

May 21, 2012

Washington State Department of Transportation  
Headquarters Environmental Services Office  
P.O. Box 47417  
Olympia, Washington 98504-7417

Attention: Jeff Sawyer

Subject: Site Summary Letter  
Remedial Investigation Field Activities and Sampling Results  
US 101 Midway Metals  
Clallam County, Washington  
File No. 0180-292-00

## **INTRODUCTION**

GeoEngineers is pleased to present this letter which summarizes the Remedial Investigation (RI) field activities and sampling results of the Midway Metals property located at 258010 Highway 101, Sequim, Washington (Site). The Site is shown with respect to surrounding physical features on the Vicinity Map, Figure 1. This letter presents the sample locations, chemical analytical results and general observations made by GeoEngineers during field activities completed in December 2011 and February 2012.

## **BACKGROUND**

Contamination was identified at the Site during an Initial Investigation performed by Washington State Department of Ecology (Ecology) and Clallam County Environmental Health (Clallam Health) (Initial Investigation Field Report, 2006) in conjunction with a Site Hazard Assessment performed by Clallam Health (Site Hazard Assessment, 2008). Washington State Department of Transportation (WSDOT) conducted a Phase II Environmental Site Assessment (ESA) in March 2011 to assess soil and groundwater conditions within a portion of the Site. Sample locations from these previous investigations are presented on the Site Plan, Figure 2.

The field activities conducted by GeoEngineers characterized the soil, groundwater, surface water and sediment on the Site. The field activities were based on the RI work plan completed and submitted to Ecology for review in December 2011. Sample locations are shown on Figure 3.

## **FIELD AND ANALYTICAL SUMMARY**

GeoEngineers completed field activities in December 2011 and February 2012. Twenty (20) soil borings were completed on site using a direct-push drill rig operated by Cascade Drilling of Woodinville,



Washington. Of the 20 soil borings, three were sampled for groundwater (DP-1, DP-12 and DP-15). The direct-push rig achieved maximum depth with sampler probe refusal at all boring locations with the exception of some shallow borings that were completed to assess surface soil. In addition to the soil and groundwater samples, GeoEngineers collected three surface water samples and nine sediment samples from drainage ditches. The drainage ditches did not contain flowing surface water at the time of our Site visits. The three surface water samples were collected from standing water observed in the ditches.

A summary of analytical results detected at concentrations greater than respective regulatory action levels (in parentheses) are listed below.

#### **Soil (Model Toxics Control Act [MTCA] Method A Unrestricted Land Use)**

- Lube Oil-Range Hydrocarbons – Sample 3 Batteries, MM-SC2 and MM-SC3.
- Cadmium – DP-4, BP-5, DP-6, HA3, Sample 1 Lawn Mower, Sample 2 Tier 2 West and Sample 3 Batteries.
- Lead – Sample 3 Batteries and MM-SC3.
- Mercury – MM-B7-0-4.
- cPAHs – DP-9.

#### **Groundwater (MTCA Method A Table Value)**

- Total Arsenic – DP1-W, DP12-W and DP15-W.
- Total Chromium – DP12-W and DP15-W.
- Total Lead – DP15-W.
- Dissolved Arsenic – DP1-W.

#### **Sediment (MTCA Method A Unrestricted Land Use)**

- cPAHs – SED5.

#### **Surface Water (ARAR – Aquatic Life -Fresh/Chronic - Ch. 173-201A WAC)<sup>1</sup>**

- Total Lead – SW1, SW2 and SW9.
- Dissolved Lead – SW1, SW2 and SW9.
- Total Aroclors (PCBs) – SW1.

#### **Surface Water (ARAR – Aquatic Life - Fresh/Chronic - Clean Water Act §304)<sup>1</sup>**

- Total Lead – SW1, SW2 and SW9.
- Total Aroclors (PCBs) – SW1.

#### **Surface Water (ARAR – Aquatic Life - Fresh/Chronic - National Toxics Rule, 40 CFR 131)<sup>1</sup>**

- Total Lead – SW1, SW2 and SW9.

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<sup>1</sup> ARARs and data gaps are currently being evaluated as part of the final RI (risk assessment).



The results listed above are presented in Table 1. All results are presented in Tables 2 through 5. The complete sets of laboratory analytical data are included in Appendix A

## **OBSERVATIONS**

### **Site Soil Conditions**

Geologic conditions observed during the RI field activities consisted of an upper unit of silty sand to sandy silt to approximately 2 feet below ground surface (bgs) (top soil), underlain by a lower unit of silty sand to silt with varying amounts of sand and gravel (weathered Glacial Till). The lower unit was relatively difficult to probe and typically resulted in sample probe refusal between 10 and 20 feet bgs.

### **Site Groundwater Conditions**

There are two groundwater wells located on Site; one in the southwest portion of the Site and one in the northeast portion of the Site. The well in the northeast portion of the Site has a steel plate welded over the top of the well casing, preventing access. The other well was available to collect a groundwater sample. See Figure 2 for the well locations.

GeoEngineers measured groundwater levels in the well located in the southwest portion of the Site in December 2011 and February 2012. This well is installed to approximately 33 feet bgs. Groundwater was 5.22 feet bgs in December and 0.25 feet above ground surface in February. Soil borings completed near this well, DP-16 and DP-17 did not contain any groundwater at the time they were completed in February 2012.

