 132	3	34
2-Chlorotoluene	50.0	43.4		ug/L		87	74 - 135	2	22
4-Chlorotoluene	50.0	45.8		ug/L		92	74 - 132	2	22
1,2-Dibromo-3-chloropropane	50.0	49.8		ug/L		100	56 - 145	4	21
1,2-Dibromoethane (EDB)	50.0	50.0		ug/L		100	80 - 135	2	10
Dibromomethane	50.0	51.1		ug/L		102	78 - 133	2	11
1,4-Dichlorobenzene	50.0	49.4		ug/L		99	80 - 120	1	10
1,3-Dichlorobenzene	50.0	48.7		ug/L		97	80 - 128	1	18
1,2-Dichlorobenzene	50.0	53.7		ug/L		107	80 - 125	0.7	11
Dichlorodifluoromethane	50.0	30.8		ug/L		62	30 - 132	1	32
1,1-Dichloroethane	50.0	48.6		ug/L		97	75 - 125	3	34
1,2-Dichloroethane	50.0	51.7		ug/L		103	70 - 134	2	25
cis-1,2-Dichloroethene	50.0	48.4		ug/L		97	71 - 132	1	32
1,1-Dichloroethene	50.0	50.1		ug/L		100	73 - 125	2	31
trans-1,2-Dichloroethene	50.0	49.6		ug/L		99	77 - 125	2	32
1,3-Dichloropropane	50.0	47.6		ug/L		95	76 - 125	0.9	20
1,2-Dichloropropane	50.0	44.0		ug/L		88	72 - 120	1	11

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-BSD1 | **Client Sample ID: Lab Control Sample Dup**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014048 | **Prep Batch: 11H1881_P**

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
2,2-Dichloropropane	50.0	51.0		ug/L		102	50 - 150	4	11
cis-1,3-Dichloropropene	50.0	50.8		ug/L		102	70 - 140	0.4	35
trans-1,3-Dichloropropene	50.0	50.2		ug/L		100	62 - 139	1	26
1,1-Dichloropropene	50.0	49.7		ug/L		99	78 - 126	2	18
Ethylbenzene	50.0	46.7		ug/L		93	78 - 133	2	12
Hexachlorobutadiene	50.0	49.8		ug/L		100	70 - 150	1	21
2-Hexanone	250	277		ug/L		111	60 - 150	1	20
Isopropylbenzene	50.0	53.2		ug/L		106	69 - 120	1	15
p-Isopropyltoluene	50.0	47.1		ug/L		94	72 - 134	2	18
Methylene Chloride	50.0	49.8		ug/L		100	80 - 133	2	36
4-Methyl-2-pentanone	250	237		ug/L		95	62 - 146	1	35
Naphthalene	50.0	49.8		ug/L		100	71 - 139	0.1	30
n-Propylbenzene	50.0	44.6		ug/L		89	70 - 143	2	23
Styrene	50.0	49.4		ug/L		99	80 - 136	0.8	29
1,1,1,2-Tetrachloroethane	50.0	53.4		ug/L		107	80 - 130	0.6	11
1,1,2,2-Tetrachloroethane	50.0	44.7		ug/L		89	73 - 131	4	28
Tetrachloroethene	50.0	51.8		ug/L		104	77 - 131	2	16
Toluene	50.0	46.6		ug/L		93	78 - 125	1	35
1,2,3-Trichlorobenzene	50.0	53.0		ug/L		106	71 - 138	0.6	28
1,2,4-Trichlorobenzene	50.0	52.5		ug/L		105	74 - 136	0.2	23
1,1,2-Trichloroethane	50.0	49.9		ug/L		100	80 - 123	0.8	21
1,1,1-Trichloroethane	50.0	54.0		ug/L		108	75 - 137	2	29
Trichloroethene	50.0	52.7		ug/L		105	74 - 139	0.9	11
Trichlorofluoromethane	50.0	46.5		ug/L		93	60 - 133	7	33
1,2,3-Trichloropropane	50.0	47.1		ug/L		94	64 - 127	2	25
1,3,5-Trimethylbenzene	50.0	46.0		ug/L		92	75 - 134	1	21
1,2,4-Trimethylbenzene	50.0	44.8		ug/L		90	77 - 134	2	20
Vinyl chloride	50.0	39.7		ug/L		79	60 - 122	2	32
Xylenes, total	150	144		ug/L		96	78 - 134	2	18

Surrogate	*LCS Dup % Recovery*	*LCS Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*103*		*63 - 140*
Dibromofluoromethane	*108*		*73 - 131*
Toluene-d8	*96*		*80 - 120*
4-Bromofluorobenzene	*90*		*79 - 125*

Lab Sample ID: 11H1881-MS1 | **Client Sample ID: MW-38 (08022011)**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014048 | **Prep Batch: 11H1881_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	ND	L	250	271		ug/L		109	56 - 150
Tert-Amyl Methyl Ether	ND		50.0	44.2		ug/L		88	69 - 139
Benzene	46.4		50.0	89.0		ug/L		85	65 - 151
Bromobenzene	ND		50.0	40.4		ug/L		81	69 - 142
Bromochloromethane	ND		50.0	55.1		ug/L		110	64 - 154
Ethyl tert-Butyl Ether	ND		50.0	42.0		ug/L		84	68 - 139
Bromodichloromethane	ND		50.0	50.4		ug/L		101	75 - 138
Diisopropyl Ether	2.31		50.0	41.3		ug/L		78	59 - 145

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-MS1 | **Client Sample ID: MW-38 (08022011)**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014048 | **Prep Batch: 11H1881_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Methyl tert-Butyl Ether	7.70		50.0	53.0		ug/L		91	56 - 152
Bromoform	ND		50.0	48.1		ug/L		96	55 - 153
Tertiary Butyl Alcohol	ND		500	535		ug/L		107	14 - 200
Bromomethane	ND		50.0	45.7		ug/L		91	13 - 176
2-Butanone	ND		250	246		ug/L		98	45 - 164
sec-Butylbenzene	4.29		50.0	47.2		ug/L		86	68 - 159
n-Butylbenzene	3.36		50.0	44.8		ug/L		83	67 - 151
tert-Butylbenzene	0.380		50.0	44.8		ug/L		89	73 - 153
Carbon disulfide	ND		50.0	42.4		ug/L		85	33 - 187
Carbon Tetrachloride	ND		50.0	56.7		ug/L		113	64 - 157
Chlorobenzene	ND		50.0	47.9		ug/L		96	78 - 136
Chlorodibromomethane	ND		50.0	55.7		ug/L		111	64 - 145
Chloroethane	ND		50.0	39.5		ug/L		79	48 - 159
Chloroform	ND		50.0	53.7		ug/L		107	72 - 145
Chloromethane	ND		50.0	28.0		ug/L		56	10 - 194
2-Chlorotoluene	ND		50.0	40.7		ug/L		81	66 - 155
4-Chlorotoluene	ND		50.0	42.7		ug/L		85	69 - 149
1,2-Dibromo-3-chloropropane	ND		50.0	45.9		ug/L		92	49 - 162
1,2-Dibromoethane (EDB)	ND		50.0	49.5		ug/L		99	70 - 152
Dibromomethane	ND		50.0	50.6		ug/L		101	75 - 141
1,4-Dichlorobenzene	ND		50.0	47.6		ug/L		95	75 - 135
1,3-Dichlorobenzene	ND		50.0	47.3		ug/L		95	72 - 146
1,2-Dichlorobenzene	ND		50.0	48.0		ug/L		96	80 - 136
Dichlorodifluoromethane	ND		50.0	33.2		ug/L		66	23 - 159
1,1-Dichloroethane	ND		50.0	47.7		ug/L		95	64 - 154
1,2-Dichloroethane	ND		50.0	51.3		ug/L		103	72 - 137
cis-1,2-Dichloroethene	ND		50.0	44.8		ug/L		90	57 - 154
1,1-Dichloroethene	ND		50.0	49.3		ug/L		99	34 - 151
trans-1,2-Dichloroethene	ND		50.0	45.4		ug/L		91	57 - 157
1,3-Dichloropropane	ND		50.0	44.9		ug/L		90	71 - 137
1,2-Dichloropropane	ND		50.0	42.9		ug/L		86	71 - 139
2,2-Dichloropropane	ND		50.0	44.8		ug/L		90	10 - 198
cis-1,3-Dichloropropene	ND		50.0	46.7		ug/L		93	56 - 156
trans-1,3-Dichloropropene	ND		50.0	45.4		ug/L		91	47 - 157
1,1-Dichloropropene	ND		50.0	47.5		ug/L		95	70 - 155
Ethylbenzene	43.2		50.0	80.7		ug/L		75	68 - 157
Hexachlorobutadiene	ND		50.0	50.5		ug/L		101	47 - 173
2-Hexanone	ND		250	212		ug/L		85	57 - 154
Isopropylbenzene	22.7		50.0	69.9		ug/L		94	69 - 139
p-Isopropyltoluene	ND		50.0	44.0		ug/L		88	69 - 151
Methylene Chloride	ND		50.0	46.0		ug/L		92	71 - 136
4-Methyl-2-pentanone	ND		250	208		ug/L		83	62 - 159
Naphthalene	1.48		50.0	49.6		ug/L		96	56 - 161
n-Propylbenzene	16.6		50.0	54.6		ug/L		76	61 - 167
Styrene	ND		50.0	47.4		ug/L		95	69 - 150
1,1,1,2-Tetrachloroethane	ND		50.0	52.9		ug/L		106	80 - 140
1,1,2,2-Tetrachloroethane	ND		50.0	42.1		ug/L		84	76 - 141
Tetrachloroethene	ND		50.0	51.2		ug/L		102	63 - 155
Toluene	0.580		50.0	46.6		ug/L		92	61 - 153
1,2,3-Trichlorobenzene	ND		50.0	53.9		ug/L		108	57 - 155

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-MS1 — **Client Sample ID: MW-38 (08022011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014048 — **Prep Batch: 11H1881_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
1,2,4-Trichlorobenzene	ND		50.0	53.7		ug/L		107	64 - 147
1,1,2-Trichloroethane	ND		50.0	54.1		ug/L		108	74 - 138
1,1,1-Trichloroethane	ND		50.0	54.3		ug/L		109	78 - 153
Trichloroethene	ND		50.0	54.6		ug/L		109	74 - 139
Trichlorofluoromethane	ND		50.0	46.0		ug/L		92	53 - 149
1,2,3-Trichloropropane	ND		50.0	42.5		ug/L		85	49 - 148
1,3,5-Trimethylbenzene	ND		50.0	43.4		ug/L		87	67 - 151
1,2,4-Trimethylbenzene	ND		50.0	42.4		ug/L		85	69 - 150
Vinyl chloride	ND		50.0	32.3		ug/L		65	53 - 137
Xylenes, total	3.00		150	138		ug/L		90	68 - 158

Surrogate	*Matrix Spike % Recovery*	*Matrix Spike Qualifier*	*Limits*
1,2-Dichloroethane-d4	*94*		*63 - 140*
Dibromofluoromethane	*109*		*73 - 131*
Toluene-d8	*95*		*80 - 120*
4-Bromofluorobenzene	*89*		*79 - 125*

Lab Sample ID: 11H1881-MSD1 — **Client Sample ID: MW-38 (08022011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014048 — **Prep Batch: 11H1881_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	ND	L	250	257		ug/L		103	56 - 150	5	31
Tert-Amyl Methyl Ether	ND		50.0	39.8		ug/L		80	69 - 139	10	16
Benzene	46.4		50.0	86.8		ug/L		81	65 - 151	2	12
Bromobenzene	ND		50.0	38.7		ug/L		77	69 - 142	4	23
Bromochloromethane	ND		50.0	51.1		ug/L		102	64 - 154	8	32
Ethyl tert-Butyl Ether	ND		50.0	38.2		ug/L		76	68 - 139	10	16
Bromodichloromethane	ND		50.0	46.5		ug/L		93	75 - 138	8	13
Diisopropyl Ether	2.31		50.0	38.2		ug/L		72	59 - 145	8	32
Methyl tert-Butyl Ether	7.70		50.0	48.4		ug/L		81	56 - 152	9	32
Bromoform	ND		50.0	44.7		ug/L		89	55 - 153	7	18
Tertiary Butyl Alcohol	ND		500	495		ug/L		99	14 - 200	8	30
Bromomethane	ND		50.0	44.1		ug/L		88	13 - 176	4	50
2-Butanone	ND		250	220		ug/L		88	45 - 164	11	37
sec-Butylbenzene	4.29		50.0	47.6		ug/L		87	68 - 159	1	21
n-Butylbenzene	3.36		50.0	45.3		ug/L		84	67 - 151	1	11
tert-Butylbenzene	0.380		50.0	44.9		ug/L		89	73 - 153	0.2	20
Carbon disulfide	ND		50.0	41.1		ug/L		82	33 - 187	3	28
Carbon Tetrachloride	ND		50.0	55.1		ug/L		110	64 - 157	3	26
Chlorobenzene	ND		50.0	46.1		ug/L		92	78 - 136	4	11
Chlorodibromomethane	ND		50.0	50.4		ug/L		101	64 - 145	10	16
Chloroethane	ND		50.0	38.2		ug/L		76	48 - 159	4	35
Chloroform	ND		50.0	50.1		ug/L		100	72 - 145	7	32
Chloromethane	ND		50.0	26.7		ug/L		53	10 - 194	5	34
2-Chlorotoluene	ND		50.0	40.2		ug/L		80	66 - 155	1	22
4-Chlorotoluene	ND		50.0	42.0		ug/L		84	69 - 149	2	22
1,2-Dibromo-3-chloropropane	ND		50.0	43.4		ug/L		87	49 - 162	6	21
1,2-Dibromoethane (EDB)	ND		50.0	45.6		ug/L		91	70 - 152	8	10

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H1881-MSD1 | **Client Sample ID: MW-38 (08022011)**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014048 | **Prep Batch: 11H1881_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Dibromomethane	ND		50.0	46.7		ug/L		93	75 - 141	8	11
1,4-Dichlorobenzene	ND		50.0	46.5		ug/L		93	75 - 135	2	10
1,3-Dichlorobenzene	ND		50.0	46.0		ug/L		92	72 - 146	3	18
1,2-Dichlorobenzene	ND		50.0	46.2		ug/L		92	80 - 136	4	11
Dichlorodifluoromethane	ND		50.0	33.2		ug/L		66	23 - 159	0.1	32
1,1-Dichloroethane	ND		50.0	44.0		ug/L		88	64 - 154	8	34
1,2-Dichloroethane	ND		50.0	48.2		ug/L		96	72 - 137	6	25
cis-1,2-Dichloroethene	ND		50.0	42.1		ug/L		84	57 - 154	6	32
1,1-Dichloroethene	ND		50.0	44.9		ug/L		90	34 - 151	9	31
trans-1,2-Dichloroethene	ND		50.0	42.7		ug/L		85	57 - 157	6	32
1,3-Dichloropropane	ND		50.0	41.0		ug/L		82	71 - 137	9	20
1,2-Dichloropropane	ND		50.0	40.1		ug/L		80	71 - 139	7	11
2,2-Dichloropropane	ND		50.0	42.1		ug/L		84	10 - 198	6	11
cis-1,3-Dichloropropene	ND		50.0	43.6		ug/L		87	56 - 156	7	35
trans-1,3-Dichloropropene	ND		50.0	41.5		ug/L		83	47 - 157	9	26
1,1-Dichloropropene	ND		50.0	46.6		ug/L		93	70 - 155	2	18
Ethylbenzene	43.2		50.0	82.5		ug/L		79	68 - 157	2	12
Hexachlorobutadiene	ND		50.0	49.7		ug/L		99	47 - 173	2	21
2-Hexanone	ND		250	196		ug/L		78	57 - 154	8	20
Isopropylbenzene	22.7		50.0	70.5		ug/L		96	69 - 139	0.8	15
p-Isopropyltoluene	ND		50.0	44.1		ug/L		88	69 - 151	0.1	18
Methylene Chloride	ND		50.0	43.6		ug/L		87	71 - 136	5	36
4-Methyl-2-pentanone	ND		250	191		ug/L		76	62 - 159	9	35
Naphthalene	1.48		50.0	47.6		ug/L		92	56 - 161	4	30
n-Propylbenzene	16.6		50.0	56.1		ug/L		79	61 - 167	3	23
Styrene	ND		50.0	45.4		ug/L		91	69 - 150	4	29
1,1,1,2-Tetrachloroethane	ND		50.0	49.5		ug/L		99	80 - 140	7	11
1,1,2,2-Tetrachloroethane	ND		50.0	39.3		ug/L		79	76 - 141	7	28
Tetrachloroethene	ND		50.0	51.0		ug/L		102	63 - 155	0.5	16
Toluene	0.580		50.0	44.4		ug/L		88	61 - 153	5	35
1,2,3-Trichlorobenzene	ND		50.0	51.5		ug/L		103	57 - 155	4	28
1,2,4-Trichlorobenzene	ND		50.0	51.8		ug/L		104	64 - 147	4	23
1,1,2-Trichloroethane	ND		50.0	50.8		ug/L		102	74 - 138	6	21
1,1,1-Trichloroethane	ND		50.0	51.6		ug/L		103	78 - 153	5	29
Trichloroethene	ND		50.0	53.1		ug/L		106	74 - 139	3	11
Trichlorofluoromethane	ND		50.0	45.3		ug/L		91	53 - 149	2	33
1,2,3-Trichloropropane	ND		50.0	39.4		ug/L		79	49 - 148	8	25
1,3,5-Trimethylbenzene	ND		50.0	43.3		ug/L		87	67 - 151	0.07	21
1,2,4-Trimethylbenzene	ND		50.0	42.3		ug/L		85	69 - 150	0.3	20
Vinyl chloride	ND		50.0	30.6		ug/L		61	53 - 137	5	32
Xylenes, total	3.00		150	134		ug/L		87	68 - 158	3	18