Shallow groundwater was observed approximately 10 to 14 feet bgs in the northeast portion of the Site and approximately 3 to 5 feet bgs in the southern portion of the Site. Groundwater observed in the soil borings completed on Site appears to be constrained to isolated perched lenses or directly related to surface water infiltration.

## **LIMITATIONS**

We have prepared this site summary letter for the exclusive use of WSDOT and their authorized agents for the US 101 Midway Metals project in Sequim, Washington.

Within the limitations of scope, schedule and budget, our services have been executed in accordance with generally accepted environmental science practices in this area at the time this report was prepared. The conclusions and opinions presented in this report are based on our professional knowledge, judgment and experience. No warranty or other conditions, express or implied, should be understood.



We appreciate the opportunity to assist you with this project. Please call if you have any questions regarding this submittal.

Respectfully Submitted,  
GeoEngineers, Inc.



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Attachments:

- Tables 1A, 1B and 1C. Chemical Analytical Data Summary
- Table 2A, 2B, 2C and 2D. Chemical Analytical Data – Soil
- Table 3. Chemical Analytical Data – Groundwater
- Table 4. Chemical Analytical Data – Sediment
- Table 5. Chemical Analytical Data – Surface Water
- Figure 1. Vicinity Map
- Figure 2. Site Plan with Previous Sample Locations
- Figure 3. Remedial Investigation Plan
- Figure 4. Contaminant Boundary
- Appendix A. Laboratory Analytical Data

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**TABLE 1A**  
**CHEMICAL ANALYTICAL DATA SUMMARY**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Sample Identification	Date Collected	Sample Depth Interval (ft)	Analytes - Soil & Sediment											
			Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx			Metals by EPA6010B/7471A/SW7196A			PCBs by SW8082				cPAHs by SW8270	
			Gasoline-range hydrocarbons (mg/kg)	Diesel-range hydrocarbons (mg/kg)	Lube Oil-range Hydrocarbons (mg/kg)	Cadmium (mg/kg)	Lead (mg/kg)	Mercury (mg/kg)	PCB-aroclor 1242 (µg/kg)	PCB-aroclor 1254 (µg/kg)	PCB-aroclor 1260 (µg/kg)	Total Aroclors (µg/kg)	Benzo(a) pyrene (µg/kg)	Total cPAH <sup>5</sup> TTEC (ND=0.5RL) (µg/kg)
Sample 1 Lawn Mower	10/10/2006	0-2	-	120	530	4.1	172	-	-	-	-	-	-	-
Sample 2 Tier 2 West	10/10/2006	0-2	-	280	1,300	3.5	136	-	-	-	-	-	-	-
Sample 3 Batteries	10/10/2006	0-2	-	1,800	10,000	7.1	3,000	-	-	-	-	-	-	-
MM-B7-0-4	3/2011	0-4	-	-	110	-	12	2.0	-	-	-	-	-	-
MM-B8-0-4	3/2011	0-4	-	-	1,900	-	18	-	-	-	-	-	-	-
MM-SC2	3/2011	0-2	24	-	4,900	-	150	-	-	-	-	-	-	-
MM-SC3	3/2011	0-2	13	-	4,300	-	300	-	-	-	-	-	-	-
DP-3-0.0-2.0	12/13/2011	0-2	6.3 U	180 U	880	0.8	40	-	59 U	59 U	59 U	59 U	28	38.15
DP-4-0.0-2.0	12/13/2011	0-2	7.6 U	100 U	580	2.5	150	-	420	67 U	67 U	420	9.5	12.92
DP-5-0.0-2.0	12/13/2011	0-2	5.7 U	200 U	1,100	2.1	130	-	89	140	58 U	229	46	62.72
DP-6-0.0-2.0	12/13/2011	0-2	5.4 U	280	1,100	3.2	160	-	140	56 U	130	270	33	45.63
DP-7-0.0-2.0	12/14/2011	0-2	7.2 U	34	130	0.81	93	-	120	63 U	79	199	8.5 U	9.24
DP-8-0.0-2.0	12/14/2011	0-2	18	30 U	150	0.55 U	24	-	55 U	58	55 U	58	7.3 U	6.06
DP-9-0.0-2.0	12/14/2011	0-2	8.4 U	73 U	360	1.6	85	-	70 U	70 U	70 U	70 U	<b>140</b>	<b>190.4</b>
DP-10-0.0-2.0	12/14/2011	0-2	42	34 U	120	0.58 U	8.3	-	58 U	58 U	58 U	58 U	7.8 U	5.89 U
DP-11-0.0-2.0	12/14/2011	0-2	5.8 U	91 U	480	0.89	26	-	58 U	58 U	58 U	58 U	7.7 U	7.11
DP-12-0.0-2.0	2/8/2012	0-2	5.3 U	28 U	130	0.55 U	19	-	55 U	55 U	55 U	55 U	21	24.31
DP-21-0.0-2.0	2/8/2012	0-2	-	100	470	0.83	48	-	-	-	-	-	-	-
HA3-0.0-2.0	2/9/2012	0-2	-	110 U	660	2.7	150	-	-	-	-	-	-	-
SED5	12/13/2011	-	7.6 U	120 U	950	0.89	120	0.33 U	69 U	160	66 U	160	270	356.9
MTCA A ULU			100	2,000	2,000	2.0	250	2.0	NE	NE	NE	1,000	100	100