Surrogate	*Matrix Spike Dup % Recovery*	*Matrix Spike Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*97*		*63 - 140*
Dibromofluoromethane	*108*		*73 - 131*
Toluene-d8	*94*		*80 - 120*
4-Bromofluorobenzene	*89*		*79 - 125*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-BLK1
Matrix: Water
Analysis Batch: U014133

Client Sample ID: Method Blank
Prep Type: Total
Prep Batch: 11H2288_P

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Benzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Bromobenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Bromochloromethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Bromodichloromethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Bromoform	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Bromomethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
2-Butanone	ND		50.0		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
n-Butylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Carbon disulfide	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Chlorobenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Chloroethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Chloroform	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Chloromethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Dibromomethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Ethylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
2-Hexanone	ND		50.0		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Isopropylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Methylene Chloride	ND		5.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-BLK1 | **Client Sample ID: Method Blank**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014133 | **Prep Batch: 11H2288_P**

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
4-Methyl-2-pentanone	ND		10.0		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Naphthalene	ND		5.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
n-Propylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Styrene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Tetrachloroethene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Toluene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Trichloroethene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Vinyl chloride	ND		1.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00
Xylenes, total	ND		3.00		ug/L		08/09/11 18:04	08/09/11 20:52	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*99*		*63 - 140*	*08/09/11 18:04*	*08/09/11 20:52*	*1.00*
Dibromofluoromethane	*104*		*73 - 131*	*08/09/11 18:04*	*08/09/11 20:52*	*1.00*
Toluene-d8	*94*		*80 - 120*	*08/09/11 18:04*	*08/09/11 20:52*	*1.00*
4-Bromofluorobenzene	*91*		*79 - 125*	*08/09/11 18:04*	*08/09/11 20:52*	*1.00*

Lab Sample ID: 11H2288-BS1 | **Client Sample ID: Lab Control Sample**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014133 | **Prep Batch: 11H2288_P**

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	250	235		ug/L		94	56 - 150
Tert-Amyl Methyl Ether	50.0	46.0		ug/L		92	70 - 133
Benzene	50.0	48.1		ug/L		96	80 - 121
Bromobenzene	50.0	43.4		ug/L		87	72 - 130
Bromochloromethane	50.0	53.0		ug/L		106	73 - 137
Ethyl tert-Butyl Ether	50.0	43.0		ug/L		86	68 - 138
Bromodichloromethane	50.0	49.4		ug/L		99	75 - 131
Diisopropyl Ether	50.0	39.0		ug/L		78	63 - 136
Methyl tert-Butyl Ether	50.0	48.5		ug/L		97	76 - 120
Bromoform	50.0	47.7		ug/L		95	65 - 140
Tertiary Butyl Alcohol	500	534		ug/L		107	60 - 140
Bromomethane	50.0	42.1		ug/L		84	50 - 150
2-Butanone	250	247		ug/L		99	70 - 144
sec-Butylbenzene	50.0	49.2		ug/L		98	72 - 140
n-Butylbenzene	50.0	48.4		ug/L		97	68 - 140
tert-Butylbenzene	50.0	49.7		ug/L		99	76 - 135
Carbon disulfide	50.0	43.6		ug/L		87	74 - 137
Carbon Tetrachloride	50.0	56.8		ug/L		114	71 - 137

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-BS1
Matrix: Water
Analysis Batch: U014133

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H2288_P

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Chlorobenzene	50.0	49.8		ug/L		100	80 - 121
Chlorodibromomethane	50.0	54.6		ug/L		109	68 - 137
Chloroethane	50.0	39.0		ug/L		78	50 - 146
Chloroform	50.0	49.9		ug/L		100	73 - 131
Chloromethane	50.0	31.5		ug/L		63	30 - 132
2-Chlorotoluene	50.0	44.8		ug/L		90	74 - 135
4-Chlorotoluene	50.0	46.7		ug/L		93	74 - 132
1,2-Dibromo-3-chloropropane	50.0	49.2		ug/L		98	56 - 145
1,2-Dibromoethane (EDB)	50.0	51.1		ug/L		102	80 - 135
Dibromomethane	50.0	50.9		ug/L		102	78 - 133
1,4-Dichlorobenzene	50.0	51.7		ug/L		103	80 - 120
1,3-Dichlorobenzene	50.0	50.6		ug/L		101	80 - 128
1,2-Dichlorobenzene	50.0	53.6		ug/L		107	80 - 125
Dichlorodifluoromethane	50.0	38.1		ug/L		76	30 - 132
1,1-Dichloroethane	50.0	46.3		ug/L		93	75 - 125
1,2-Dichloroethane	50.0	48.7		ug/L		97	70 - 134
cis-1,2-Dichloroethene	50.0	46.3		ug/L		93	71 - 132
1,1-Dichloroethene	50.0	49.8		ug/L		100	73 - 125
trans-1,2-Dichloroethene	50.0	46.9		ug/L		94	77 - 125
1,3-Dichloropropane	50.0	45.9		ug/L		92	76 - 125
1,2-Dichloropropane	50.0	42.7		ug/L		85	72 - 120
2,2-Dichloropropane	50.0	50.8		ug/L		102	50 - 150
cis-1,3-Dichloropropene	50.0	48.8		ug/L		98	70 - 140
trans-1,3-Dichloropropene	50.0	47.1		ug/L		94	62 - 139
1,1-Dichloropropene	50.0	49.7		ug/L		99	78 - 126
Ethylbenzene	50.0	48.0		ug/L		96	78 - 133
Hexachlorobutadiene	50.0	54.8		ug/L		110	70 - 150
2-Hexanone	250	231		ug/L		93	60 - 150
Isopropylbenzene	50.0	54.7		ug/L		109	69 - 120
p-Isopropyltoluene	50.0	49.3		ug/L		99	72 - 134
Methylene Chloride	50.0	48.2		ug/L		96	80 - 133
4-Methyl-2-pentanone	250	225		ug/L		90	62 - 146
Naphthalene	50.0	53.6		ug/L		107	71 - 139
n-Propylbenzene	50.0	46.3		ug/L		93	70 - 143
Styrene	50.0	49.6		ug/L		99	80 - 136
1,1,1,2-Tetrachloroethane	50.0	52.6		ug/L		105	80 - 130
1,1,2,2-Tetrachloroethane	50.0	45.1		ug/L		90	73 - 131
Tetrachloroethene	50.0	54.4		ug/L		109	77 - 131
Toluene	50.0	48.4		ug/L		97	78 - 125
1,2,3-Trichlorobenzene	50.0	56.0		ug/L		112	71 - 138
1,2,4-Trichlorobenzene	50.0	56.3		ug/L		113	74 - 136
1,1,2-Trichloroethane	50.0	49.7		ug/L		99	80 - 123
1,1,1-Trichloroethane	50.0	53.3		ug/L		107	75 - 137
Trichloroethene	50.0	55.4		ug/L		111	74 - 139
Trichlorofluoromethane	50.0	46.5		ug/L		93	60 - 133
1,2,3-Trichloropropane	50.0	44.9		ug/L		90	64 - 127
1,3,5-Trimethylbenzene	50.0	48.2		ug/L		96	75 - 134
1,2,4-Trimethylbenzene	50.0	47.1		ug/L		94	77 - 134
Vinyl chloride	50.0	38.5		ug/L		77	60 - 122
Xylenes, total	150	147		ug/L		98	78 - 134

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-BS1
Matrix: Water
Analysis Batch: U014133

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H2288_P

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
1,2-Dichloroethane-d4	*96*		*63 - 140*
Dibromofluoromethane	*105*		*73 - 131*
Toluene-d8	*95*		*80 - 120*
4-Bromofluorobenzene	*92*		*79 - 125*

Lab Sample ID: 11H2288-BSD1
Matrix: Water
Analysis Batch: U014133

Client Sample ID: Lab Control Sample Dup
Prep Type: Total
Prep Batch: 11H2288_P

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	250	250		ug/L		100	56 - 150	7	31
Tert-Amyl Methyl Ether	50.0	45.2		ug/L		90	70 - 133	2	16
Benzene	50.0	47.2		ug/L		94	80 - 121	2	12
Bromobenzene	50.0	42.5		ug/L		85	72 - 130	2	23
Bromochloromethane	50.0	52.8		ug/L		106	73 - 137	0.5	32
Ethyl tert-Butyl Ether	50.0	42.5		ug/L		85	68 - 138	1	16
Bromodichloromethane	50.0	49.6		ug/L		99	75 - 131	0.4	13
Diisopropyl Ether	50.0	38.7		ug/L		77	63 - 136	0.9	32
Methyl tert-Butyl Ether	50.0	47.6		ug/L		95	76 - 120	2	32
Bromoform	50.0	47.8		ug/L		96	65 - 140	0.2	18
Tertiary Butyl Alcohol	500	553		ug/L		111	60 - 140	4	30
Bromomethane	50.0	43.0		ug/L		86	50 - 150	2	50
2-Butanone	250	254		ug/L		101	70 - 144	3	37
sec-Butylbenzene	50.0	47.9		ug/L		96	72 - 140	3	21
n-Butylbenzene	50.0	48.4		ug/L		97	68 - 140	0.08	11
tert-Butylbenzene	50.0	47.8		ug/L		96	76 - 135	4	20
Carbon disulfide	50.0	42.6		ug/L		85	74 - 137	2	28
Carbon Tetrachloride	50.0	55.8		ug/L		112	71 - 137	2	26
Chlorobenzene	50.0	49.2		ug/L		98	80 - 121	1	11
Chlorodibromomethane	50.0	54.4		ug/L		109	68 - 137	0.3	16
Chloroethane	50.0	38.4		ug/L		77	50 - 146	2	35
Chloroform	50.0	49.2		ug/L		98	73 - 131	1	32
Chloromethane	50.0	30.9		ug/L		62	30 - 132	2	34
2-Chlorotoluene	50.0	43.8		ug/L		88	74 - 135	2	22
4-Chlorotoluene	50.0	45.6		ug/L		91	74 - 132	2	22
1,2-Dibromo-3-chloropropane	50.0	49.4		ug/L		99	56 - 145	0.4	21
1,2-Dibromoethane (EDB)	50.0	49.9		ug/L		100	80 - 135	2	10
Dibromomethane	50.0	50.6		ug/L		101	78 - 133	0.5	11
1,4-Dichlorobenzene	50.0	50.0		ug/L		100	80 - 120	3	10
1,3-Dichlorobenzene	50.0	49.4		ug/L		99	80 - 128	2	18
1,2-Dichlorobenzene	50.0	52.4		ug/L		105	80 - 125	2	11
Dichlorodifluoromethane	50.0	37.6		ug/L		75	30 - 132	1	32
1,1-Dichloroethane	50.0	45.7		ug/L		91	75 - 125	1	34
1,2-Dichloroethane	50.0	48.3		ug/L		97	70 - 134	0.8	25
cis-1,2-Dichloroethene	50.0	45.6		ug/L		91	71 - 132	2	32
1,1-Dichloroethene	50.0	49.2		ug/L		98	73 - 125	1	31
trans-1,2-Dichloroethene	50.0	46.8		ug/L		94	77 - 125	0.4	32
1,3-Dichloropropane	50.0	45.0		ug/L		90	76 - 125	2	20
1,2-Dichloropropane	50.0	41.9		ug/L		84	72 - 120	2	11

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-BSD1 | **Client Sample ID: Lab Control Sample Dup**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014133 | **Prep Batch: 11H2288_P**

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
2,2-Dichloropropane	50.0	49.2		ug/L		98	50 - 150	3	11
cis-1,3-Dichloropropene	50.0	48.1		ug/L		96	70 - 140	1	35
trans-1,3-Dichloropropene	50.0	46.7		ug/L		93	62 - 139	1	26
1,1-Dichloropropene	50.0	48.6		ug/L		97	78 - 126	2	18
Ethylbenzene	50.0	47.5		ug/L		95	78 - 133	1	12
Hexachlorobutadiene	50.0	52.9		ug/L		106	70 - 150	3	21
2-Hexanone	250	231		ug/L		93	60 - 150	0.07	20
Isopropylbenzene	50.0	53.9		ug/L		108	69 - 120	1	15
p-Isopropyltoluene	50.0	47.6		ug/L		95	72 - 134	3	18
Methylene Chloride	50.0	48.8		ug/L		98	80 - 133	1	36
4-Methyl-2-pentanone	250	223		ug/L		89	62 - 146	0.6	35
Naphthalene	50.0	51.6		ug/L		103	71 - 139	4	30
n-Propylbenzene	50.0	45.3		ug/L		91	70 - 143	2	23
Styrene	50.0	49.4		ug/L		99	80 - 136	0.4	29
1,1,1,2-Tetrachloroethane	50.0	52.1		ug/L		104	80 - 130	0.9	11
1,1,2,2-Tetrachloroethane	50.0	44.6		ug/L		89	73 - 131	1	28
Tetrachloroethene	50.0	53.8		ug/L		108	77 - 131	1	16
Toluene	50.0	47.6		ug/L		95	78 - 125	2	35
1,2,3-Trichlorobenzene	50.0	55.0		ug/L		110	71 - 138	2	28
1,2,4-Trichlorobenzene	50.0	55.4		ug/L		111	74 - 136	2	23
1,1,2-Trichloroethane	50.0	48.8		ug/L		98	80 - 123	2	21
1,1,1-Trichloroethane	50.0	52.5		ug/L		105	75 - 137	2	29
Trichloroethene	50.0	54.4		ug/L		109	74 - 139	2	11
Trichlorofluoromethane	50.0	45.7		ug/L		91	60 - 133	2	33
1,2,3-Trichloropropane	50.0	45.0		ug/L		90	64 - 127	0.3	25
1,3,5-Trimethylbenzene	50.0	46.9		ug/L		94	75 - 134	3	21
1,2,4-Trimethylbenzene	50.0	45.6		ug/L		91	77 - 134	3	20
Vinyl chloride	50.0	37.2		ug/L		74	60 - 122	3	32
Xylenes, total	150	145		ug/L		96	78 - 134	1	18

Surrogate	*LCS Dup % Recovery*	*LCS Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*96*		*63 - 140*
Dibromofluoromethane	*105*		*73 - 131*
Toluene-d8	*94*		*80 - 120*
4-Bromofluorobenzene	*91*		*79 - 125*

Lab Sample ID: 11H2288-MS1 | **Client Sample ID: Matrix Spike**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014133 | **Prep Batch: 11H2288_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	ND		25000	21300		ug/L		85	56 - 150
Tert-Amyl Methyl Ether	ND		5000	4660		ug/L		93	69 - 139
Benzene	18500		5000	21800		ug/L		65	65 - 151
Bromobenzene	ND		5000	4500		ug/L		90	69 - 142
Bromochloromethane	ND		5000	5860		ug/L		117	64 - 154
Ethyl tert-Butyl Ether	ND		5000	4330		ug/L		87	68 - 139
Bromodichloromethane	ND		5000	5420		ug/L		108	75 - 138
Diisopropyl Ether	ND		5000	4130		ug/L		83	59 - 145

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-MS1
Matrix: Water
Analysis Batch: U014133

Client Sample ID: Matrix Spike
Prep Type: Total
Prep Batch: 11H2288_P

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Methyl tert-Butyl Ether	ND		5000	4870		ug/L		97	56 - 152
Bromoform	ND		5000	5100		ug/L		102	55 - 153
Tertiary Butyl Alcohol	ND		50000	47100		ug/L		94	14 - 200
Bromomethane	ND		5000	4160		ug/L		83	13 - 176
2-Butanone	ND		25000	25200		ug/L		101	45 - 164
sec-Butylbenzene	ND		5000	5120		ug/L		102	68 - 159
n-Butylbenzene	ND		5000	5130		ug/L		103	67 - 151
tert-Butylbenzene	ND		5000	5240		ug/L		105	73 - 153
Carbon disulfide	ND		5000	4580		ug/L		92	33 - 187
Carbon Tetrachloride	ND		5000	6320		ug/L		126	64 - 157
Chlorobenzene	ND		5000	5350		ug/L		107	78 - 136
Chlorodibromomethane	ND		5000	5950		ug/L		119	64 - 145
Chloroethane	ND		5000	4290		ug/L		86	48 - 159
Chloroform	ND		5000	5410		ug/L		108	72 - 145
Chloromethane	ND		5000	3490		ug/L		70	10 - 194
2-Chlorotoluene	ND		5000	4740		ug/L		95	66 - 155
4-Chlorotoluene	ND		5000	4910		ug/L		98	69 - 149
1,2-Dibromo-3-chloropropane	ND		5000	5050		ug/L		101	49 - 162
1,2-Dibromoethane (EDB)	ND		5000	5250		ug/L		105	70 - 152
Dibromomethane	ND		5000	5470		ug/L		109	75 - 141
1,4-Dichlorobenzene	ND		5000	5440		ug/L		109	75 - 135
1,3-Dichlorobenzene	ND		5000	5390		ug/L		108	72 - 146
1,2-Dichlorobenzene	ND		5000	5730		ug/L		115	80 - 136
Dichlorodifluoromethane	ND		5000	3670		ug/L		73	23 - 159
1,1-Dichloroethane	ND		5000	5030		ug/L		101	64 - 154
1,2-Dichloroethane	ND		5000	5900		ug/L		118	72 - 137
cis-1,2-Dichloroethene	ND		5000	4920		ug/L		98	57 - 154
1,1-Dichloroethene	ND		5000	5360		ug/L		107	34 - 151
trans-1,2-Dichloroethene	ND		5000	5060		ug/L		101	57 - 157
1,3-Dichloropropane	ND		5000	4730		ug/L		95	71 - 137
1,2-Dichloropropane	ND		5000	4550		ug/L		91	71 - 139
2,2-Dichloropropane	ND		5000	4980		ug/L		100	10 - 198
cis-1,3-Dichloropropene	ND		5000	5050		ug/L		101	56 - 156
trans-1,3-Dichloropropene	ND		5000	4870		ug/L		97	47 - 157
1,1-Dichloropropene	ND		5000	5310		ug/L		106	70 - 155
Ethylbenzene	7770		5000	12100		ug/L		87	68 - 157
Hexachlorobutadiene	ND		5000	5340		ug/L		107	47 - 173
2-Hexanone	ND		25000	22300		ug/L		89	57 - 154
Isopropylbenzene	ND		5000	5820		ug/L		116	69 - 139
p-Isopropyltoluene	ND		5000	5160		ug/L		103	69 - 151
Methylene Chloride	ND		5000	5370		ug/L		107	71 - 136
4-Methyl-2-pentanone	ND		25000	22200		ug/L		89	62 - 159
Naphthalene	ND		5000	5270		ug/L		105	56 - 161
n-Propylbenzene	ND		5000	4860		ug/L		97	61 - 167
Styrene	ND		5000	5420		ug/L		108	69 - 150
1,1,1,2-Tetrachloroethane	ND		5000	5740		ug/L		115	80 - 140
1,1,2,2-Tetrachloroethane	ND		5000	4570		ug/L		91	76 - 141
Tetrachloroethene	ND		5000	5880		ug/L		118	63 - 155
Toluene	73.0		5000	5130		ug/L		101	61 - 153
1,2,3-Trichlorobenzene	ND		5000	5580		ug/L		112	57 - 155