**Bold Type** Exceeds MTCA A (Soil - Unrestricted Land Use Cleanup Level)

**TABLE 1B**  
DRAFT CHEMICAL ANALYTICAL DATA SUMMARY  
WSDOT MIDWAY METALS  
SEQUIM, WASHINGTON

Sample Identification	Date Collected	Sample Depth Interval (ft)	Analytes - Groundwater			
			Metals by EPA6010B/7471A/SW7196A			
			Total Arsenic (µg/l)	Dissolved Arsenic (µg/l)	Dissolved Chromium (µg/l)	Dissolved Lead (µg/l)
DP1-W	12/13/2011	-	18	21	26	1.1 U
DP12-W	2/8/2012	-	12	3.0 U	140	12
DP15-W	2/8/2012	-	27	3.0 U	180	16
		MTCA A	5.0	5.0	50	15
<b>Bold Type</b> Exceeds MTCA A (Groundwater - Table Value)						

**TABLE 1C**  
DRAFT CHEMICAL ANALYTICAL DATA SUMMARY  
WSDOT MIDWAY METALS  
SEQUIM, WASHINGTON

Sample Identification	Date Collected	Sample Depth Interval (ft)	Analytes - Surface Water		
			Metals by EPA6010B/7471A/SW7196A		PCBs by SW8082
			Total Lead (µg/l)	Dissolved Lead (µg/l)	Total Aroclors (µg/l)
SW1	12/13/2011	-	7.4	1.4	0.074 T
SW2	12/13/2011	-	4.0	1.4	0.048 UT
SW9	2/8/2012	-	8.3	1.6	0.053 UT
Surface Water ARAR - Aquatic Life - Fresh/Chronic - Ch. 173-201A WAC			0.54	0.54	0.014
Surface Water ARAR - Aquatic Life - Fresh/Chronic - Clean Water Act §304			2.5	2.5	0.014
Surface Water ARAR - Aquatic Life - Fresh/Chronic - National Toxics Rule, 40 CFR 131			2.5	2.5	0.14
<b>Bold Type</b> Exceeds One or More ARAR Screening Level					

**TABLE 2A**  
**CHEMICAL ANALYTICAL DATA - SOIL<sup>1</sup>**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-1-0-2 12/13/11 0 - 2	DP-1-13-15 12/13/11 13 - 15	DP-3-0-2 12/13/11 0 - 2	DP-3-3-4 <sup>3</sup> 12/13/11 3 - 4	DP-4-0-2 12/13/11 0 - 2	DP-4-5-6 <sup>3</sup> 12/13/11 5 - 6	DP-5-0-2 12/13/11 0 - 2	DP-5-6-7 <sup>3</sup> 12/13/11 6 - 7	DP-6-0-2 12/13/11 0 - 2	DP-6-3.5-4.5 <sup>3</sup> 12/13/11 3.5 - 4.5	DP-6-11.5-12.5 12/13/11 11.5 - 12.5	MTCA <sup>2</sup> A ULU
<b>Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx</b>													
Gasoline-range hydrocarbons	mg/kg	6.2 U	5.9 U	6.3 U	--	7.6 U	--	5.7 U	--	5.4 U	--	5.7 U	100
Diesel-range hydrocarbons	mg/kg	29 U	30 U	180 U	29 U	100 U	32 U	200 U	30 U	280	29 U	30 U	2,000
Lube Oil-range Hydrocarbons	mg/kg	58 U	59 U	880	58 U	580	64 U	1,100	60 U	1,100	59 U	60 U	2,000
<b>Metals by EPA6010B/7471A/SW7196A</b>													
Arsenic	mg/kg	12 U	12 U	12 U	--	13 U	13 U	12 U	12 U	11 U	12 U	12 U	20
Barium	mg/kg	70	30	89	--	99	54	120	72	72	46	42	NE
Cadmium	mg/kg	0.58 U	0.59 U	0.8	--	2.5	0.64 U	2.1	0.6 U	3.2	0.59 U	0.6 U	2.0
Chromium <sup>4</sup>	mg/kg	34	29	36	--	42	46	35	50	41	34	26	NE
Lead	mg/kg	5.8 U	5.9 U	40	--	150	6.4 U	130	6 U	160	5.9 U	6 U	250
Mercury	mg/kg	0.29 U	0.3 U	0.29 U	--	0.67	0.32 U	0.32	0.3 U	0.28 U	0.29 U	0.3 U	2.0
Selenium	mg/kg	12 U	12 U	12 U	--	13 U	13 U	12 U	12 U	11 U	12 U	12 U	NE
Silver	mg/kg	0.58 U	0.59 U	0.59 U	--	0.67 U	0.64 U	0.58 U	0.6 U	0.56 U	0.59 U	0.6 U	NE
<b>PCBs by SW8082</b>													
PCB-aroclor 1016	µg/kg	58 U	59 U	59 U	--	67 U	64 U	58 U	60 U	56 U	59 U	60 U	NE
PCB-aroclor 1221	µg/kg	58 U	59 U	59 U	--	67 U	64 U	58 U	60 U	56 U	59 U	60 U	NE
PCB-aroclor 1232	µg/kg	58 U	59 U	59 U	--	67 U	64 U	58 U	60 U	56 U	59 U	60 U	NE
PCB-aroclor 1242	µg/kg	58 U	59 U	59 U	--	420	64 U	89	60 U	140	59 U	60 U	NE
PCB-aroclor 1248	µg/kg	58 U	59 U	59 U	--	67 U	64 U	58 U	60 U	56 U	59 U	60 U	NE
PCB-aroclor 1254	µg/kg	58 U	59 U	59 U	--	67 U	64 U	140	60 U	56 U	59 U	60 U	NE
PCB-aroclor 1260	µg/kg	58 U	59 U	59 U	--	67 U	64 U	58 U	60 U	130	59 U	60 U	NE
Total Aroclors	µg/kg	58 U	59 U	59 U	--	420	64 U	229	60 U	270	59 U	60 U	1,000
<b>PAHs by SW8270</b>													
1-Methylnaphthalene	µg/kg	7.8 U	7.9 U	10	--	8.9 U	--	15	--	16	--	8 U	NE
2-Methylnaphthalene	µg/kg	7.8 U	7.9 U	20	--	18	--	32	--	45	--	8 U	NE
Acenaphthene	µg/kg	7.8 U	7.9 U	7.8 U	--	8.9 U	--	7.7 U	--	8.5	--	8 U	NE
Acenaphthylene	µg/kg	7.8 U	7.9 U	11	--	8.9 U	--	7.7 U	--	9.9	--	8 U	NE
Anthracene	µg/kg	7.8 U	7.9 U	11	--	8.9 U	--	14	--	15	--	8 U	NE
Benzo(a)anthracene	µg/kg	7.8 U	7.9 U	24	--	8.9 U	--	36	--	26	--	8 U	NE
Benzo(a)pyrene	µg/kg	7.8 U	7.9 U	28	--	9.5	--	46	--	33	--	8 U	100
Benzo(b)fluoranthene	µg/kg	7.8 U	7.9 U	39	--	15	--	62	--	47	--	8 U	NE
Benzo(ghi)perylene	µg/kg	7.8 U	7.9 U	25	--	12	--	45	--	38	--	8 U	NE
Benzo(j,k)fluoranthene	µg/kg	7.8 U	7.9 U	11	--	8.9 U	--	17	--	16	--	8 U	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-1-0-2 12/13/11 0 - 2	DP-1-13-15 12/13/11 13 - 15	DP-3-0-2 12/13/11 0 - 2	DP-3-3-4 <sup>3</sup> 12/13/11 3 - 4	DP-4-0-2 12/13/11 0 - 2	DP-4-5-6 <sup>3</sup> 12/13/11 5 - 6	DP-5-0-2 12/13/11 0 - 2	DP-5-6-7 <sup>3</sup> 12/13/11 6 - 7	DP-6-0-2 12/13/11 0 - 2	DP-6-3.5-4.5 <sup>3</sup> 12/13/11 3.5 - 4.5	DP-6-11.5-12.5 12/13/11 11.5 - 12.5	MTCA <sup>2</sup> A ULU
Chrysene	µg/kg	7.8 U	7.9 U	36	--	14	--	50	--	35	--	8 U	NE
Dibenzo(a,h)anthracene	µg/kg	7.8 U	7.9 U	7.8 U	--	8.9 U	--	9.2	--	7.5 U	--	8 U	NE
Dibenzofuran	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
Fluoranthene	µg/kg	7.8 U	7.9 U	45	--	17	--	63	--	45	--	8 U	NE
Fluorene	µg/kg	7.8 U	7.9 U	7.8 U	--	8.9 U	--	8.2	--	13	--	8 U	NE
Indeno(1,2,3-cd)pyrene	µg/kg	7.8 U	7.9 U	20	--	8.9 U	--	38	--	30	--	8 U	NE
Naphthalene	µg/kg	7.8 U	7.9 U	21	--	15	--	33	--	38	--	8 U	5,000
Phenanthrene	µg/kg	7.8 U	7.9 U	34	--	17	--	49	--	40	--	8 U	NE
Pyrene	µg/kg	13	7.9 U	47	--	20	--	68	--	63	--	8 U	NE
Total cPAH <sup>5</sup> TTEC (ND=0.5RL)	µg/kg	5.89 U	5.96 U	38.15	--	12.92	--	62.72	--	45.63	--	6.04 U	100
<b>VOCs by SW8260</b>													
1,1,1,2-Tetrachloroethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,1,1-Trichloroethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	2,000
1,1,2,2-Tetrachloroethane	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,1,2-Trichloroethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,1-Dichloroethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,1-Dichloroethene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,1-Dichloropropene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,2,3-Trichlorobenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,2,3-Trichloropropane	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,2,4-Trichlorobenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,2,4-Trimethylbenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,2-Dibromo-3-Chloropropane	µg/kg	4.7 U	5.3 U	5.5 U	--	380 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
1,2-dibromoethane (EDB)	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	5.0
1,2-Dichlorobenzene (o-Dichlorobenzene)	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	1.9	0.95 U	NE
1,2-Dichloroethane (EDC)	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,2-Dichloropropane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,3,5-Trimethylbenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,3-Dichlorobenzene (m-Dichlorobenzene)	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,3-Dichloropropane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
1,4-Dichlorobenzene (p-Dichlorobenzene)	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
2,2-Dichloropropane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
2-Butanone (MEK)	µg/kg	35	5.3 U	100	--	5.6 U	5.6 U	5.2 U	5.4 U	6.0	4.8 U	4.8 U	NE
2-Chloroethyl vinyl ether	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
2-Chlorotoluene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
2-Hexanone	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-1-0-2 12/13/11 0 - 2	DP-1-13-15 12/13/11 13 - 15	DP-3-0-2 12/13/11 0 - 2	DP-3-3-4 <sup>3</sup> 12/13/11 3 - 4	DP-4-0-2 12/13/11 0 - 2	DP-4-5-6 <sup>3</sup> 12/13/11 5 - 6	DP-5-0-2 12/13/11 0 - 2	DP-5-6-7 <sup>3</sup> 12/13/11 6 - 7	DP-6-0-2 12/13/11 0 - 2	DP-6-3.5-4.5 <sup>3</sup> 12/13/11 3.5 - 4.5	DP-6-11.5-12.5 12/13/11 11.5 - 12.5	MTCA <sup>2</sup> A ULU
4-Chlorotoluene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
4-Methyl-2-Pentanone (Methyl isobutyl ketone)	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
Acetone	µg/kg	200 E	46	520	--	5.6 U	5.6 U	5.2 U	5.4 U	53	16	4.8 U	NE
Benzene	µg/kg	1.4	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	30
Bromobenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Bromochloromethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Bromodichloromethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Bromoform (Tribromomethane)	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Bromomethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Carbon Disulfide	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Carbon Tetrachloride	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Chlorobenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Chloroethane	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
Chloroform	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Chloromethane	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
Cis-1,2-Dichloroethene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Cis-1,3-Dichloropropene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Dibromochloromethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Dibromomethane	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Dichlorodifluoromethane (CFC-12)	µg/kg	2.2	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Ethylbenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	6,000
Hexachlorobutadiene	µg/kg	4.7 U	5.3 U	5.5 U	--	380 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
Isopropylbenzene (Cumene)	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Methyl Iodide (Iodomethane)	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
Methyl t-butyl ether	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	100
Methylene Chloride	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	20
Naphthalene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	4.2	1.1 U	0.95 U	0.97 U	0.95 U	5,000
n-Butylbenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
n-Propylbenzene	µg/kg	4.3	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
p-Isopropyltoluene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Sec-Butylbenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Styrene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Tert-Butylbenzene	µg/kg	0.94 U	1.1 U	1.1 U	--	75 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Tetrachloroethene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	4.0	1.1 U	1.6	0.97 U	0.95 U	50
Toluene	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	7,000
Trans-1,2-Dichloroethene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Trans-1,3-Dichloropropene	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Trichloroethene (TCE)	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	30