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-MS1 — **Client Sample ID: Matrix Spike**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014133 — **Prep Batch: 11H2288_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
1,2,4-Trichlorobenzene	ND		5000	5750		ug/L		115	64 - 147
1,1,2-Trichloroethane	ND		5000	5150		ug/L		103	74 - 138
1,1,1-Trichloroethane	ND		5000	5840		ug/L		117	78 - 153
Trichloroethene	ND		5000	6030		ug/L		121	74 - 139
Trichlorofluoromethane	ND		5000	5170		ug/L		103	53 - 149
1,2,3-Trichloropropane	ND		5000	4570		ug/L		91	49 - 148
1,3,5-Trimethylbenzene	ND		5000	5050		ug/L		101	67 - 151
1,2,4-Trimethylbenzene	47.0		5000	4950		ug/L		98	69 - 150
Vinyl chloride	ND		5000	4080		ug/L		82	53 - 137
Xylenes, total	41400		15000	49600	M8	ug/L		54	68 - 158

Surrogate	*Matrix Spike % Recovery*	*Matrix Spike Qualifier*	*Limits*
1,2-Dichloroethane-d4	*95*		*63 - 140*
Dibromofluoromethane	*109*		*73 - 131*
Toluene-d8	*94*		*80 - 120*
4-Bromofluorobenzene	*88*		*79 - 125*

Lab Sample ID: 11H2288-MSD1 — **Client Sample ID: Matrix Spike Duplicate**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014133 — **Prep Batch: 11H2288_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	ND		25000	22100		ug/L		89	56 - 150	4	31
Tert-Amyl Methyl Ether	ND		5000	4600		ug/L		92	69 - 139	1	16
Benzene	18500		5000	21300	M8	ug/L		55	65 - 151	2	12
Bromobenzene	ND		5000	4500		ug/L		90	69 - 142	0	23
Bromochloromethane	ND		5000	5660		ug/L		113	64 - 154	3	32
Ethyl tert-Butyl Ether	ND		5000	4340		ug/L		87	68 - 139	0.3	16
Bromodichloromethane	ND		5000	5310		ug/L		106	75 - 138	2	13
Diisopropyl Ether	ND		5000	4010		ug/L		80	59 - 145	3	32
Methyl tert-Butyl Ether	ND		5000	4840		ug/L		97	56 - 152	0.7	32
Bromoform	ND		5000	5070		ug/L		101	55 - 153	0.6	18
Tertiary Butyl Alcohol	ND		50000	54400		ug/L		109	14 - 200	14	30
Bromomethane	ND		5000	4480		ug/L		90	13 - 176	7	50
2-Butanone	ND		25000	25000		ug/L		100	45 - 164	0.5	37
sec-Butylbenzene	ND		5000	5090		ug/L		102	68 - 159	0.6	21
n-Butylbenzene	ND		5000	4870		ug/L		97	67 - 151	5	11
tert-Butylbenzene	ND		5000	5110		ug/L		102	73 - 153	2	20
Carbon disulfide	ND		5000	4510		ug/L		90	33 - 187	2	28
Carbon Tetrachloride	ND		5000	6230		ug/L		125	64 - 157	1	26
Chlorobenzene	ND		5000	5220		ug/L		104	78 - 136	2	11
Chlorodibromomethane	ND		5000	5800		ug/L		116	64 - 145	3	16
Chloroethane	ND		5000	4150		ug/L		83	48 - 159	3	35
Chloroform	ND		5000	5370		ug/L		107	72 - 145	0.8	32
Chloromethane	ND		5000	3320		ug/L		66	10 - 194	5	34
2-Chlorotoluene	ND		5000	4690		ug/L		94	66 - 155	0.9	22
4-Chlorotoluene	ND		5000	4930		ug/L		99	69 - 149	0.6	22
1,2-Dibromo-3-chloropropane	ND		5000	5080		ug/L		102	49 - 162	0.7	21
1,2-Dibromoethane (EDB)	ND		5000	5180		ug/L		104	70 - 152	1	10

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2288-MSD1
Matrix: Water
Analysis Batch: U014133

Client Sample ID: Matrix Spike Duplicate
Prep Type: Total
Prep Batch: 11H2288_P

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Dibromomethane	ND		5000	5320		ug/L		106	75 - 141	3	11
1,4-Dichlorobenzene	ND		5000	5380		ug/L		108	75 - 135	1	10
1,3-Dichlorobenzene	ND		5000	5340		ug/L		107	72 - 146	0.8	18
1,2-Dichlorobenzene	ND		5000	5420		ug/L		108	80 - 136	6	11
Dichlorodifluoromethane	ND		5000	3560		ug/L		71	23 - 159	3	32
1,1-Dichloroethane	ND		5000	4870		ug/L		97	64 - 154	3	34
1,2-Dichloroethane	ND		5000	5790		ug/L		116	72 - 137	2	25
cis-1,2-Dichloroethene	ND		5000	4770		ug/L		95	57 - 154	3	32
1,1-Dichloroethene	ND		5000	5260		ug/L		105	34 - 151	2	31
trans-1,2-Dichloroethene	ND		5000	5010		ug/L		100	57 - 157	1	32
1,3-Dichloropropane	ND		5000	4690		ug/L		94	71 - 137	0.8	20
1,2-Dichloropropane	ND		5000	4450		ug/L		89	71 - 139	2	11
2,2-Dichloropropane	ND		5000	4880		ug/L		98	10 - 198	2	11
cis-1,3-Dichloropropene	ND		5000	4970		ug/L		99	56 - 156	2	35
trans-1,3-Dichloropropene	ND		5000	4780		ug/L		96	47 - 157	2	26
1,1-Dichloropropene	ND		5000	5080		ug/L		102	70 - 155	4	18
Ethylbenzene	7770		5000	11900		ug/L		82	68 - 157	2	12
Hexachlorobutadiene	ND		5000	5580		ug/L		112	47 - 173	4	21
2-Hexanone	ND		25000	22400		ug/L		90	57 - 154	0.6	20
Isopropylbenzene	ND		5000	5680		ug/L		114	69 - 139	2	15
p-Isopropyltoluene	ND		5000	5080		ug/L		102	69 - 151	2	18
Methylene Chloride	ND		5000	5160		ug/L		103	71 - 136	4	36
4-Methyl-2-pentanone	ND		25000	22000		ug/L		88	62 - 159	0.8	35
Naphthalene	ND		5000	5350		ug/L		107	56 - 161	1	30
n-Propylbenzene	ND		5000	4820		ug/L		96	61 - 167	1	23
Styrene	ND		5000	5360		ug/L		107	69 - 150	1	29
1,1,1,2-Tetrachloroethane	ND		5000	5690		ug/L		114	80 - 140	0.8	11
1,1,2,2-Tetrachloroethane	ND		5000	4600		ug/L		92	76 - 141	0.7	28
Tetrachloroethene	ND		5000	5690		ug/L		114	63 - 155	3	16
Toluene	73.0		5000	5020		ug/L		99	61 - 153	2	35
1,2,3-Trichlorobenzene	ND		5000	5890		ug/L		118	57 - 155	5	28
1,2,4-Trichlorobenzene	ND		5000	5890		ug/L		118	64 - 147	2	23
1,1,2-Trichloroethane	ND		5000	5080		ug/L		102	74 - 138	1	21
1,1,1-Trichloroethane	ND		5000	5750		ug/L		115	78 - 153	2	29
Trichloroethene	ND		5000	5890		ug/L		118	74 - 139	2	11
Trichlorofluoromethane	ND		5000	5030		ug/L		101	53 - 149	3	33
1,2,3-Trichloropropane	ND		5000	4590		ug/L		92	49 - 148	0.4	25
1,3,5-Trimethylbenzene	ND		5000	5000		ug/L		100	67 - 151	0.8	21
1,2,4-Trimethylbenzene	47.0		5000	4880		ug/L		97	69 - 150	1	20
Vinyl chloride	ND		5000	4040		ug/L		81	53 - 137	1	32
Xylenes, total	41400		15000	49100	M8	ug/L		51	68 - 158	0.9	18

Surrogate	*Matrix Spike Dup % Recovery*	*Matrix Spike Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*95*		*63 - 140*
Dibromofluoromethane	*107*		*73 - 131*
Toluene-d8	*94*		*80 - 120*
4-Bromofluorobenzene	*91*		*79 - 125*

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-BLK1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Method Blank
Prep Type: Total
Prep Batch: 11H2289_P

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Acetone	ND		50.0		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Tert-Amyl Methyl Ether	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Benzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Bromobenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Bromochloromethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Ethyl tert-Butyl Ether	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Bromodichloromethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Diisopropyl Ether	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Methyl tert-Butyl Ether	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Bromoform	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Tertiary Butyl Alcohol	ND		20.0		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Bromomethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
2-Butanone	ND		50.0		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
sec-Butylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
n-Butylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
tert-Butylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Carbon disulfide	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Carbon Tetrachloride	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Chlorobenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Chlorodibromomethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Chloroethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Chloroform	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Chloromethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
2-Chlorotoluene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
4-Chlorotoluene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2-Dibromo-3-chloropropane	ND		5.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2-Dibromoethane (EDB)	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Dibromomethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,4-Dichlorobenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,3-Dichlorobenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2-Dichlorobenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Dichlorodifluoromethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,1-Dichloroethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2-Dichloroethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
cis-1,2-Dichloroethene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,1-Dichloroethene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
trans-1,2-Dichloroethene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,3-Dichloropropane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2-Dichloropropane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
2,2-Dichloropropane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
cis-1,3-Dichloropropene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
trans-1,3-Dichloropropene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,1-Dichloropropene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Ethylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Hexachlorobutadiene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
2-Hexanone	ND		50.0		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Isopropylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
p-Isopropyltoluene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Methylene Chloride	ND		5.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-BLK1 | **Client Sample ID: Method Blank**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014193 | **Prep Batch: 11H2289_P**

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
4-Methyl-2-pentanone	ND		10.0		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Naphthalene	ND		5.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
n-Propylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Styrene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,1,1,2-Tetrachloroethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,1,2,2-Tetrachloroethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Tetrachloroethene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Toluene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2,3-Trichlorobenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2,4-Trichlorobenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,1,2-Trichloroethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,1,1-Trichloroethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Trichloroethene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Trichlorofluoromethane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2,3-Trichloropropane	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,3,5-Trimethylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
1,2,4-Trimethylbenzene	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Vinyl chloride	ND		1.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00
Xylenes, total	ND		3.00		ug/L		08/10/11 06:14	08/10/11 09:02	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
1,2-Dichloroethane-d4	*101*		*63 - 140*	*08/10/11 06:14*	*08/10/11 09:02*	*1.00*
Dibromofluoromethane	*107*		*73 - 131*	*08/10/11 06:14*	*08/10/11 09:02*	*1.00*
Toluene-d8	*92*		*80 - 120*	*08/10/11 06:14*	*08/10/11 09:02*	*1.00*
4-Bromofluorobenzene	*89*		*79 - 125*	*08/10/11 06:14*	*08/10/11 09:02*	*1.00*

Lab Sample ID: 11H2289-BS1 | **Client Sample ID: Lab Control Sample**
Matrix: Water | **Prep Type: Total**
Analysis Batch: U014193 | **Prep Batch: 11H2289_P**

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	250	234		ug/L		93	56 - 150
Tert-Amyl Methyl Ether	50.0	47.6		ug/L		95	70 - 133
Benzene	50.0	51.3		ug/L		103	80 - 121
Bromobenzene	50.0	44.6		ug/L		89	72 - 130
Bromochloromethane	50.0	58.6		ug/L		117	73 - 137
Ethyl tert-Butyl Ether	50.0	44.9		ug/L		90	68 - 138
Bromodichloromethane	50.0	55.3		ug/L		111	75 - 131
Diisopropyl Ether	50.0	41.3		ug/L		83	63 - 136
Methyl tert-Butyl Ether	50.0	49.8		ug/L		100	76 - 120
Bromoform	50.0	55.2		ug/L		110	65 - 140
Tertiary Butyl Alcohol	500	546		ug/L		109	60 - 140
Bromomethane	50.0	45.1		ug/L		90	50 - 150
2-Butanone	250	264		ug/L		105	70 - 144
sec-Butylbenzene	50.0	50.4		ug/L		101	72 - 140
n-Butylbenzene	50.0	47.1		ug/L		94	68 - 140
tert-Butylbenzene	50.0	50.9		ug/L		102	76 - 135
Carbon disulfide	50.0	46.8		ug/L		94	74 - 137
Carbon Tetrachloride	50.0	64.4		ug/L		129	71 - 137

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-BS1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H2289_P

Analyte	Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
Chlorobenzene	50.0	53.6		ug/L		107	80 - 121
Chlorodibromomethane	50.0	61.4		ug/L		123	68 - 137
Chloroethane	50.0	39.8		ug/L		80	50 - 146
Chloroform	50.0	55.2		ug/L		110	73 - 131
Chloromethane	50.0	30.8		ug/L		62	30 - 132
2-Chlorotoluene	50.0	46.7		ug/L		93	74 - 135
4-Chlorotoluene	50.0	48.7		ug/L		97	74 - 132
1,2-Dibromo-3-chloropropane	50.0	51.4		ug/L		103	56 - 145
1,2-Dibromoethane (EDB)	50.0	54.7		ug/L		109	80 - 135
Dibromomethane	50.0	55.9		ug/L		112	78 - 133
1,4-Dichlorobenzene	50.0	53.9		ug/L		108	80 - 120
1,3-Dichlorobenzene	50.0	53.0		ug/L		106	80 - 128
1,2-Dichlorobenzene	50.0	54.2		ug/L		108	80 - 125
Dichlorodifluoromethane	50.0	39.4		ug/L		79	30 - 132
1,1-Dichloroethane	50.0	49.8		ug/L		100	75 - 125
1,2-Dichloroethane	50.0	53.5		ug/L		107	70 - 134
cis-1,2-Dichloroethene	50.0	48.2		ug/L		96	71 - 132
1,1-Dichloroethene	50.0	53.9		ug/L		108	73 - 125
trans-1,2-Dichloroethene	50.0	50.6		ug/L		101	77 - 125
1,3-Dichloropropane	50.0	48.5		ug/L		97	76 - 125
1,2-Dichloropropane	50.0	46.4		ug/L		93	72 - 120
2,2-Dichloropropane	50.0	41.9		ug/L		84	50 - 150
cis-1,3-Dichloropropene	50.0	50.2		ug/L		100	70 - 140
trans-1,3-Dichloropropene	50.0	49.1		ug/L		98	62 - 139
1,1-Dichloropropene	50.0	52.3		ug/L		105	78 - 126
Ethylbenzene	50.0	50.5		ug/L		101	78 - 133
Hexachlorobutadiene	50.0	54.6		ug/L		109	70 - 150
2-Hexanone	250	238		ug/L		95	60 - 150
Isopropylbenzene	50.0	58.1		ug/L		116	69 - 120
p-Isopropyltoluene	50.0	50.4		ug/L		101	72 - 134
Methylene Chloride	50.0	53.7		ug/L		107	80 - 133
4-Methyl-2-pentanone	250	232		ug/L		93	62 - 146
Naphthalene	50.0	54.0		ug/L		108	71 - 139
n-Propylbenzene	50.0	47.3		ug/L		95	70 - 143
Styrene	50.0	54.1		ug/L		108	80 - 136
1,1,1,2-Tetrachloroethane	50.0	59.1		ug/L		118	80 - 130
1,1,2,2-Tetrachloroethane	50.0	46.2		ug/L		92	73 - 131
Tetrachloroethene	50.0	58.3		ug/L		117	77 - 131
Toluene	50.0	50.9		ug/L		102	78 - 125
1,2,3-Trichlorobenzene	50.0	58.2		ug/L		116	71 - 138
1,2,4-Trichlorobenzene	50.0	58.5		ug/L		117	74 - 136
1,1,2-Trichloroethane	50.0	52.9		ug/L		106	80 - 123
1,1,1-Trichloroethane	50.0	58.8		ug/L		118	75 - 137
Trichloroethene	50.0	60.6		ug/L		121	74 - 139
Trichlorofluoromethane	50.0	48.9		ug/L		98	60 - 133
1,2,3-Trichloropropane	50.0	46.2		ug/L		92	64 - 127
1,3,5-Trimethylbenzene	50.0	49.9		ug/L		100	75 - 134
1,2,4-Trimethylbenzene	50.0	48.4		ug/L		97	77 - 134
Vinyl chloride	50.0	39.2		ug/L		78	60 - 122
Xylenes, total	150	157		ug/L		105	78 - 134