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-1-0-2 12/13/11 0 - 2	DP-1-13-15 12/13/11 13 - 15	DP-3-0-2 12/13/11 0 - 2	DP-3-3-4 <sup>3</sup> 12/13/11 3 - 4	DP-4-0-2 12/13/11 0 - 2	DP-4-5-6 <sup>3</sup> 12/13/11 5 - 6	DP-5-0-2 12/13/11 0 - 2	DP-5-6-7 <sup>3</sup> 12/13/11 6 - 7	DP-6-0-2 12/13/11 0 - 2	DP-6-3.5-4.5 <sup>3</sup> 12/13/11 3.5 - 4.5	DP-6-11.5-12.5 12/13/11 11.5 - 12.5	MTCA <sup>2</sup> A ULU
Trichlorofluoromethane (CFC-11)	µg/kg	0.94 U	1.1 U	640	--	3,600	1.1 U	11	1.1 U	9.0	0.97 U	0.95 U	NE
Vinyl Acetate	µg/kg	4.7 U	5.3 U	5.5 U	--	5.6 U	5.6 U	5.2 U	5.4 U	4.8 U	4.8 U	4.8 U	NE
Vinyl Chloride	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Xylene, m-,p-	µg/kg	1.9 U	2.1 U	2.2 U	--	2.2 U	2.3 U	2.1 U	2.2 U	1.9 U	1.9 U	1.9 U	NE
Xylene, o-	µg/kg	0.94 U	1.1 U	1.1 U	--	1.1 U	1.1 U	1.0 U	1.1 U	0.95 U	0.97 U	0.95 U	NE
Xylene, Total	µg/kg	2.8 U	3.2 U	3.3 U	--	3.3 U	3.4 U	3.1 U	3.3 U	2.9 U	2.9 U	2.9 U	9,000
<b>SVOCs by SW8270</b>													
1,2-Diphenylhydrazine	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
1,3-Dinitrobenzene	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,3,4,6-Tetrachlorophenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,3,5,6-Tetrachlorophenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,3-DICHLOROANILINE	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,4,5-Trichlorophenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,4,6-Trichlorophenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,4-Dichlorophenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,4-Dimethylphenol	µg/kg	390 U	--	--	--	--	--	--	--	1,900 U	--	--	NE
2,4-Dinitrophenol	µg/kg	190 U	--	--	--	--	--	--	--	940 U	--	--	NE
2,4-Dinitrotoluene	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2,6-Dinitrotoluene	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2-Chloronaphthalene	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2-Chlorophenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2-Nitroaniline	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
2-Nitrophenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
3,3'-Dichlorobenzidine	µg/kg	390 U	--	--	--	--	--	--	--	1,900 U	--	--	NE
3-Nitroaniline	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
4,6-Dinitro-2-Methylphenol	µg/kg	190 U	--	--	--	--	--	--	--	940 U	--	--	NE
4-Bromophenyl phenyl ether	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
4-Chloro-3-Methylphenol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
4-Chloroaniline	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
4-Chlorophenyl-Phenylether	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
4-Nitroaniline	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
4-Nitrophenol (p-Nitrophenol)	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
Aniline	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
Benzene, 1,4-Dinitro-	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
Benzidine	µg/kg	390 U	--	--	--	--	--	--	--	1,900 U	--	--	NE
Benzyl Alcohol	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
Bis(2-Chloroethoxy)Methane	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE
Bis(2-Chloroethyl)Ether	µg/kg	39 U	--	--	--	--	--	--	--	190 U	--	--	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-1-0-2 12/13/11 0 - 2	DP-1-13-15 12/13/11 13 - 15	DP-3-0-2 12/13/11 0 - 2	DP-3-3-4 <sup>3</sup> 12/13/11 3 - 4	DP-4-0-2 12/13/11 0 - 2	DP-4-5-6 <sup>3</sup> 12/13/11 5 - 6	DP-5-0-2 12/13/11 0 - 2	DP-5-6-7 <sup>3</sup> 12/13/11 6 - 7	DP-6-0-2 12/13/11 0 - 2	DP-6-3.5-4.5 <sup>3</sup> 12/13/11 3.5 - 4.5	DP-6-11.5-12.5 12/13/11 11.5 - 12.5	MTCA <sup>2</sup> A ULU
Bis(2-chloroisopropyl) ether	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Bis(2-Ethylhexyl) Phthalate	µg/kg	190 U	-	-	-	-	-	-	-	970	-	-	NE
Butyl benzyl phthalate	µg/kg	390 U	-	-	-	-	-	-	-	1,900 U	-	-	NE
Carbazole	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Dibutyl phthalate	µg/kg	390 U	-	-	-	-	-	-	-	1,900 U	-	-	NE
Diethyl phthalate	µg/kg	190 U	-	-	-	-	-	-	-	940 U	-	-	NE
Dimethyl phthalate	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Di-N-Octyl Phthalate	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Hexachlorobenzene	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Hexachlorocyclopentadiene	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Hexachloroethane	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Hexanedioic Acid, Bis(2-Ethylhexyl) Ester	µg/kg	190 U	-	-	-	-	-	-	-	940 U	-	-	NE
Isophorone	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
m,p-Cresol	µg/kg	44	-	-	-	-	-	-	-	190 U	-	-	NE
Naphthalene	µg/kg	7.8 U	-	-	-	-	-	-	-	38	-	-	5,000
Nitrobenzene	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
N-Nitrosodimethylamine	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
N-Nitrosodi-n-propylamine	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
N-Nitrosodiphenylamine	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
o-Cresol (2-methylphenol)	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
O-DINITROBENZENE	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Pentachlorophenol	µg/kg	190 U	-	-	-	-	-	-	-	940 U	-	-	NE
Phenol	µg/kg	39 U	-	-	-	-	-	-	-	190 U	-	-	NE
Pyridine	µg/kg	390 U	-	-	-	-	-	-	-	1,900 U	-	-	NE