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-BS1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Lab Control Sample
Prep Type: Total
Prep Batch: 11H2289_P

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
1,2-Dichloroethane-d4	*96*		*63 - 140*
Dibromofluoromethane	*108*		*73 - 131*
Toluene-d8	*94*		*80 - 120*
4-Bromofluorobenzene	*89*		*79 - 125*

Lab Sample ID: 11H2289-BSD1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Lab Control Sample Dup
Prep Type: Total
Prep Batch: 11H2289_P

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	250	211		ug/L		84	56 - 150	10	31
Tert-Amyl Methyl Ether	50.0	39.7	R2	ug/L		79	70 - 133	18	16
Benzene	50.0	43.6	R2	ug/L		87	80 - 121	16	12
Bromobenzene	50.0	38.5		ug/L		77	72 - 130	15	23
Bromochloromethane	50.0	50.3		ug/L		101	73 - 137	15	32
Ethyl tert-Butyl Ether	50.0	37.5	R2	ug/L		75	68 - 138	18	16
Bromodichloromethane	50.0	46.9	R2	ug/L		94	75 - 131	16	13
Diisopropyl Ether	50.0	35.2		ug/L		70	63 - 136	16	32
Methyl tert-Butyl Ether	50.0	41.6		ug/L		83	76 - 120	18	32
Bromoform	50.0	46.0		ug/L		92	65 - 140	18	18
Tertiary Butyl Alcohol	500	489		ug/L		98	60 - 140	11	30
Bromomethane	50.0	40.4		ug/L		81	50 - 150	11	50
2-Butanone	250	219		ug/L		88	70 - 144	18	37
sec-Butylbenzene	50.0	42.8		ug/L		86	72 - 140	16	21
n-Butylbenzene	50.0	40.3	R2	ug/L		81	68 - 140	16	11
tert-Butylbenzene	50.0	43.6		ug/L		87	76 - 135	15	20
Carbon disulfide	50.0	39.3		ug/L		79	74 - 137	17	28
Carbon Tetrachloride	50.0	53.6		ug/L		107	71 - 137	18	26
Chlorobenzene	50.0	45.9	R2	ug/L		92	80 - 121	16	11
Chlorodibromomethane	50.0	51.6	R2	ug/L		103	68 - 137	17	16
Chloroethane	50.0	37.7		ug/L		75	50 - 146	5	35
Chloroform	50.0	46.4		ug/L		93	73 - 131	17	32
Chloromethane	50.0	29.6		ug/L		59	30 - 132	4	34
2-Chlorotoluene	50.0	40.1		ug/L		80	74 - 135	15	22
4-Chlorotoluene	50.0	41.6		ug/L		83	74 - 132	16	22
1,2-Dibromo-3-chloropropane	50.0	44.4		ug/L		89	56 - 145	15	21
1,2-Dibromoethane (EDB)	50.0	45.9	R2	ug/L		92	80 - 135	18	10
Dibromomethane	50.0	47.1	R2	ug/L		94	78 - 133	17	11
1,4-Dichlorobenzene	50.0	46.5	R2	ug/L		93	80 - 120	15	10
1,3-Dichlorobenzene	50.0	46.1		ug/L		92	80 - 128	14	18
1,2-Dichlorobenzene	50.0	46.4	R2	ug/L		93	80 - 125	15	11
Dichlorodifluoromethane	50.0	36.8		ug/L		74	30 - 132	7	32
1,1-Dichloroethane	50.0	42.7		ug/L		85	75 - 125	15	34
1,2-Dichloroethane	50.0	45.1		ug/L		90	70 - 134	17	25
cis-1,2-Dichloroethene	50.0	41.2		ug/L		82	71 - 132	16	32
1,1-Dichloroethene	50.0	45.7		ug/L		91	73 - 125	16	31
trans-1,2-Dichloroethene	50.0	42.7		ug/L		85	77 - 125	17	32
1,3-Dichloropropane	50.0	41.5		ug/L		83	76 - 125	16	20
1,2-Dichloropropane	50.0	38.8	R2	ug/L		78	72 - 120	18	11

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-BSD1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Lab Control Sample Dup
Prep Type: Total
Prep Batch: 11H2289_P

Analyte	Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
2,2-Dichloropropane	50.0	33.9	R2	ug/L		68	50 - 150	21	11
cis-1,3-Dichloropropene	50.0	41.8		ug/L		84	70 - 140	18	35
trans-1,3-Dichloropropene	50.0	40.9		ug/L		82	62 - 139	18	26
1,1-Dichloropropene	50.0	44.8		ug/L		90	78 - 126	15	18
Ethylbenzene	50.0	43.3	R2	ug/L		87	78 - 133	15	12
Hexachlorobutadiene	50.0	47.3		ug/L		95	70 - 150	14	21
2-Hexanone	250	198		ug/L		79	60 - 150	19	20
Isopropylbenzene	50.0	49.6	R2	ug/L		99	69 - 120	16	15
p-Isopropyltoluene	50.0	43.2		ug/L		86	72 - 134	16	18
Methylene Chloride	50.0	45.2		ug/L		90	80 - 133	17	36
4-Methyl-2-pentanone	250	195		ug/L		78	62 - 146	17	35
Naphthalene	50.0	46.4		ug/L		93	71 - 139	15	30
n-Propylbenzene	50.0	40.6		ug/L		81	70 - 143	15	23
Styrene	50.0	46.0		ug/L		92	80 - 136	16	29
1,1,1,2-Tetrachloroethane	50.0	50.1	R2	ug/L		100	80 - 130	16	11
1,1,2,2-Tetrachloroethane	50.0	39.4		ug/L		79	73 - 131	16	28
Tetrachloroethene	50.0	49.3	R2	ug/L		99	77 - 131	17	16
Toluene	50.0	43.2		ug/L		86	78 - 125	16	35
1,2,3-Trichlorobenzene	50.0	51.2		ug/L		102	71 - 138	13	28
1,2,4-Trichlorobenzene	50.0	50.2		ug/L		100	74 - 136	15	23
1,1,2-Trichloroethane	50.0	45.0		ug/L		90	80 - 123	16	21
1,1,1-Trichloroethane	50.0	49.4		ug/L		99	75 - 137	18	29
Trichloroethene	50.0	51.7	R2	ug/L		103	74 - 139	16	11
Trichlorofluoromethane	50.0	45.4		ug/L		91	60 - 133	8	33
1,2,3-Trichloropropane	50.0	39.3		ug/L		79	64 - 127	16	25
1,3,5-Trimethylbenzene	50.0	42.3		ug/L		85	75 - 134	16	21
1,2,4-Trimethylbenzene	50.0	41.6		ug/L		83	77 - 134	15	20
Vinyl chloride	50.0	37.2		ug/L		74	60 - 122	5	32
Xylenes, total	150	133		ug/L		89	78 - 134	17	18

Surrogate	*LCS Dup % Recovery*	*LCS Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*96*		*63 - 140*
Dibromofluoromethane	*106*		*73 - 131*
Toluene-d8	*94*		*80 - 120*
4-Bromofluorobenzene	*89*		*79 - 125*

Lab Sample ID: 11H2289-MS1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Matrix Spike
Prep Type: Total
Prep Batch: 11H2289_P

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Acetone	ND		250	214		ug/L		86	56 - 150
Tert-Amyl Methyl Ether	ND		50.0	43.9		ug/L		88	69 - 139
Benzene	ND		50.0	50.2		ug/L		100	65 - 151
Bromobenzene	ND		50.0	40.4		ug/L		81	69 - 142
Bromochloromethane	ND		50.0	57.5		ug/L		115	64 - 154
Ethyl tert-Butyl Ether	ND		50.0	42.0		ug/L		84	68 - 139
Bromodichloromethane	ND		50.0	55.6		ug/L		111	75 - 138
Diisopropyl Ether	ND		50.0	39.4		ug/L		79	59 - 145

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-MS1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Matrix Spike
Prep Type: Total
Prep Batch: 11H2289_P

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
Methyl tert-Butyl Ether	ND		50.0	47.6		ug/L		95	56 - 152
Bromoform	ND		50.0	52.5		ug/L		105	55 - 153
Tertiary Butyl Alcohol	ND		500	466		ug/L		93	14 - 200
Bromomethane	ND		50.0	40.1		ug/L		80	13 - 176
2-Butanone	ND		250	236		ug/L		95	45 - 164
sec-Butylbenzene	ND		50.0	47.1		ug/L		94	68 - 159
n-Butylbenzene	ND		50.0	43.5		ug/L		87	67 - 151
tert-Butylbenzene	ND		50.0	47.7		ug/L		95	73 - 153
Carbon disulfide	ND		50.0	44.4		ug/L		89	33 - 187
Carbon Tetrachloride	ND		50.0	68.9		ug/L		138	64 - 157
Chlorobenzene	ND		50.0	51.4		ug/L		103	78 - 136
Chlorodibromomethane	ND		50.0	59.3		ug/L		119	64 - 145
Chloroethane	ND		50.0	41.9		ug/L		84	48 - 159
Chloroform	ND		50.0	55.9		ug/L		112	72 - 145
Chloromethane	ND		50.0	35.6		ug/L		71	10 - 194
2-Chlorotoluene	ND		50.0	42.5		ug/L		85	66 - 155
4-Chlorotoluene	ND		50.0	45.0		ug/L		90	69 - 149
1,2-Dibromo-3-chloropropane	ND		50.0	45.6		ug/L		91	49 - 162
1,2-Dibromoethane (EDB)	ND		50.0	50.7		ug/L		101	70 - 152
Dibromomethane	ND		50.0	55.2		ug/L		110	75 - 141
1,4-Dichlorobenzene	ND		50.0	50.3		ug/L		101	75 - 135
1,3-Dichlorobenzene	ND		50.0	49.2		ug/L		98	72 - 146
1,2-Dichlorobenzene	ND		50.0	50.0		ug/L		100	80 - 136
Dichlorodifluoromethane	ND		50.0	38.5		ug/L		77	23 - 159
1,1-Dichloroethane	ND		50.0	49.7		ug/L		99	64 - 154
1,2-Dichloroethane	ND		50.0	55.2		ug/L		110	72 - 137
cis-1,2-Dichloroethene	ND		50.0	48.0		ug/L		96	57 - 154
1,1-Dichloroethene	ND		50.0	53.1		ug/L		106	34 - 151
trans-1,2-Dichloroethene	ND		50.0	50.2		ug/L		100	57 - 157
1,3-Dichloropropane	ND		50.0	45.9		ug/L		92	71 - 137
1,2-Dichloropropane	ND		50.0	44.5		ug/L		89	71 - 139
2,2-Dichloropropane	ND		50.0	39.8		ug/L		80	10 - 198
cis-1,3-Dichloropropene	ND		50.0	46.1		ug/L		92	56 - 156
trans-1,3-Dichloropropene	ND		50.0	45.2		ug/L		90	47 - 157
1,1-Dichloropropene	ND		50.0	52.7		ug/L		105	70 - 155
Ethylbenzene	ND		50.0	49.3		ug/L		99	68 - 157
Hexachlorobutadiene	ND		50.0	50.9		ug/L		102	47 - 173
2-Hexanone	ND		250	215		ug/L		86	57 - 154
Isopropylbenzene	ND		50.0	57.6		ug/L		115	69 - 139
p-Isopropyltoluene	ND		50.0	47.5		ug/L		95	69 - 151
Methylene Chloride	ND		50.0	51.2		ug/L		102	71 - 136
4-Methyl-2-pentanone	ND		250	210		ug/L		84	62 - 159
Naphthalene	ND		50.0	46.4		ug/L		93	56 - 161
n-Propylbenzene	ND		50.0	43.5		ug/L		87	61 - 167
Styrene	ND		50.0	47.6		ug/L		95	69 - 150
1,1,1,2-Tetrachloroethane	ND		50.0	58.0		ug/L		116	80 - 140
1,1,2,2-Tetrachloroethane	ND		50.0	40.5		ug/L		81	76 - 141
Tetrachloroethene	0.760		50.0	58.2		ug/L		115	63 - 155
Toluene	ND		50.0	48.8		ug/L		98	61 - 153
1,2,3-Trichlorobenzene	ND		50.0	51.9		ug/L		104	57 - 155

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-MS1 — **Client Sample ID: Matrix Spike**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014193 — **Prep Batch: 11H2289_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
1,2,4-Trichlorobenzene	ND		50.0	51.3		ug/L		103	64 - 147
1,1,2-Trichloroethane	ND		50.0	50.6		ug/L		101	74 - 138
1,1,1-Trichloroethane	ND		50.0	61.9		ug/L		124	78 - 153
Trichloroethene	ND		50.0	59.8		ug/L		120	74 - 139
Trichlorofluoromethane	ND		50.0	55.2		ug/L		110	53 - 149
1,2,3-Trichloropropane	ND		50.0	41.0		ug/L		82	49 - 148
1,3,5-Trimethylbenzene	ND		50.0	45.8		ug/L		92	67 - 151
1,2,4-Trimethylbenzene	ND		50.0	44.7		ug/L		89	69 - 150
Vinyl chloride	ND		50.0	40.2		ug/L		80	53 - 137
Xylenes, total	ND		150	153		ug/L		102	68 - 158

Surrogate	*Matrix Spike % Recovery*	*Matrix Spike Qualifier*	*Limits*
1,2-Dichloroethane-d4	*105*		*63 - 140*
Dibromofluoromethane	*114*		*73 - 131*
Toluene-d8	*93*		*80 - 120*
4-Bromofluorobenzene	*83*		*79 - 125*

Lab Sample ID: 11H2289-MSD1 — **Client Sample ID: Matrix Spike Duplicate**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014193 — **Prep Batch: 11H2289_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Acetone	ND		250	178		ug/L		71	56 - 150	18	31
Tert-Amyl Methyl Ether	ND		50.0	36.8	R2	ug/L		74	69 - 139	17	16
Benzene	ND		50.0	42.8	R2	ug/L		86	65 - 151	16	12
Bromobenzene	ND		50.0	34.8		ug/L		70	69 - 142	15	23
Bromochloromethane	ND		50.0	48.1		ug/L		96	64 - 154	18	32
Ethyl tert-Butyl Ether	ND		50.0	35.8		ug/L		72	68 - 139	16	16
Bromodichloromethane	ND		50.0	46.2	R2	ug/L		92	75 - 138	18	13
Diisopropyl Ether	ND		50.0	34.2		ug/L		68	59 - 145	14	32
Methyl tert-Butyl Ether	ND		50.0	40.8		ug/L		82	56 - 152	15	32
Bromoform	ND		50.0	44.0		ug/L		88	55 - 153	18	18
Tertiary Butyl Alcohol	ND		500	481		ug/L		96	14 - 200	3	30
Bromomethane	ND		50.0	38.4		ug/L		77	13 - 176	4	50
2-Butanone	ND		250	204		ug/L		82	45 - 164	14	37
sec-Butylbenzene	ND		50.0	40.4		ug/L		81	68 - 159	15	21
n-Butylbenzene	ND		50.0	37.7	R2	ug/L		75	67 - 151	14	11
tert-Butylbenzene	ND		50.0	41.1		ug/L		82	73 - 153	15	20
Carbon disulfide	ND		50.0	38.2		ug/L		76	33 - 187	15	28
Carbon Tetrachloride	ND		50.0	57.8		ug/L		116	64 - 157	18	26
Chlorobenzene	ND		50.0	43.9	R2	ug/L		88	78 - 136	16	11
Chlorodibromomethane	ND		50.0	50.1	R2	ug/L		100	64 - 145	17	16
Chloroethane	ND		50.0	36.7		ug/L		73	48 - 159	13	35
Chloroform	ND		50.0	47.4		ug/L		95	72 - 145	17	32
Chloromethane	ND		50.0	30.2		ug/L		60	10 - 194	17	34
2-Chlorotoluene	ND		50.0	36.8		ug/L		74	66 - 155	14	22
4-Chlorotoluene	ND		50.0	38.6		ug/L		77	69 - 149	15	22
1,2-Dibromo-3-chloropropane	ND		50.0	38.4		ug/L		77	49 - 162	17	21
1,2-Dibromoethane (EDB)	ND		50.0	43.6	R2	ug/L		87	70 - 152	15	10

Method: SW846 8260B - Volatile Organic Compounds by EPA Method 8260B (Continued)