**Notes:**

<sup>1</sup> Chemical analysis performed by OnSite Environmental Inc. of Redmond, Washington.

<sup>2</sup> Model Toxics Control Act (MTCA) Cleanup Regulation Chapter 173-340 WAC. Method A Unrestricted Land Use (ULU) clean-up levels.

<sup>3</sup> Sample holding time was exceeded by one day for follow up analysis.

<sup>4</sup> The published natural background concentration for chromium in the Puget Sound region is 48 mg/kg (*Natural Background Soil Metals Concentrations in Washington State*. Publication #94-115. October 1994.)

<sup>5</sup> cPAH testing and regulatory evaluation is completed for individual carcinogenic compounds as well as the for the summation of the mixture of the seven carcinogenic PAHs (known as Ecology's toxicity equivalency methodology). The summation

PCBs = Polychlorinated biphenyls

PAHs = Polycyclic aromatic hydrocarbons

VOCs = Volatile organic compound

SVOCs = Semivolatile organic compound

µg/kg = microgram per kilogram

mg/kg = miligrams per kilogram

- = Not analyzed

U = The analyte was not detected at a concentration greater than the given RL or MDL.

E = The concentration reported exceeds the quantitation range and is an estimate.

H = The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and may be impacting the sample result

J = The value reported was below the practical quantitation limit. The value is an estimate.

NE = A cleanup level has not been established by Ecology.