Lab Sample ID: 11H2289-MSD1
Matrix: Water
Analysis Batch: U014193

Client Sample ID: Matrix Spike Duplicate
Prep Type: Total
Prep Batch: 11H2289_P

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Dibromomethane	ND		50.0	46.4	R2	ug/L		93	75 - 141	17	11
1,4-Dichlorobenzene	ND		50.0	43.4	R2	ug/L		87	75 - 135	15	10
1,3-Dichlorobenzene	ND		50.0	42.1		ug/L		84	72 - 146	16	18
1,2-Dichlorobenzene	ND		50.0	42.6	R2	ug/L		85	80 - 136	16	11
Dichlorodifluoromethane	ND		50.0	32.2		ug/L		64	23 - 159	18	32
1,1-Dichloroethane	ND		50.0	42.4		ug/L		85	64 - 154	16	34
1,2-Dichloroethane	ND		50.0	46.6		ug/L		93	72 - 137	17	25
cis-1,2-Dichloroethene	ND		50.0	40.8		ug/L		82	57 - 154	16	32
1,1-Dichloroethene	ND		50.0	44.1		ug/L		88	34 - 151	18	31
trans-1,2-Dichloroethene	ND		50.0	43.2		ug/L		86	57 - 157	15	32
1,3-Dichloropropane	ND		50.0	39.5		ug/L		79	71 - 137	15	20
1,2-Dichloropropane	ND		50.0	37.8	R2	ug/L		76	71 - 139	16	11
2,2-Dichloropropane	ND		50.0	33.8	R2	ug/L		68	10 - 198	16	11
cis-1,3-Dichloropropene	ND		50.0	39.4		ug/L		79	56 - 156	16	35
trans-1,3-Dichloropropene	ND		50.0	38.4		ug/L		77	47 - 157	16	26
1,1-Dichloropropene	ND		50.0	44.7		ug/L		89	70 - 155	16	18
Ethylbenzene	ND		50.0	42.1	R2	ug/L		84	68 - 157	16	12
Hexachlorobutadiene	ND		50.0	45.4		ug/L		91	47 - 173	11	21
2-Hexanone	ND		250	183		ug/L		73	57 - 154	16	20
Isopropylbenzene	ND		50.0	48.2	R2	ug/L		96	69 - 139	18	15
p-Isopropyltoluene	ND		50.0	40.8		ug/L		82	69 - 151	15	18
Methylene Chloride	ND		50.0	44.4		ug/L		89	71 - 136	14	36
4-Methyl-2-pentanone	ND		250	180		ug/L		72	62 - 159	15	35
Naphthalene	ND		50.0	39.2		ug/L		78	56 - 161	17	30
n-Propylbenzene	ND		50.0	37.5		ug/L		75	61 - 167	15	23
Styrene	ND		50.0	39.6		ug/L		79	69 - 150	18	29
1,1,1,2-Tetrachloroethane	ND		50.0	49.1	R2	ug/L		98	80 - 140	17	11
1,1,2,2-Tetrachloroethane	ND		50.0	35.4	M8	ug/L		71	76 - 141	13	28
Tetrachloroethene	0.760		50.0	49.3	R2	ug/L		97	63 - 155	17	16
Toluene	ND		50.0	42.0		ug/L		84	61 - 153	15	35
1,2,3-Trichlorobenzene	ND		50.0	45.4		ug/L		91	57 - 155	13	28
1,2,4-Trichlorobenzene	ND		50.0	44.4		ug/L		89	64 - 147	14	23
1,1,2-Trichloroethane	ND		50.0	42.9		ug/L		86	74 - 138	17	21
1,1,1-Trichloroethane	ND		50.0	52.4		ug/L		105	78 - 153	17	29
Trichloroethene	ND		50.0	49.9	R2	ug/L		100	74 - 139	18	11
Trichlorofluoromethane	ND		50.0	45.9		ug/L		92	53 - 149	18	33
1,2,3-Trichloropropane	ND		50.0	35.4		ug/L		71	49 - 148	15	25
1,3,5-Trimethylbenzene	ND		50.0	39.3		ug/L		79	67 - 151	15	21
1,2,4-Trimethylbenzene	ND		50.0	38.4		ug/L		77	69 - 150	15	20
Vinyl chloride	ND		50.0	35.6		ug/L		71	53 - 137	12	32
Xylenes, total	ND		150	130		ug/L		86	68 - 158	17	18

Surrogate	*Matrix Spike Dup % Recovery*	*Matrix Spike Dup Qualifier*	*Limits*
1,2-Dichloroethane-d4	*101*		*63 - 140*
Dibromofluoromethane	*110*		*73 - 131*
Toluene-d8	*92*		*80 - 120*
4-Bromofluorobenzene	*86*		*79 - 125*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons

Lab Sample ID: 11H2053-BLK1 — **Client Sample ID: Method Blank**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/10/11 00:00	08/10/11 12:56	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*99*		*60 - 150*	*08/10/11 00:00*	*08/10/11 12:56*	*1.00*

Lab Sample ID: 11H2053-BLK2 — **Client Sample ID: Method Blank**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/10/11 00:00	08/10/11 23:52	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*96*		*60 - 150*	*08/10/11 00:00*	*08/10/11 23:52*	*1.00*

Lab Sample ID: 11H2053-BS1 — **Client Sample ID: Lab Control Sample**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits		
GRO as Gasoline			1000	1090		ug/L		109	64 - 134		

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
a,a,a-Trifluorotoluene	*96*		*60 - 150*

Lab Sample ID: 11H2053-BS2 — **Client Sample ID: Lab Control Sample**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits		
GRO as Gasoline			1000	1020		ug/L		102	64 - 134		

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
a,a,a-Trifluorotoluene	*89*		*60 - 150*

Lab Sample ID: 11H2053-MS1 — **Client Sample ID: MW-48 (08022011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits		
GRO as Gasoline	38.6		1000	1340		ug/L		131	49 - 161		

Surrogate	*Matrix Spike % Recovery*	*Matrix Spike Qualifier*	*Limits*
a,a,a-Trifluorotoluene	*91*		*60 - 150*

Method: SW846 8015B - Purgeable Petroleum Hydrocarbons (Continued)

Lab Sample ID: 11H2053-MS2 — **Client Sample ID: RB-2 (08032011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Result	Matrix Spike Qualifier	Unit	D	% Rec	% Rec. Limits
GRO as Gasoline	ND		1000	1230		ug/L		123	49 - 161

Surrogate	*Matrix Spike % Recovery*	*Matrix Spike Qualifier*	*Limits*
a,a,a-Trifluorotoluene	*89*		*60 - 150*

Lab Sample ID: 11H2053-MSD1 — **Client Sample ID: MW-48 (08022011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
GRO as Gasoline	38.6		1000	1190		ug/L		115	49 - 161	12	33

Surrogate	*Matrix Spike Dup % Recovery*	*Matrix Spike Dup Qualifier*	*Limits*
a,a,a-Trifluorotoluene	*94*		*60 - 150*

Lab Sample ID: 11H2053-MSD2 — **Client Sample ID: RB-2 (08032011)**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014275 — **Prep Batch: 11H2053_P**

Analyte	Sample Result	Sample Qualifier	Spike Added	Matrix Spike Dup Result	Matrix Spike Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
GRO as Gasoline	ND		1000	1250		ug/L		125	49 - 161	2	33

Surrogate	*Matrix Spike Dup % Recovery*	*Matrix Spike Dup Qualifier*	*Limits*
a,a,a-Trifluorotoluene	*91*		*60 - 150*

Lab Sample ID: 11H2720-BLK1 — **Client Sample ID: Method Blank**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014376 — **Prep Batch: 11H2720_P**

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
GRO as Gasoline	ND		100		ug/L		08/11/11 00:00	08/11/11 13:02	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
a,a,a-Trifluorotoluene	*94*		*60 - 150*	*08/11/11 00:00*	*08/11/11 13:02*	*1.00*

Lab Sample ID: 11H2720-BS1 — **Client Sample ID: Lab Control Sample**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014376 — **Prep Batch: 11H2720_P**

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits
GRO as Gasoline			1000	1160		ug/L		116	64 - 134

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
a,a,a-Trifluorotoluene	*90*		*60 - 150*

Method: SW846 8015B - Extractable Petroleum Hydrocarbons

Lab Sample ID: 11H1829-BLK1 — **Client Sample ID: Method Blank**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014235 — **Prep Batch: 11H1829_P**

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	ND		100		ug/L		08/08/11 11:05	08/10/11 13:35	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*86*		*53 - 140*	*08/08/11 11:05*	*08/10/11 13:35*	*1.00*

Lab Sample ID: 11H1829-BS1 — **Client Sample ID: Lab Control Sample**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014235 — **Prep Batch: 11H1829_P**

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits		
Diesel			1000	718	MNR1	ug/L		72	55 - 130		

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
o-Terphenyl	*112*		*53 - 140*

Lab Sample ID: 11H1829-BSD1 — **Client Sample ID: Lab Control Sample Dup**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014235 — **Prep Batch: 11H1829_P**

Analyte			Spike Added	LCS Dup Result	LCS Dup Qualifier	Unit	D	% Rec	% Rec. Limits	RPD	RPD Limit
Diesel			1000	742		ug/L		74	55 - 130	3	41

Surrogate	*LCS Dup % Recovery*	*LCS Dup Qualifier*	*Limits*
o-Terphenyl	*113*		*53 - 140*

Lab Sample ID: 11H1832-BLK1 — **Client Sample ID: Method Blank**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014117 — **Prep Batch: 11H1832_P**

Analyte	Blank Result	Blank Qualifier	RL	MDL	Unit	D	Prepared	Analyzed	Dil Fac
Diesel	ND		100		ug/L		08/08/11 12:05	08/09/11 06:24	1.00

Surrogate	*Blank % Recovery*	*Blank Qualifier*	*Limits*	*Prepared*	*Analyzed*	*Dil Fac*
o-Terphenyl	*100*		*53 - 140*	*08/08/11 12:05*	*08/09/11 06:24*	*1.00*

Lab Sample ID: 11H1832-BS1 — **Client Sample ID: Lab Control Sample**
Matrix: Water — **Prep Type: Total**
Analysis Batch: U014117 — **Prep Batch: 11H1832_P**

Analyte			Spike Added	LCS Result	LCS Qualifier	Unit	D	% Rec	% Rec. Limits		
Diesel			1000	1000	MNR1	ug/L		100	55 - 130		

Surrogate	*LCS % Recovery*	*LCS Qualifier*	*Limits*
o-Terphenyl	*116*		*53 - 140*

GCMS Volatiles

Analysis Batch: U014048

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1881-BS1	Lab Control Sample	Total	Water	SW846 8260B	11H1881_P
11H1881-BSD1	Lab Control Sample Dup	Total	Water	SW846 8260B	11H1881_P
11H1881-BLK1	Method Blank	Total	Water	SW846 8260B	11H1881_P
NUH0938-27	TB-1 (08032011)	Total	Water	SW846 8260B	11H1881_P
NUH0938-28	TB-2 (08032011)	Total	Water	SW846 8260B	11H1881_P
NUH0938-25	RB-1 (08022011)	Total	Ground Water	SW846 8260B	11H1881_P
NUH0938-26	RB-2 (08032011)	Total	Ground Water	SW846 8260B	11H1881_P
NUH0938-03	MW-13A (08022011)	Total	Ground Water	SW846 8260B	11H1881_P
NUH0938-04	MW-14A (08022011)	Total	Ground Water	SW846 8260B	11H1881_P
NUH0938-05	MW-37 (08022011)	Total	Ground Water	SW846 8260B	11H1881_P
NUH0938-06	MW-38 (08022011)	Total	Ground Water	SW846 8260B	11H1881_P
11H1881-MS1	MW-38 (08022011)	Total	Water	SW846 8260B	11H1881_P
11H1881-MSD1	MW-38 (08022011)	Total	Water	SW846 8260B	11H1881_P

Analysis Batch: U014132

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1622-BS1	Lab Control Sample	Total	Water	SW846 8260B	11H1622_P
11H1622-BSD1	Lab Control Sample Dup	Total	Water	SW846 8260B	11H1622_P
11H1622-BLK1	Method Blank	Total	Water	SW846 8260B	11H1622_P
NUH0938-07	MW-46 (08022011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-08	MW-48 (08022011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-09	MW-50 (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-10	MW-51 (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-11	MW-53D (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-12	MW-53I (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-13	MW-53S (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-14	MW-54D (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-16	MW-55S (08022011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-17	RW-52 (08022011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-18	VW-1 (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-20	VW-3 (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-21	MP-1 (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-22	DPE-1 (08032011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-23	DUP-1 (08022011)	Total	Ground Water	SW846 8260B	11H1622_P
NUH0938-29	MW-55I (08022011)	Total	Ground Water	SW846 8260B	11H1622_P
11H1622-MS1	MW-46 (08022011)	Total	Water	SW846 8260B	11H1622_P
11H1622-MSD1	MW-46 (08022011)	Total	Water	SW846 8260B	11H1622_P

Analysis Batch: U014133

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2288-BS1	Lab Control Sample	Total	Water	SW846 8260B	11H2288_P
11H2288-BSD1	Lab Control Sample Dup	Total	Water	SW846 8260B	11H2288_P
11H2288-BLK1	Method Blank	Total	Water	SW846 8260B	11H2288_P
NUH0938-01 - RE1	MW-5 (08032011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-02 - RE1	MW-6 (08032011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-15 - RE1	MW-55D (08022011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-19 - RE1	VW-2 (08032011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-24 - RE1	DUP-2 (08032011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-03 - RE1	MW-13A (08022011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-04 - RE1	MW-14A (08022011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-12 - RE1	MW-53I (08032011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-14 - RE1	MW-54D (08032011)	Total	Ground Water	SW846 8260B	11H2288_P

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Analysis Batch: U014133 (Continued)

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
NUH0938-16 - RE1	MW-55S (08022011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-23 - RE1	DUP-1 (08022011)	Total	Ground Water	SW846 8260B	11H2288_P
NUH0938-18 - RE1	VW-1 (08032011)	Total	Ground Water	SW846 8260B	11H2288_P
11H2288-MS1	Matrix Spike	Total	Water	SW846 8260B	11H2288_P
11H2288-MSD1	Matrix Spike Duplicate	Total	Water	SW846 8260B	11H2288_P

Analysis Batch: U014193

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2289-BS1	Lab Control Sample	Total	Water	SW846 8260B	11H2289_P
11H2289-BSD1	Lab Control Sample Dup	Total	Water	SW846 8260B	11H2289_P
11H2289-BLK1	Method Blank	Total	Water	SW846 8260B	11H2289_P
NUH0938-13 - RE1	MW-53S (08032011)	Total	Ground Water	SW846 8260B	11H2289_P
NUH0938-13 - RE2	MW-53S (08032011)	Total	Ground Water	SW846 8260B	11H2289_P
NUH0938-21 - RE1	MP-1 (08032011)	Total	Ground Water	SW846 8260B	11H2289_P
NUH0938-21 - RE2	MP-1 (08032011)	Total	Ground Water	SW846 8260B	11H2289_P
NUH0938-22 - RE1	DPE-1 (08032011)	Total	Ground Water	SW846 8260B	11H2289_P
NUH0938-22 - RE2	DPE-1 (08032011)	Total	Ground Water	SW846 8260B	11H2289_P
NUH0938-29 - RE1	MW-55I (08022011)	Total	Ground Water	SW846 8260B	11H2289_P
NUH0938-29 - RE2	MW-55I (08022011)	Total	Ground Water	SW846 8260B	11H2289_P
11H2289-MS1	Matrix Spike	Total	Water	SW846 8260B	11H2289_P
11H2289-MSD1	Matrix Spike Duplicate	Total	Water	SW846 8260B	11H2289_P

Prep Batch: 11H1622_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1622-BS1	Lab Control Sample	Total	Water	EPA 5030B	
11H1622-BSD1	Lab Control Sample Dup	Total	Water	EPA 5030B	
11H1622-BLK1	Method Blank	Total	Water	EPA 5030B	
NUH0938-07	MW-46 (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-08	MW-48 (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-09	MW-50 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-10	MW-51 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-11	MW-53D (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-12	MW-53I (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-13	MW-53S (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-14	MW-54D (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-16	MW-55S (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-17	RW-52 (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-18	VW-1 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-20	VW-3 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-21	MP-1 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-22	DPE-1 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-23	DUP-1 (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-29	MW-55I (08022011)	Total	Ground Water	EPA 5030B	
11H1622-MS1	MW-46 (08022011)	Total	Water	EPA 5030B	
11H1622-MSD1	MW-46 (08022011)	Total	Water	EPA 5030B	

Prep Batch: 11H1881_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1881-BS1	Lab Control Sample	Total	Water	EPA 5030B	
11H1881-BSD1	Lab Control Sample Dup	Total	Water	EPA 5030B	
11H1881-BLK1	Method Blank	Total	Water	EPA 5030B	
NUH0938-27	TB-1 (08032011)	Total	Water	EPA 5030B	

GCMS Volatiles (Continued)

Prep Batch: 11H1881_P (Continued)

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
NUH0938-28	TB-2 (08032011)	Total	Water	EPA 5030B	
NUH0938-25	RB-1 (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-26	RB-2 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-03	MW-13A (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-04	MW-14A (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-05	MW-37 (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-06	MW-38 (08022011)	Total	Ground Water	EPA 5030B	
11H1881-MS1	MW-38 (08022011)	Total	Water	EPA 5030B	
11H1881-MSD1	MW-38 (08022011)	Total	Water	EPA 5030B	

Prep Batch: 11H2288_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2288-BS1	Lab Control Sample	Total	Water	EPA 5030B	
11H2288-BSD1	Lab Control Sample Dup	Total	Water	EPA 5030B	
11H2288-BLK1	Method Blank	Total	Water	EPA 5030B	
NUH0938-01 - RE1	MW-5 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-02 - RE1	MW-6 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-15 - RE1	MW-55D (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-19 - RE1	VW-2 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-24 - RE1	DUP-2 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-03 - RE1	MW-13A (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-04 - RE1	MW-14A (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-12 - RE1	MW-53I (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-14 - RE1	MW-54D (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-16 - RE1	MW-55S (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-23 - RE1	DUP-1 (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-18 - RE1	VW-1 (08032011)	Total	Ground Water	EPA 5030B	
11H2288-MS1	Matrix Spike	Total	Water	EPA 5030B	
11H2288-MSD1	Matrix Spike Duplicate	Total	Water	EPA 5030B	

Prep Batch: 11H2289_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2289-BS1	Lab Control Sample	Total	Water	EPA 5030B	
11H2289-BSD1	Lab Control Sample Dup	Total	Water	EPA 5030B	
11H2289-BLK1	Method Blank	Total	Water	EPA 5030B	
NUH0938-13 - RE1	MW-53S (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-13 - RE2	MW-53S (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-21 - RE1	MP-1 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-21 - RE2	MP-1 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-22 - RE1	DPE-1 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-22 - RE2	DPE-1 (08032011)	Total	Ground Water	EPA 5030B	
NUH0938-29 - RE1	MW-55I (08022011)	Total	Ground Water	EPA 5030B	
NUH0938-29 - RE2	MW-55I (08022011)	Total	Ground Water	EPA 5030B	
11H2289-MS1	Matrix Spike	Total	Water	EPA 5030B	
11H2289-MSD1	Matrix Spike Duplicate	Total	Water	EPA 5030B	

GC Volatiles

Analysis Batch: U014275

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2053-BS1	Lab Control Sample	Total	Water	SW846 8015B	11H2053_P
11H2053-BLK1	Method Blank	Total	Water	SW846 8015B	11H2053_P

GC Volatiles (Continued)

Analysis Batch: U014275 (Continued)

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
NUH0938-01	MW-5 (08032011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-02	MW-6 (08032011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-05	MW-37 (08022011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-06	MW-38 (08022011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-07	MW-46 (08022011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-08	MW-48 (08022011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-09	MW-50 (08032011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-10	MW-51 (08032011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-16	MW-55S (08022011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-17	RW-52 (08022011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-19	VW-2 (08032011)	Total	Ground Water	SW846 8015B	11H2053_P
11H2053-BLK2	Method Blank	Total	Water	SW846 8015B	11H2053_P
NUH0938-25	RB-1 (08022011)	Total	Ground Water	SW846 8015B	11H2053_P
NUH0938-26	RB-2 (08032011)	Total	Ground Water	SW846 8015B	11H2053_P
11H2053-BS2	Lab Control Sample	Total	Water	SW846 8015B	11H2053_P
11H2053-MS1	MW-48 (08022011)	Total	Water	SW846 8015B	11H2053_P
11H2053-MSD1	MW-48 (08022011)	Total	Water	SW846 8015B	11H2053_P
11H2053-MS2	RB-2 (08032011)	Total	Water	SW846 8015B	11H2053_P
11H2053-MSD2	RB-2 (08032011)	Total	Water	SW846 8015B	11H2053_P

Analysis Batch: U014376

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2720-BLK1	Method Blank	Total	Water	SW846 8015B	11H2720_P
NUH0938-11 - RE1	MW-53D (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-20 - RE1	VW-3 (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-24 - RE1	DUP-2 (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-15 - RE1	MW-55D (08022011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-13 - RE1	MW-53S (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-21 - RE1	MP-1 (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-22 - RE1	DPE-1 (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-29 - RE1	MW-55I (08022011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-14 - RE1	MW-54D (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-12 - RE1	MW-53I (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-18 - RE1	VW-1 (08032011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-03 - RE1	MW-13A (08022011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-04 - RE1	MW-14A (08022011)	Total	Ground Water	SW846 8015B	11H2720_P
NUH0938-23 - RE1	DUP-1 (08022011)	Total	Ground Water	SW846 8015B	11H2720_P
11H2720-BS1	Lab Control Sample	Total	Water	SW846 8015B	11H2720_P

Prep Batch: 11H2053_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2053-BS1	Lab Control Sample	Total	Water	EPA 5030B (GC)	
11H2053-BLK1	Method Blank	Total	Water	EPA 5030B (GC)	
NUH0938-01	MW-5 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-02	MW-6 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-05	MW-37 (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-06	MW-38 (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-07	MW-46 (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-08	MW-48 (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-09	MW-50 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-10	MW-51 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-16	MW-55S (08022011)	Total	Ground Water	EPA 5030B (GC)	

GC Volatiles (Continued)

Prep Batch: 11H2053_P (Continued)

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
NUH0938-17	RW-52 (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-19	VW-2 (08032011)	Total	Ground Water	EPA 5030B (GC)	
11H2053-BLK2	Method Blank	Total	Water	EPA 5030B (GC)	
NUH0938-25	RB-1 (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-26	RB-2 (08032011)	Total	Ground Water	EPA 5030B (GC)	
11H2053-BS2	Lab Control Sample	Total	Water	EPA 5030B (GC)	
11H2053-MS1	MW-48 (08022011)	Total	Water	EPA 5030B (GC)	
11H2053-MSD1	MW-48 (08022011)	Total	Water	EPA 5030B (GC)	
11H2053-MS2	RB-2 (08032011)	Total	Water	EPA 5030B (GC)	
11H2053-MSD2	RB-2 (08032011)	Total	Water	EPA 5030B (GC)	

Prep Batch: 11H2720_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H2720-BLK1	Method Blank	Total	Water	EPA 5030B (GC)	
NUH0938-11 - RE1	MW-53D (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-20 - RE1	VW-3 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-24 - RE1	DUP-2 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-15 - RE1	MW-55D (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-13 - RE1	MW-53S (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-21 - RE1	MP-1 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-22 - RE1	DPE-1 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-29 - RE1	MW-55I (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-14 - RE1	MW-54D (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-12 - RE1	MW-53I (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-18 - RE1	VW-1 (08032011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-03 - RE1	MW-13A (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-04 - RE1	MW-14A (08022011)	Total	Ground Water	EPA 5030B (GC)	
NUH0938-23 - RE1	DUP-1 (08022011)	Total	Ground Water	EPA 5030B (GC)	
11H2720-BS1	Lab Control Sample	Total	Water	EPA 5030B (GC)	

GC Semivolatiles

Analysis Batch: U014117

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1832-BLK1	Method Blank	Total	Water	SW846 8015B	11H1832_P
11H1832-BS1	Lab Control Sample	Total	Water	SW846 8015B	11H1832_P
NUH0938-01	MW-5 (08032011)	Total	Ground Water	SW846 8015B	11H1832_P
NUH0938-02	MW-6 (08032011)	Total	Ground Water	SW846 8015B	11H1832_P
NUH0938-04	MW-14A (08022011)	Total	Ground Water	SW846 8015B	11H1832_P

Analysis Batch: U014159

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
NUH0938-05	MW-37 (08022011)	Total	Ground Water	SW846 8015B	11H1832_P
NUH0938-06	MW-38 (08022011)	Total	Ground Water	SW846 8015B	11H1832_P
NUH0938-03	MW-13A (08022011)	Total	Ground Water	SW846 8015B	11H1832_P
NUH0938-07	MW-46 (08022011)	Total	Ground Water	SW846 8015B	11H1832_P
NUH0938-08	MW-48 (08022011)	Total	Ground Water	SW846 8015B	11H1832_P

Analysis Batch: U014235

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1829-BLK1	Method Blank	Total	Water	SW846 8015B	11H1829_P
11H1829-BS1	Lab Control Sample	Total	Water	SW846 8015B	11H1829_P

GC Semivolatiles (Continued)

Analysis Batch: U014235 (Continued)

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1829-BSD1	Lab Control Sample Dup	Total	Water	SW846 8015B	11H1829_P
NUH0938-09	MW-50 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-10	MW-51 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-11	MW-53D (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-12	MW-53I (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-15	MW-55D (08022011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-16	MW-55S (08022011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-17	RW-52 (08022011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-19	VW-2 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-20	VW-3 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-23	DUP-1 (08022011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-24	DUP-2 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-25	RB-1 (08022011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-26	RB-2 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-13 - RE1	MW-53S (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-14 - RE1	MW-54D (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-18 - RE1	VW-1 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-21 - RE1	MP-1 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-22 - RE1	DPE-1 (08032011)	Total	Ground Water	SW846 8015B	11H1829_P
NUH0938-29 - RE1	MW-55I (08022011)	Total	Ground Water	SW846 8015B	11H1829_P

Prep Batch: 11H1829_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1829-BLK1	Method Blank	Total	Water	EPA 3510C	
11H1829-BS1	Lab Control Sample	Total	Water	EPA 3510C	
11H1829-BSD1	Lab Control Sample Dup	Total	Water	EPA 3510C	
NUH0938-09	MW-50 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-10	MW-51 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-11	MW-53D (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-12	MW-53I (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-15	MW-55D (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-16	MW-55S (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-17	RW-52 (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-19	VW-2 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-20	VW-3 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-23	DUP-1 (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-24	DUP-2 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-25	RB-1 (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-26	RB-2 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-13 - RE1	MW-53S (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-14 - RE1	MW-54D (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-18 - RE1	VW-1 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-21 - RE1	MP-1 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-22 - RE1	DPE-1 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-29 - RE1	MW-55I (08022011)	Total	Ground Water	EPA 3510C	

Prep Batch: 11H1832_P

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
11H1832-BLK1	Method Blank	Total	Water	EPA 3510C	
11H1832-BS1	Lab Control Sample	Total	Water	EPA 3510C	
NUH0938-01	MW-5 (08032011)	Total	Ground Water	EPA 3510C	
NUH0938-02	MW-6 (08032011)	Total	Ground Water	EPA 3510C	

GC Semivolatiles (Continued)

Prep Batch: 11H1832_P (Continued)

Lab Sample ID	Client Sample ID	Prep Type	Matrix	Method	Prep Batch
NUH0938-04	MW-14A (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-05	MW-37 (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-06	MW-38 (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-03	MW-13A (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-07	MW-46 (08022011)	Total	Ground Water	EPA 3510C	
NUH0938-08	MW-48 (08022011)	Total	Ground Water	EPA 3510C	

Client Sample ID: MW-5 (08032011)

Lab Sample ID: NUH0938-01

Date Collected: 08/03/11 12:40

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	1.00	U014133	08/09/11 21:49	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/03/11 12:40	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 13:54	GWM	TAL NSH
Total	Prep	EPA 3510C		0.990	11H1832_P	08/08/11 12:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014117	08/09/11 07:36	KKH	TAL NSH

Client Sample ID: MW-6 (08032011)

Lab Sample ID: NUH0938-02

Date Collected: 08/03/11 15:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	1.00	U014133	08/09/11 22:17	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/03/11 15:05	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 14:24	GWM	TAL NSH
Total	Prep	EPA 3510C		0.943	11H1832_P	08/08/11 12:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014117	08/09/11 08:00	KKH	TAL NSH

Client Sample ID: MW-13A (08022011)

Lab Sample ID: NUH0938-03

Date Collected: 08/02/11 20:15

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 04:02	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014133	08/10/11 00:09	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/02/11 20:15	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	10.0	U014376	08/11/11 19:01	GWM	TAL NSH
Total	Prep	EPA 3510C		0.943	11H1832_P	08/08/11 12:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014159	08/09/11 21:18	KKH	TAL NSH

Client Sample ID: MW-14A (08022011)

Lab Sample ID: NUH0938-04

Date Collected: 08/02/11 20:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 04:30	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014133	08/10/11 00:37	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/02/11 20:05	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	10.0	U014376	08/11/11 19:31	GWM	TAL NSH
Total	Prep	EPA 3510C		0.943	11H1832_P	08/08/11 12:05	JJR	TAL NSH

Client Sample ID: MW-14A (08022011)

Lab Sample ID: NUH0938-04

Date Collected: 08/02/11 20:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Analysis	SW846 8015B		1.00	U014117	08/09/11 08:24	KKH	TAL NSH

Client Sample ID: MW-37 (08022011)

Lab Sample ID: NUH0938-05

Date Collected: 08/02/11 15:20

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 04:58	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/02/11 15:20	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 15:54	GWM	TAL NSH
Total	Prep	EPA 3510C		1.11	11H1832_P	08/08/11 12:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014159	08/09/11 18:43	KKH	TAL NSH

Client Sample ID: MW-38 (08022011)

Lab Sample ID: NUH0938-06

Date Collected: 08/02/11 16:35

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 05:26	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/02/11 16:35	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 16:24	GWM	TAL NSH
Total	Prep	EPA 3510C		0.971	11H1832_P	08/08/11 12:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014159	08/09/11 19:02	KKH	TAL NSH

Client Sample ID: MW-46 (08022011)

Lab Sample ID: NUH0938-07

Date Collected: 08/02/11 12:45

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 09:10	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/02/11 12:45	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 16:54	GWM	TAL NSH
Total	Prep	EPA 3510C		0.952	11H1832_P	08/08/11 12:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014159	08/09/11 21:38	KKH	TAL NSH

Client Sample ID: MW-48 (08022011)

Lab Sample ID: NUH0938-08

Date Collected: 08/02/11 18:00

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 09:38	CMM	TAL NSH

Client Sample ID: MW-48 (08022011)

Lab Sample ID: NUH0938-08

Date Collected: 08/02/11 18:00

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/02/11 18:00	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 17:24	GWM	TAL NSH
Total	Prep	EPA 3510C		0.990	11H1832_P	08/08/11 12:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014159	08/09/11 21:57	KKH	TAL NSH

Client Sample ID: MW-50 (08032011)

Lab Sample ID: NUH0938-09

Date Collected: 08/03/11 09:30

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 10:06	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/03/11 09:30	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 17:53	GWM	TAL NSH
Total	Prep	EPA 3510C		1.00	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 14:48	KKH	TAL NSH

Client Sample ID: MW-51 (08032011)

Lab Sample ID: NUH0938-10

Date Collected: 08/03/11 18:00

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 10:34	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/03/11 18:00	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 18:23	GWM	TAL NSH
Total	Prep	EPA 3510C		0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 15:12	KKH	TAL NSH

Client Sample ID: MW-53D (08032011)

Lab Sample ID: NUH0938-11

Date Collected: 08/03/11 15:10

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 11:02	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 15:10	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	1.00	U014376	08/11/11 13:32	GWM	TAL NSH
Total	Prep	EPA 3510C		0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 15:36	KKH	TAL NSH

Client Sample ID: MW-53I (08032011)

Lab Sample ID: NUH0938-12

Date Collected: 08/03/11 12:10

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 11:30	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014133	08/10/11 01:05	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 12:10	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	20.0	U014376	08/11/11 18:01	GWM	TAL NSH
Total	Prep	EPA 3510C		0.971	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 16:00	KKH	TAL NSH

Client Sample ID: MW-53S (08032011)

Lab Sample ID: NUH0938-13

Date Collected: 08/03/11 11:15

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 11:58	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014193	08/10/11 09:30	CMM	TAL NSH
Total	Prep	EPA 5030B	RE2	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE2	100	U014193	08/10/11 09:59	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 11:15	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	100	U014376	08/11/11 15:32	GWM	TAL NSH
Total	Prep	EPA 3510C	RE1	0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B	RE1	2.00	U014235	08/11/11 06:23	KKH	TAL NSH

Client Sample ID: MW-54D (08032011)

Lab Sample ID: NUH0938-14

Date Collected: 08/03/11 10:15

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 12:26	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014133	08/10/11 01:33	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 10:15	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	25.0	U014376	08/11/11 17:31	GWM	TAL NSH
Total	Prep	EPA 3510C	RE1	1.05	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B	RE1	2.00	U014235	08/11/11 06:47	KKH	TAL NSH

Client Sample ID: MW-55D (08022011)

Lab Sample ID: NUH0938-15

Date Collected: 08/02/11 12:20

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH

Client Sample ID: MW-55D (08022011)

Lab Sample ID: NUH0938-15

Date Collected: 08/02/11 12:20

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Analysis	SW846 8260B	RE1	1.00	U014133	08/09/11 22:45	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/02/11 12:20	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	1.00	U014376	08/11/11 15:02	GWM	TAL NSH
Total	Prep	EPA 3510C		1.00	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 17:12	KKH	TAL NSH

Client Sample ID: MW-55S (08022011)

Lab Sample ID: NUH0938-16

Date Collected: 08/02/11 13:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 13:23	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014133	08/10/11 02:01	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/02/11 13:05	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 21:23	GWM	TAL NSH
Total	Prep	EPA 3510C		0.971	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 17:36	KKH	TAL NSH

Client Sample ID: RW-52 (08022011)

Lab Sample ID: NUH0938-17

Date Collected: 08/02/11 18:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 13:51	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/02/11 18:05	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 21:52	GWM	TAL NSH
Total	Prep	EPA 3510C		0.952	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 18:00	KKH	TAL NSH

Client Sample ID: VW-1 (08032011)

Lab Sample ID: NUH0938-18

Date Collected: 08/03/11 15:45

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 14:19	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	20.0	U014133	08/10/11 02:57	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 15:45	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	20.0	U014376	08/11/11 18:31	GWM	TAL NSH
Total	Prep	EPA 3510C	RE1	0.952	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B	RE1	2.00	U014235	08/11/11 07:11	KKH	TAL NSH

Client Sample ID: VW-2 (08032011)

Lab Sample ID: NUH0938-19

Date Collected: 08/03/11 16:05

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	1.00	U014133	08/09/11 23:13	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/03/11 16:05	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/10/11 22:52	GWM	TAL NSH
Total	Prep	EPA 3510C		0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 18:48	KKH	TAL NSH

Client Sample ID: VW-3 (08032011)

Lab Sample ID: NUH0938-20

Date Collected: 08/03/11 10:40

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 15:15	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 10:40	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	1.00	U014376	08/11/11 14:02	GWM	TAL NSH
Total	Prep	EPA 3510C		0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 20:00	KKH	TAL NSH

Client Sample ID: MP-1 (08032011)

Lab Sample ID: NUH0938-21

Date Collected: 08/03/11 16:50

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 15:43	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014193	08/10/11 10:27	CMM	TAL NSH
Total	Prep	EPA 5030B	RE2	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE2	100	U014193	08/10/11 10:55	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 16:50	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	100	U014376	08/11/11 16:02	GWM	TAL NSH
Total	Prep	EPA 3510C	RE1	0.990	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B	RE1	2.00	U014235	08/11/11 07:35	KKH	TAL NSH

Client Sample ID: DPE-1 (08032011)

Lab Sample ID: NUH0938-22

Date Collected: 08/03/11 17:45

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 16:11	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014193	08/10/11 11:23	CMM	TAL NSH
Total	Prep	EPA 5030B	RE2	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH

Client Sample ID: DPE-1 (08032011)

Lab Sample ID: NUH0938-22

Date Collected: 08/03/11 17:45

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Analysis	SW846 8260B	RE2	100	U014193	08/10/11 11:51	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 17:45	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	100	U014376	08/11/11 16:32	GWM	TAL NSH
Total	Prep	EPA 3510C	RE1	0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B	RE1	2.00	U014235	08/11/11 07:59	KKH	TAL NSH

Client Sample ID: DUP-1 (08022011)

Lab Sample ID: NUH0938-23

Date Collected: 08/02/11 12:00

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 16:39	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014133	08/10/11 02:29	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/02/11 12:00	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	25.0	U014376	08/11/11 20:01	GWM	TAL NSH
Total	Prep	EPA 3510C		1.00	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 21:12	KKH	TAL NSH

Client Sample ID: DUP-2 (08032011)

Lab Sample ID: NUH0938-24

Date Collected: 08/03/11 12:00

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B	RE1	1.00	11H2288_P	08/09/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	1.00	U014133	08/09/11 23:41	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/03/11 12:00	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	1.00	U014376	08/11/11 14:32	GWM	TAL NSH
Total	Prep	EPA 3510C		0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 21:36	KKH	TAL NSH

Client Sample ID: RB-1 (08022011)

Lab Sample ID: NUH0938-25

Date Collected: 08/02/11 15:20

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 01:42	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/02/11 15:20	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/11/11 02:21	GWM	TAL NSH
Total	Prep	EPA 3510C		0.980	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 22:00	KKH	TAL NSH

Client Sample ID: RB-2 (08032011)

Lab Sample ID: NUH0938-26

Date Collected: 08/03/11 17:55

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 02:10	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)		1.00	11H2053_P	08/03/11 17:55	GWM	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014275	08/11/11 02:51	GWM	TAL NSH
Total	Prep	EPA 3510C		0.943	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B		1.00	U014235	08/10/11 22:24	KKH	TAL NSH

Client Sample ID: TB-1 (08032011)

Lab Sample ID: NUH0938-27

Date Collected: 08/03/11 00:01

Matrix: Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 00:46	CMM	TAL NSH

Client Sample ID: TB-2 (08032011)

Lab Sample ID: NUH0938-28

Date Collected: 08/03/11 00:01

Matrix: Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1881_P	08/08/11 17:45	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014048	08/09/11 01:14	CMM	TAL NSH

Client Sample ID: MW-55I (08022011)

Lab Sample ID: NUH0938-29

Date Collected: 08/02/11 15:40

Matrix: Ground Water

Date Received: 08/06/11 08:25

Prep Type	Batch Type	Batch Method	Run	Dilution Factor	Batch Number	Prepared Or Analyzed	Analyst	Lab
Total	Prep	EPA 5030B		1.00	11H1622_P	08/09/11 05:54	TSP	TAL NSH
Total	Analysis	SW846 8260B		1.00	U014132	08/09/11 17:36	CMM	TAL NSH
Total	Prep	EPA 5030B	RE1	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE1	10.0	U014193	08/10/11 12:19	CMM	TAL NSH
Total	Prep	EPA 5030B	RE2	1.00	11H2289_P	08/10/11 06:14	TSP	TAL NSH
Total	Analysis	SW846 8260B	RE2	100	U014193	08/10/11 12:47	CMM	TAL NSH
Total	Prep	EPA 5030B (GC)	RE1	1.00	11H2720_P	08/02/11 15:40	GWM	TAL NSH
Total	Analysis	SW846 8015B	RE1	100	U014376	08/11/11 17:01	GWM	TAL NSH
Total	Prep	EPA 3510C	RE1	0.990	11H1829_P	08/08/11 11:05	JJR	TAL NSH
Total	Analysis	SW846 8015B	RE1	2.00	U014235	08/11/11 08:23	KKH	TAL NSH

Laboratory References:
TAL NSH = TestAmerica Nashville, 2960 Foster Creighton Road, Nashville, TN 37204, TEL 800-765-0980

Method Summary

Client: ARCADIS Lexington - Exxon
Project/Site: 25553 - Thurmont, MD

TestAmerica Job ID: NUH0938

Method	Method Description	Protocol	Laboratory
SW846 8260B	Volatile Organic Compounds by EPA Method 8260B		TAL NSH
SW846 8015B	Purgeable Petroleum Hydrocarbons		TAL NSH
SW846 8015B	Extractable Petroleum Hydrocarbons		TAL NSH

Protocol References:

Laboratory References:

TAL NSH = TestAmerica Nashville, 2960 Foster Creighton Road, Nashville, TN 37204, TEL 800-765-0980

Certification Summary

Client: ARCADIS Lexington - Exxon
Project/Site: 25553 - Thurmont, MD

TestAmerica Job ID: NUH0938

Laboratory	Authority	Program	EPA Region	Certification ID
TestAmerica Nashville	A2LA	ISO/IEC 17025		0453.07
TestAmerica Nashville	A2LA	WY UST		453.07
TestAmerica Nashville	AIHA	IHLAP		100790
TestAmerica Nashville	Alabama	State Program	4	41150
TestAmerica Nashville	Alaska	Alaska UST	10	UST-087
TestAmerica Nashville	Arizona	State Program	9	AZ0473
TestAmerica Nashville	Arkansas	State Program	6	88-0737
TestAmerica Nashville	CALA	CALA		3744
TestAmerica Nashville	California	NELAC	9	1168CA
TestAmerica Nashville	Colorado	State Program	8	N/A
TestAmerica Nashville	Connecticut	State Program	1	PH-0220
TestAmerica Nashville	Florida	NELAC	4	E87358
TestAmerica Nashville	Illinois	NELAC	5	200010
TestAmerica Nashville	Iowa	State Program	7	131
TestAmerica Nashville	Kansas	NELAC	7	E-10229
TestAmerica Nashville	Kentucky	Kentucky UST	4	19
TestAmerica Nashville	Kentucky	State Program	4	90038
TestAmerica Nashville	Louisiana	NELAC	6	LA100011
TestAmerica Nashville	Louisiana	NELAC	6	30613
TestAmerica Nashville	Maryland	State Program	3	316
TestAmerica Nashville	Massachusetts	State Program	1	M-TN032
TestAmerica Nashville	Minnesota	NELAC	5	047-999-345
TestAmerica Nashville	Mississippi	State Program	4	N/A
TestAmerica Nashville	Montana	MT DEQ UST	8	NA
TestAmerica Nashville	Nevada	State Program	9	TN00032
TestAmerica Nashville	New Hampshire	NELAC	1	2963
TestAmerica Nashville	New Jersey	NELAC	2	TN965
TestAmerica Nashville	New York	NELAC	2	11342
TestAmerica Nashville	North Carolina	North Carolina DENR	4	387
TestAmerica Nashville	North Dakota	State Program	8	R-146
TestAmerica Nashville	Ohio	OVAP	5	CL0033
TestAmerica Nashville	Oklahoma	State Program	6	9412
TestAmerica Nashville	Oregon	NELAC	10	TN200001
TestAmerica Nashville	Pennsylvania	NELAC	3	68-00585
TestAmerica Nashville	Rhode Island	State Program	1	LAO00268
TestAmerica Nashville	South Carolina	State Program	4	84009
TestAmerica Nashville	South Carolina	State Program	4	84009
TestAmerica Nashville	Tennessee	State Program	4	2008
TestAmerica Nashville	Texas	NELAC	6	T104704077-09-TX
TestAmerica Nashville	USDA	USDA		S-48469
TestAmerica Nashville	Utah	NELAC	8	TAN
TestAmerica Nashville	Virginia	NELAC Secondary AB	3	460152
TestAmerica Nashville	Virginia	State Program	3	00323
TestAmerica Nashville	Washington	State Program	10	C789
TestAmerica Nashville	West Virginia	West Virginia DEP	3	219
TestAmerica Nashville	Wisconsin	State Program	5	998020430

Accreditation may not be offered or required for all methods and analytes reported in this package. Please contact your project manager for the laboratory's current list of certified methods and analytes.

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
Nashville, TN

COOLER R

NUH0938

Cooler Received/Opened On 8/6/2011 @ 8:25

1. Tracking # 7128 (last 4 digits, FedE

Courier: FEDEX IR Gun ID 12080142

2. Temperature of rep. sample or temp blank when opened: 2.1 *Degrees Celsius*

3. If Item #2 temperature is 0°C or less, was the representative sample or temp blank frozen? YES...NO...NA

4. Were custody seals on outside of cooler? YES...NO...NA

   If yes, how many and where: 1 fr.

5. Were the seals intact, signed, and dated correctly? YES...NO...NA

6. Were custody papers inside cooler? YES...NO...NA

I certify that I opened the cooler and answered questions 1-6 (intial) JG

7. Were custody seals on containers: YES NO and Intact YES...NO...NA

   Were these signed and dated correctly? YES...NO...NA

8. Packing mat'l used? Bubblewrap Plastic bag Peanuts Vermiculite Foam Insert Paper Other None

9. Cooling process: Ice Ice-pack Ice (direct contact) Dry ice Other None

10. Did all containers arrive in good condition (unbroken)? YES...NO...NA

11. Were all container labels complete (#, date, signed, pres., etc)? YES...NO...NA

12. Did all container labels and tags agree with custody papers? YES...NO...NA

13a. Were VOA vials received? YES...NO...NA

   b. Was there any observable headspace present in any VOA vial? YES...NO...NA

14. Was there a Trip Blank in this cooler? YES...NO...NA If multiple coolers, sequence # 2

I certify that I unloaded the cooler and answered questions 7-14 (intial) GB

15a. On pres'd bottles, did pH test strips suggest preservation reached the correct pH level? YES...NO...NA

   b. Did the bottle labels indicate that the correct preservatives were used YES...NO...NA

16. Was residual chlorine present? YES...NO...NA

I certify that I checked for chlorine and pH as per SOP and answered questions 15-16 (intial) G

17. Were custody papers properly filled out (ink, signed, etc)? YES...NO...NA

18. Did you sign the custody papers in the appropriate place? YES...NO...NA

19. Were correct containers used for the analysis requested? YES...NO...NA

20. Was sufficient amount of sample sent in each container? YES...NO...NA

I certify that I entered this project into LIMS and answered questions 17-20 (intial) G

I certify that I attached a label with the unique LIMS number to each container (intial) G

21. Were there Non-Conformance issues at login? YES...NO Was a PIPE generated? YES...NO...#

DUP 2, One Broken Vial
VW-3 581 VP-3 on 3U bottles but I matched the Times up

BIS = Broken in shipment
Cooler Receipt Form.doc

LF-1
End of Form

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Nashville, TN

COOLER REC

NUH0936

Cooler Received/Opened On 8/6/2011 @ 0825

1. Tracking # 6977 (last 4 digits, FedEx)

Courier: FedEx IR Gun ID 14740456

2. Temperature of rep. sample or temp blank when opened: 0.7 *Degrees Celsius*
3. If Item #2 temperature is 0°C or less, was the representative sample or temp blank frozen? YES NO...NA
4. Were custody seals on outside of cooler? YES...NO...NA

   If yes, how many and where: 1 Front
5. Were the seals intact, signed, and dated correctly? YES...NO...NA
6. Were custody papers inside cooler? YES...NO...NA

I certify that I opened the cooler and answered questions 1-6 (intial) F

7. Were custody seals on containers: YES NO and Intact YES...NO...NA

   Were these signed and dated correctly? YES...NO...NA
8. Packing mat'l used? Bubblewrap Plastic bag Peanuts Vermiculite Foam Insert Paper Other None
9. Cooling process: Ice Ice-pack Ice (direct contact) Dry ice Other None
10. Did all containers arrive in good condition (unbroken)? YES...NO...NA
11. Were all container labels complete (#, date, signed, pres., etc)? YES...NO...NA
12. Did all container labels and tags agree with custody papers? YES...NO...NA

13a. Were VOA vials received? YES...NO...NA

   b. Was there any observable headspace present in any VOA vial? YES...NO...NA
14. Was there a Trip Blank in this cooler? YES...NO...NA If multiple coolers, sequence # 1

I certify that I unloaded the cooler and answered questions 7-14 (intial) EB

15a. On pres'd bottles, did pH test strips suggest preservation reached the correct pH level? YES...NO...NA

   b. Did the bottle labels indicate that the correct preservatives were used YES...NO...NA
16. Was residual chlorine present? YES...NO...NA

I certify that I checked for chlorine and pH as per SOP and answered questions 15-16 (intial) F

17. Were custody papers properly filled out (ink, signed, etc)? YES...NO...NA
18. Did you sign the custody papers in the appropriate place? YES...NO...NA
19. Were correct containers used for the analysis requested? YES...NO...NA
20. Was sufficient amount of sample sent in each container? YES...NO...NA

I certify that I entered this project into LIMS and answered questions 17-20 (intial) F

I certify that I attached a label with the unique LIMS number to each container (intial) F

21. Were there Non-Conformance issues at login? YES...NO Was a PIPE generated? YES...NO...# _____

One Amber Broken, MW-551

BIS = Broken in shipment
Cooler Receipt Form.doc

LF-1
End of Form

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08/16/2011

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Nashville, TN

COOLER RE(

NUH0938

Cooler Received/Opened On 8/6/2011 @ 0825

1. Tracking # 6825 (last 4 digits, FedEx)

Courier: FedEx IR Gun ID 94660220

2. Temperature of rep. sample or temp blank when opened: 2.2 *Degrees Celsius*

3. If Item #2 temperature is 0°C or less, was the representative sample or temp blank frozen? YES NO NA

4. Were custody seals on outside of cooler? YES...NO...NA

   If yes, how many and where: (1) Front

5. Were the seals intact, signed, and dated correctly? YES...NO...NA

6. Were custody papers inside cooler? YES...NO...NA

I certify that I opened the cooler and answered questions 1-6 (intial) _____

7. Were custody seals on containers: YES NO and Intact YES...NO...NA

   Were these signed and dated correctly? YES...NO...NA

8. Packing mat'l used? Bubblewrap Plastic bag Peanuts Vermiculite Foam Insert Paper Other None

9. Cooling process: Ice Ice-pack Ice (direct contact) Dry ice Other None

10. Did all containers arrive in good condition (unbroken)? YES...NO...NA

11. Were all container labels complete (#, date, signed, pres., etc)? YES...NO...NA

12. Did all container labels and tags agree with custody papers? YES...NO...NA

13a. Were VOA vials received? YES...NO...NA

   b. Was there any observable headspace present in any VOA vial? YES...NO...NA

14. Was there a Trip Blank in this cooler? YES...NO...NA If multiple coolers, sequence # 7

I certify that I unloaded the cooler and answered questions 7-14 (intial) _____

15a. On pres'd bottles, did pH test strips suggest preservation reached the correct pH level? YES..NO..NA

   b. Did the bottle labels indicate that the correct preservatives were used YES...NO...NA

16. Was residual chlorine present? YES...NO...NA

I certify that I checked for chlorine and pH as per SOP and answered questions 15-16 (intial) _____

17. Were custody papers properly filled out (ink, signed, etc)? YES...NO...NA

18. Did you sign the custody papers in the appropriate place? YES...NO...NA

19. Were correct containers used for the analysis requested? YES...NO...NA

20. Was sufficient amount of sample sent in each container? YES...NO...NA

I certify that I entered this project into LIMS and answered questions 17-20 (intial) _____

I certify that I attached a label with the unique LIMS number to each container (intial) _____

21. Were there Non-Conformance issues at login? YES...NO Was a PIPE generated? YES...NO...# _____

BIS = Broken in shipment
Cooler Receipt Form.doc

LF-1
End of Form

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Nashville, TN

COOLER RE

NUH0938

Cooler Received/Opened On 8/6/2011 @ 08:25

1. Tracking # 6685 (last 4 digits, FedEx)

Courier: FEDEX IR Gun ID 96210146

2. Temperature of rep. sample or temp blank when opened: 1.4 *Degrees Celsius*

3. If Item #2 temperature is 0°C or less, was the representative sample or temp blank frozen? YES...NO...NA

4. Were custody seals on outside of cooler? YES...NO...NA

   If yes, how many and where: N/A

5. Were the seals intact, signed, and dated correctly? YES...NO...NA

6. Were custody papers inside cooler? YES...NO...NA

I certify that I opened the cooler and answered questions 1-6 (intial) P.H.

7. Were custody seals on containers: YES NO and Intact YES...NO...NA

   Were these signed and dated correctly? YES...NO...NA

8. Packing mat'l used? Bubblewrap Plastic bag Peanuts Vermiculite Foam Insert Paper Other None

9. Cooling process: Ice Ice-pack Ice (direct contact) Dry ice Other None

10. Did all containers arrive in good condition (unbroken)? YES...NO...NA

11. Were all container labels complete (#, date, signed, pres., etc)? YES...NO...NA

12. Did all container labels and tags agree with custody papers? YES...NO...NA

13a. Were VOA vials received? YES...NO...NA

   b. Was there any observable headspace present in any VOA vial? YES...NO...NA

14. Was there a Trip Blank in this cooler? YES...NO...NA If multiple coolers, sequence # 5

I certify that I unloaded the cooler and answered questions 7-14 (intial) EB

15a. On pres'd bottles, did pH test strips suggest preservation reached the correct pH level? YES...NO...NA

   b. Did the bottle labels indicate that the correct preservatives were used YES...NO...NA

16. Was residual chlorine present? YES...NO...NA

I certify that I checked for chlorine and pH as per SOP and answered questions 15-16 (intial) [illegible]

17. Were custody papers properly filled out (ink, signed, etc)? YES...NO...NA

18. Did you sign the custody papers in the appropriate place? YES...NO...NA

19. Were correct containers used for the analysis requested? YES...NO...NA

20. Was sufficient amount of sample sent in each container? YES...NO...NA

I certify that I entered this project into LIMS and answered questions 17-20 (intial) [illegible]

I certify that I attached a label with the unique LIMS number to each container (intial) [illegible]

21. Were there Non-Conformance issues at login? YES...NO Was a PIPE generated? YES...NO...#

BIS = Broken in shipment
Cooler Receipt Form.doc

LF-1
End of Form

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Nashville, TN

COOLER RI

NUH0938

Cooler Received/Opened On 8/6/2011 @ 08:25

1. Tracking # 6917 (last 4 digits, FedEx

Courier: FEDEX IR Gun ID 96210146

2. Temperature of rep. sample or temp blank when opened: 0.3 *Degrees Celsius*

3. If Item #2 temperature is 0°C or less, was the representative sample or temp blank frozen? YES...NO...NA

4. Were custody seals on outside of cooler? YES...NO...NA

   If yes, how many and where: 1- Front

5. Were the seals intact, signed, and dated correctly? YES...NO...NA

6. Were custody papers inside cooler? YES...NO...NA

I certify that I opened the cooler and answered questions 1-6 (intial) P.H.

7. Were custody seals on containers: YES NO and Intact YES...NO...NA

   Were these signed and dated correctly? YES...NO...NA

8. Packing mat'l used? Bubblewrap Plastic bag Peanuts Vermiculite Foam Insert Paper Other None

9. Cooling process: Ice Ice-pack Ice (direct contact) Dry ice Other None

10. Did all containers arrive in good condition (unbroken)? YES...NO...NA

11. Were all container labels complete (#, date, signed, pres., etc)? YES...NO...NA

12. Did all container labels and tags agree with custody papers? YES...NO...NA

13a. Were VOA vials received? YES...NO...NA

   b. Was there any observable headspace present in any VOA vial? YES...NO...NA

14. Was there a Trip Blank in this cooler? YES...NO...NA If multiple coolers, sequence # 4

I certify that I unloaded the cooler and answered questions 7-14 (intial) GB

15a. On pres'd bottles, did pH test strips suggest preservation reached the correct pH level? YES...NO...NA

   b. Did the bottle labels indicate that the correct preservatives were used YES...NO...NA

16. Was residual chlorine present? YES...NO...NA

I certify that I checked for chlorine and pH as per SOP and answered questions 15-16 (intial) F

17. Were custody papers properly filled out (ink, signed, etc)? YES...NO...NA

18. Did you sign the custody papers in the appropriate place? YES...NO...NA

19. Were correct containers used for the analysis requested? YES...NO...NA

20. Was sufficient amount of sample sent in each container? YES...NO...NA

I certify that I entered this project into LIMS and answered questions 17-20 (intial) F

I certify that I attached a label with the unique LIMS number to each container (intial) F

21. Were there Non-Conformance issues at login? YES...NO Was a PIPE generated? YES...NO #_____

BIS = Broken in shipment
Cooler Receipt Form.doc

LF 1
End of Form

Revised 6/24/09

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING
Nashville, TN

COOLER RE

NUH0936

Cooler Received/Opened On 8/6/2011 @ 08:25

1. Tracking # 7220 (last 4 digits, FedEx)

Courier: FEDEX IR Gun ID 96210146

2. Temperature of rep. sample or temp blank when opened: 0.9 *Degrees Celsius*

3. If Item #2 temperature is 0°C or less, was the representative sample or temp blank frozen? YES...NO...NA

4. Were custody seals on outside of cooler? YES...NO...NA

   If yes, how many and where: 1 front

5. Were the seals intact, signed, and dated correctly? YES...NO...NA

6. Were custody papers inside cooler? YES...NO...NA

I certify that I opened the cooler and answered questions 1-6 (intial) GB

7. Were custody seals on containers: YES NO and Intact YES...NO...NA

   Were these signed and dated correctly? YES...NO...NA

8. Packing mat'l used? Bubblewrap Plastic bag Peanuts Vermiculite Foam Insert Paper Other None

9. Cooling process: Ice Ice-pack Ice (direct contact) Dry ice Other None

10. Did all containers arrive in good condition (unbroken)? YES...NO...NA

11. Were all container labels complete (#, date, signed, pres., etc)? YES...NO...NA

12. Did all container labels and tags agree with custody papers? YES...NO...NA

13a. Were VOA vials received? YES...NO...NA

   b. Was there any observable headspace present in any VOA vial? YES...NO...NA

14. Was there a Trip Blank in this cooler? YES...NO...NA If multiple coolers, sequence # 3

I certify that I unloaded the cooler and answered questions 7-14 (intial) GB

15a. On pres'd bottles, did pH test strips suggest preservation reached the correct pH level? YES...NO...NA

   b. Did the bottle labels indicate that the correct preservatives were used YES...NO...NA

16. Was residual chlorine present? YES...NO...NA

I certify that I checked for chlorine and pH as per SOP and answered questions 15-16 (intial) G

17. Were custody papers properly filled out (ink, signed, etc)? YES...NO...NA

18. Did you sign the custody papers in the appropriate place? YES...NO...NA

19. Were correct containers used for the analysis requested? YES...NO...NA

20. Was sufficient amount of sample sent in each container? YES...NO...NA

I certify that I entered this project into LIMS and answered questions 17-20 (intial) G

I certify that I attached a label with the unique LIMS number to each container (intial) G

21. Were there Non-Conformance issues at login? YES...NO Was a PIPE generated? YES...NO...#

BIS = Broken in shipment
Cooler Receipt Form.doc

LF-1
End of Form

Revised 6/24/09

TestAmerica
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Nashville, TN

NUH0938
08/22/11 23:59

COOLER RECEIPT FORM

Cooler Received/Opened On 8/6/2011 @ 8:25

1. Tracking # 6825 (last 4 digits, FedEx)

Courier: FEDEX IR Gun ID 12080142

2. Temperature of rep. sample or temp blank when opened: 2.8 *Degrees Celsius*

3. If Item #2 temperature is 0°C or less, was the representative sample or temp blank frozen? YES...NO...NA

4. Were custody seals on outside of cooler? YES...NO...NA

   If yes, how many and where: 1f.

5. Were the seals intact, signed, and dated correctly? YES...NO...NA

6. Were custody papers inside cooler? YES...NO...NA

I certify that I opened the cooler and answered questions 1-6 (intial) J.G.

7. Were custody seals on containers: YES NO and Intact YES...NO...NA

   Were these signed and dated correctly? YES...NO...NA

8. Packing mat'l used? Bubblewrap Plastic bag Peanuts Vermiculite Foam Insert Paper Other None

9. Cooling process: Ice Ice-pack Ice (direct contact) Dry ice Other None

10. Did all containers arrive in good condition (unbroken)? YES...NO...NA

11. Were all container labels complete (#, date, signed, pres., etc)? YES...NO...NA

12. Did all container labels and tags agree with custody papers? YES...NO...NA

13a. Were VOA vials received? YES...NO...NA

   b. Was there any observable headspace present in any VOA vial? YES...NO...NA

14. Was there a Trip Blank in this cooler? YES...NO...NA If multiple coolers, sequence # 6

I certify that I unloaded the cooler and answered questions 7-14 (intial) GB

15a. On pres'd bottles, did pH test strips suggest preservation reached the correct pH level? YES...NO...NA

   b. Did the bottle labels indicate that the correct preservatives were used YES...NO...NA

16. Was residual chlorine present? YES...NO...NA

I certify that I checked for chlorine and pH as per SOP and answered questions 15-16 (intial) J

17. Were custody papers properly filled out (ink, signed, etc)? YES...NO...NA

18. Did you sign the custody papers in the appropriate place? YES...NO...NA

19. Were correct containers used for the analysis requested? YES...NO...NA

20. Was sufficient amount of sample sent in each container? YES...NO...NA

I certify that I entered this project into LIMS and answered questions 17-20 (intial) J

I certify that I attached a label with the unique LIMS number to each container (intial) J

21. Were there Non-Conformance issues at login? YES...NO Was a PIPE generated? YES...NO...#

BIS = Broken in shipment
Cooler Receipt Form.doc

LF-1
End of Form

Revised 6/24/09

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THE LEADER IN ENVIRONMENTAL TESTING

NUH0938

08/22/11 23:59

Nashville Division
2960 Foster Creighton Drive * Nashville TN 37204
Phone: (800) 765-0980 / (615) 726-0177 Fax:(615) 726-3404
","Reg District (CA)")

ExxonMobil

Consultant: ARCADIS Lexington - Exxon
Address: 1114 Benfield Boulevard, Suite A
City, State, Zip: Millersville MD 21108
ExxonMobil Project Mgr: Account Payable
Consultant Project Mgr: Rusty Kahl
Consultant Telephone #r: (410) 987-0032 Fax: -
Sampler Name (Print) Stephanie Amodeo
SamplerSignature: Stephanie Amodeo

TA Account #: 1409738 PO #: B0085851.0014
Invoice to: ARCADIS Lexington - Exxon
Report to: Rusty Kahl
Project Name: Exxon 25553
Retail Project (MRN): 25553
Major Project (AFE):
Site Address: 143 Frederick Road
City,State,Zip: Thurmont Maryland

Sample ID	Date Sampled	Time Sampled	# Containers Shipped	Grab	Composite	Field Filtered	Preservative: Methanol	Sodium Bisulfate	(Blue Label) HCL	(Orange Label) NaOH	(Yellow Label) Plastic H2SO4	(Yellow Label) Glass H2SO4	(Red Label) HNO3	(Black Label) None	Matrix: Groundwater	Wastewater	Drinking Water	Sludge	Soil	(specify) Other	Analyze for: 8260B VOCs + Oxygenates (5)	TPH - Diesel Range SW846 8015B	TPH-Gasoline Range SW8015	RUSH TAT (Pre Schedule)*
MP-1 (08032011)	8/3/11	1650	8	X					X						X						X	X	X	
DPE-1 (08032011)	8/3/11	1745	8	X					X						X						X	X	X	
DUP-1 (08022011)	8/2/11	1200	8	X					X						X						X	X	X	
DUP-2 (08032011)	8/3/11	1200	8	X					X						X						X	X	X	
RB-1 (08022011)	8/2/11	1520	8	X					X						X						X	X	X	
RB-2 (08032011)	8/3/11	1755	8	X					X						X						X	X	X	
TB-1 (08032011)	8/3/11	—	2	X					X						X						X		X	
TB-2 (08032011)	8/3/11	—	2	X					X						X						X		X	
MW-55I (08022011)	8/2/11	1540	8	X					X						X						X	X	X	

COMMENTS: All turn around times are calculated from the time of receipt at TestAmerica.
* It will be the responsibility of Exxon Mobil or its consultant to notify the TestAmerica Project Manager by phone or fax that a rush sample will be submitted. TA Project manager_____ Date:_____
There may be a charge assessed for TestAmerica disposing of sample remainders.

NOTES/SPECIAL INSTRUCTIONS: *BO # 22985*

Relinquished by:	Date:	Time:	Received by:	Date:	Time:	Relinquished by:	Date:	Time:
Stephanie Amodeo	8/4/11	0700						

Shipped Via: Lab Courier

Received for TestAmerica by: Date: 8-6-11 Time: 8:25

Shipped Via:

Temperature Upon Receipt: ambient

Sample Containers Intact? Y N
VOCs Free of Headspace? Y N

QC Deliverables (Please Circle One) :
Level 2 Level 3 Level 4 Site Specific
(If site specific, please pre-schedule w/ TestAmerica Project Manager or attach specific instructions)

Date Due of Report:

08/16/2011

Te

NUH0938
08/22/11 23 59

THE LEADER IN ENVIRONMENTAL TESTING

Nashville Division
2960 Foster Creighton Drive * Nashville TN 37204
Phone: (800) 765-0980 / (615) 726-0177 Fax:(615) 726-3404
',"Reg District (CA)")

ExxonMobil

08/16/2011

Consultant: ARCADIS Lexington - Exxon
Address: 1114 Benfield Boulevard, Suite A
City, State, Zip: Millersville MD 21108
ExxonMobil Project Mgr: Account Payable
Consultant Project Mgr: Rusty Kahl
Consultant Telephone #r: (410) 987-0032 Fax: -
Sampler Name (Print) Stephanie Amodeo
SamplerSignature: Stephanie Amodeo

TA Account #: 1409738 PO #: B0085851.0014
Invoice to: ARCADIS Lexington - Exxon
Report to: Rusty Kahl
Project Name: Exxon 25553
Retail Project (MRN): 25553
Major Project (AFE):
Site Address: 143 Frederick Road
City,State,Zip: Thurmont Maryland

Sample ID	Date Sampled	Time Sampled	# Containers Shipped	Grab	Composite	Field Filtered	Preservative: Methanol	Sodium Bisulfate	(Blue Label) HCL	(Orange Label) NaOH	(Yellow Label) Plastic H2SO4	(Yellow Label) Glass H2SO4	(Red Label) HNO3	(Black Label) None	Matrix: Groundwater	Wastewater	Drinking Water	Sludge	Soil	(specify) Other	Analyze for: 8260B VOCs + Oxygenates (5)	TPH - Diesel Range SW846 8015B	TPH-Gasoline Range SW8015	RUSH TAT (Pre Schedule)*
MW-53D(08032011)	8/3/11	1510	8	X					X						X						X	X	X	
MW-53I(08032011)	8/3/11	1210	8	X					X						X						X	X	X	
MW-53S(08032011)	8/3/11	1115	8	X					X						X						X	X	X	
MW-54D(08032011)	8/3/11	1015	8	X					X						X						X	X	X	
MW-55D(08022011)	8/2/11	1220	8	X					X						X						X	X	X	
MW-55S(08022011)	8/2/11	1305	8	X					X						X						X	X	X	
RW-52(08022011)	8/2/11	1805	8	X					X						X						X	X	X	
VW-1(08032011)	8/3/11	1545	8	X					X						X						X	X	X	
VW-2(08032011)	8/3/11	1605	7	X					X						X						X	X	X	
VW-3(08032011)	8/3/11	1040	8	X					X						X						X	X	X	

COMMENTS: All turn around times are calculated from the time of receipt at TestAmerica.
* It will be the responsibility of Exxon Mobil or its consultant to notify the TestAmerica Project Manager by phone or fax that a rush sample will be submitted. TA Project manager _____ Date: _____
There may be a charge assessed for TestAmerica disposing of sample remainders.

NOTES/SPECIAL INSTRUCTIONS: *BO # 22985*

Relinquished by:	Date:	Time:	Received by:	Date:	Time:	Relinquished by:	Date:	Time:
Stephanie Amodeo	08/4/11	0700						

Shipped Via: Lab Courier
Shipped Via:

Received for TestAmerica by: Date: 8-4-11 Time: 8:25
Temperature Upon Receipt: Multiple
Sample Containers Intact? Y N
VOCs Free of Headspace? Y N

QC Deliverables (Please Circle One) :
Level 2 Level 3 Level 4 Site Specific
(If site specific, please pre-schedule w/ TestAmerica Project Manager or attach specific instructions)

Date Due of Report:

TestAmerica
THE LEADER IN ENVIRONMENTAL TESTING

NUH0938
08/22/11 23 59

Nashville Division
2960 Foster Creighton Drive * Nashville TN 37204
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ExxonMobil

Consultant: ARCADIS Lexington - Exxon
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City, State, Zip: Millersville MD 21108
ExxonMobil Project Mgr: Account Payable
Consultant Project Mgr: Rusty Kahl
Consultant Telephone #r: (410) 987-0032 Fax: -
Sampler Name (Print) Stephanie Amadeo
SamplerSignature: Stephanie Amadeo

TA Account #: 1409738 PO #: B0085851.0014
Invoice to: ARCADIS Lexington - Exxon
Report to: Rusty Kahl
Project Name: Exxon 25553
Retail Project (MRN): 25553
Major Project (AFE):
Site Address: 143 Frederick Road
City,State,Zip: Thurmont Maryland

Sample ID	Date Sampled	Time Sampled	# Containers Shipped	Grab	Composite	Field Filtered	Preservative: Methanol	Sodium Bisulfate	(Blue Label) HCL	(Orange Label) NaOH	(Yellow Label) Plastic H2SO4	(Yellow Label) Glass H2SO4	(Red Label) HNO3	(Black Label) None	Matrix: Groundwater	Wastewater	Drinking Water	Sludge	Soil	(specify) Other	Analyze for: 8260B VOCs + Oxygenates (5)	TPH - Diesel Range SW846 8015B	TPH-Gasoline Range SW8015	RUSH TAT (Pre Schedule)*
MW-5 (08032011)	08/3/11	1240	8	X					X						X						X	X	X	
MW-6 (08032011)	8/3/11	1505	8	X					X						X						X	X	X	
MW-13A (08022011)	8/2/11	2015	8	X					X						X						X	X	X	
MW-14A (08022011)	8/2/11	2005	8	X					X						X						X	X	X	
MW-37 (08022011)	8/2/11	1520	8	X					X						X						X	X	X	
MW-38 (08022011)	8/2/11	1635	8	X					X						X						X	X	X	
MW-46 (08022011)	8/2/11	1245	8	X					X						X						X	X	X	
MW-48 (08022011)	8/2/11	1800	8	X					X						X						X	X	X	
MW-50 (08032011)	8/3/11	0930	8	X					X						X						X	X	X	
MW-51 (08032011)	8/3/11	1800	8	X					X						X						X	X	X	

COMMENTS: All turn around times are calculated from the time of receipt at TestAmerica.

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There may be a charge assessed for TestAmerica disposing of sample remainders.

NOTES/SPECIAL INSTRUCTIONS: *BO # 22985*

Relinquished by:	Date:	Time:	Received by:	Date:	Time:	Relinquished by:	Date:	Time:
Stephanie Amadeo	8/4/11	0700	Paul Gaylord	8/5/11	1011	Paul Gaylord	8/5/11	1710
Shipped Via: Lab Courier			Shipped Via:					
Received for TestAmerica by:	Date:	Time:	Temperature Upon Receipt: Multiple	Sample Containers Intact? Y N		QC Deliverables (Please Circle One): Level 2 Level 3 Level 4 Site Specific (If site specific, please pre-schedule w/ TestAmerica Project Manager or attach specific instructions)	Date Due of Report:	
	8-6-11	8:25		VOCs Free of Headspace? Y N				

08/16/2011