**Bold Type** Exceeds MTCA A (Soil - Unrestricted Land Use Cleanup Level)

**TABLE 2B**  
**CHEMICAL ANALYTICAL DATA - SOIL<sup>1</sup>**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-7-0-2 12/14/11 0 - 2	DP-7-5-6 12/14/11 5 - 6	DP-8-0-2 12/14/11 0 - 2	DP-8-2-3 <sup>3</sup> 12/14/11 2 - 3	DP-9-0-2 12/14/11 0 - 2	DP-9-6-7 <sup>3</sup> 12/14/11 6 - 7	DP-10-0-2 12/14/11 0 - 2	DP-10-13-15 12/14/11 13 - 15	DP-DUPE-1 12/14/11 13 - 15	DP-11-0-2 12/14/11 2 - 2	DP-11-2-3.5 12/14/11 2 - 3.5	MTCA <sup>2</sup> A ULU
<b>Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx</b>													
Gasoline-range hydrocarbons	mg/kg	7.2 U	-	18	5.2 U	8.4 U	-	42	5.1 U	5.7 U	5.8 U	5.1 U	100
Diesel-range hydrocarbons	mg/kg	34	-	30 U	-	73 U	-	34 U	28 U	28 U	91 U	29 U	2,000
Lube Oil-range Hydrocarbons	mg/kg	130	-	150	-	360	-	120	56 U	56 U	480	58 U	2,000
<b>Metals by EPA6010B/7471A/SW7196A</b>													
Arsenic	mg/kg	13 U	-	11 U	-	14 U	-	12 U	11 U	11 U	12 U	12 U	20
Barium	mg/kg	88	-	46	-	84	-	70	58	56	82	74	NE
Cadmium	mg/kg	0.81	-	0.55 U	-	1.6	-	0.58 U	0.56 U	0.56 U	0.89	0.58 U	2.0
Chromium <sup>4</sup>	mg/kg	31	-	29	-	46	-	34	38	33	37	45	NE
Chromium, Hexavalent	mg/kg	-	-	-	-	-	-	-	-	-	-	-	19
Lead	mg/kg	93	-	24	-	85	-	8.3	5.6 U	5.6 U	26	5.8 U	250
Mercury	mg/kg	0.32 U	-	0.27 U	-	0.35 U	-	0.29 U	0.28 U	0.28 U	0.29 U	0.29 U	2.0
Selenium	mg/kg	13 U	-	11 U	-	14 U	-	12 U	11 U	11 U	12 U	12 U	NE
Silver	mg/kg	0.63 U	-	0.55 U	-	0.7 U	-	0.58 U	0.56 U	0.56 U	0.58 U	0.58 U	NE
<b>PCBs by SW8082</b>													
PCB-aroclor 1016	µg/kg	63 U	61 U	55 U	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	NE
PCB-aroclor 1221	µg/kg	63 U	61 U	55 U	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	NE
PCB-aroclor 1232	µg/kg	63 U	61 U	55 U	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	NE
PCB-aroclor 1242	µg/kg	120	61 U	55 U	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	NE
PCB-aroclor 1248	µg/kg	63 U	61 U	55 U	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	NE
PCB-aroclor 1254	µg/kg	63 U	61 U	58	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	NE
PCB-aroclor 1260	µg/kg	79	61 U	55 U	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	NE
Total Aroclors	µg/kg	199	61 U	58	57 U	70 U	-	58 U	56 U	56 U	58 U	58 U	1,000
<b>PAHs by SW8270</b>													
1-Methylnaphthalene	µg/kg	8.5 U	-	7.3 U	-	9.4 U	8.3 U	20	7.5 U	7.5 U	7.7 U	7.7 U	NE
2-Methylnaphthalene	µg/kg	8.8	-	7.3 U	-	13	8.3 U	37	7.5 U	7.5 U	7.7 U	7.7 U	NE
Acenaphthene	µg/kg	8.5 U	-	7.3 U	-	19	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE
Acenaphthylene	µg/kg	8.5 U	-	7.3 U	-	37	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE
Anthracene	µg/kg	8.5 U	-	7.3 U	-	72	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE
Benzo(a)anthracene	µg/kg	27	-	7.3 U	-	160	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE
Benzo(a)pyrene	µg/kg	8.5 U	-	7.3 U	-	140	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	100
Benzo(b)fluoranthene	µg/kg	9.7	-	9.1	-	170	8.3 U	7.8 U	7.5 U	7.5 U	9.9	7.7 U	NE
Benzo(ghi)perylene	µg/kg	8.5 U	-	7.3 U	-	76	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE



Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-7-0-2 12/14/11 0 - 2	DP-7-5-6 12/14/11 5 - 6	DP-8-0-2 12/14/11 0 - 2	DP-8-2-3 <sup>3</sup> 12/14/11 2 - 3	DP-9-0-2 12/14/11 0 - 2	DP-9-6-7 <sup>3</sup> 12/14/11 6 - 7	DP-10-0-2 12/14/11 0 - 2	DP-10-13-15 12/14/11 13 - 15	DP-DUPE-1 12/14/11 13 - 15	DP-11-0-2 12/14/11 2 - 2	DP-11-2-3.5 12/14/11 2 - 3.5	MTCA <sup>2</sup> A ULU
Benzo(j,k)fluoranthene	µg/kg	8.5 U	-	7.3 U	-	57	8.3 U	7.8 U	7.5 U	7.5 U	10	7.7 U	NE
Chrysene	µg/kg	8.5 U	-	7.3 U	-	150	8.3 U	7.8 U	7.5 U	7.5 U	11	7.7 U	NE
Dibenzo(a,h)anthracene	µg/kg	8.5 U	-	7.3 U	-	20	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE
Dibenzofuran	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Fluoranthene	µg/kg	16	-	8.5	-	350	8.3 U	10	7.5 U	7.5 U	15	7.7 U	NE
Fluorene	µg/kg	8.5 U	-	7.3 U	-	28	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE
Indeno(1,2,3-cd)pyrene	µg/kg	8.5 U	-	7.3 U	-	82	8.3 U	7.8 U	7.5 U	7.5 U	7.7 U	7.7 U	NE
Naphthalene	µg/kg	8.5 U	-	8.1	-	16	8.3 U	29	7.5 U	7.5 U	26	7.7 U	5,000
Phenanthrene	µg/kg	12	-	8.2	-	220	8.3 U	13	7.5 U	7.5 U	10	7.7 U	NE
Pyrene	µg/kg	15	-	10	-	280	8.3 U	11	7.5 U	7.5 U	22	7.7 U	NE
Total cPAH <sup>5</sup> TTEC (ND=0.5RL)	µg/kg	9.24	-	6.06	-	190.4	6.27 U	5.89 U	5.66 U	5.66 U	7.11	5.81 U	100
<b>VOCs by SW8260</b>													
1,1,1,2-Tetrachloroethane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,1,1-Trichloroethane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	2,000
1,1,2,2-Tetrachloroethane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,1,2-Trichloroethane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,1-Dichloroethane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,1-Dichloroethene	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,1-Dichloropropene	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,2,3-Trichlorobenzene	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,2,3-Trichloropropane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,2,4-Trichlorobenzene	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,2,4-Trimethylbenzene	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	2,500	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,2-Dibromo-3-Chloropropane	µg/kg	5.9 U	-	4.6 U	-	6.7 U	-	290 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
1,2-dibromoethane (EDB)	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	5.0
1,2-Dichlorobenzene (o-Dichlorobenzene)	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,2-Dichloroethane (EDC)	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,2-Dichloropropane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,3,5-Trimethylbenzene	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	840	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,3-Dichlorobenzene (m-Dichlorobenzene)	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,3-Dichloropropane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
1,4-Dichlorobenzene (p-Dichlorobenzene)	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
2,2-Dichloropropane	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
2-Butanone (MEK)	µg/kg	5.9 U	-	4.6 U	-	9.1	-	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
2-Chloroethyl vinyl ether	µg/kg	5.9 U	-	4.6 U	-	6.7 U	-	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
2-Chlorotoluene	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
2-Hexanone	µg/kg	5.9 U	-	4.6 U	-	6.7 U	-	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-7-0-2 12/14/11 0 - 2	DP-7-5-6 12/14/11 5 - 6	DP-8-0-2 12/14/11 0 - 2	DP-8-2-3 <sup>3</sup> 12/14/11 2 - 3	DP-9-0-2 12/14/11 0 - 2	DP-9-6-7 <sup>3</sup> 12/14/11 6 - 7	DP-10-0-2 12/14/11 0 - 2	DP-10-13-15 12/14/11 13 - 15	DP-DUPE-1 12/14/11 13 - 15	DP-11-0-2 12/14/11 2 - 2	DP-11-2-3.5 12/14/11 2 - 3.5	MTCA <sup>2</sup> A ULU
4-Chlorotoluene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
4-Methyl-2-Pentanone (Methyl isobutyl ketone)	µg/kg	5.9 U	–	4.6 U	–	6.7 U	–	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
Acetone	µg/kg	5.9 U	–	4.6 U	–	97	–	4.8 U	4.5 U	4.9 U	49	4.4 U	NE
Benzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	13	0.9 U	0.97 U	1.1 U	0.89 U	30
Bromobenzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Bromochloromethane	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Bromodichloromethane	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Bromoform (Tribromomethane)	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Bromomethane	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Carbon Disulfide	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Carbon Tetrachloride	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Chlorobenzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Chloroethane	µg/kg	5.9 U	–	4.6 U	–	6.7 U	–	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
Chloroform	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Chloromethane	µg/kg	5.9 U	–	4.6 U	–	6.7 U	–	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
Cis-1,2-Dichloroethene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Cis-1,3-Dichloropropene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Dibromochloromethane	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Dibromomethane	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Dichlorodifluoromethane (CFC-12)	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Ethylbenzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	230	0.9 U	0.97 U	1.1 U	0.89 U	6,000
Hexachlorobutadiene	µg/kg	5.9 U	–	4.6 U	–	6.7 U	–	290 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
Isopropylbenzene (Cumene)	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	40	0.9 U	0.97 U	1.1 U	0.89 U	NE
Methyl Iodide (Iodomethane)	µg/kg	5.9 U	–	4.6 U	–	6.7 U	–	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
Methyl t-butyl ether	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	100
Methylene Chloride	µg/kg	5.9 U	–	4.6 U	–	6.7 U	–	9.5 H	4.5 U	4.9 U	5.4 U	4.4 U	20
Naphthalene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	660	0.9 U	0.97 U	1.1 U	0.89 U	5,000
n-Butylbenzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
n-Propylbenzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	270	0.9 U	0.97 U	1.1 U	0.89 U	NE
p-Isopropyltoluene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Sec-Butylbenzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	98	0.9 U	0.97 U	1.1 U	0.89 U	NE
Styrene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	5.5	0.9 U	0.97 U	1.1 U	0.89 U	NE
Tert-Butylbenzene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	59 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Tetrachloroethene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	19	0.9 U	0.97 U	1.1 U	0.89 U	50
Toluene	µg/kg	5.9 U	–	4.6 U	–	6.7 U	–	410	4.5 U	4.9 U	5.4 U	4.4 U	7,000
Trans-1,2-Dichloroethene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Trans-1,3-Dichloropropene	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Trichloroethene (TCE)	µg/kg	1.2 U	–	0.92 U	–	1.3 U	–	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	30

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-7-0-2 12/14/11 0 - 2	DP-7-5-6 12/14/11 5 - 6	DP-8-0-2 12/14/11 0 - 2	DP-8-2-3 <sup>3</sup> 12/14/11 2 - 3	DP-9-0-2 12/14/11 0 - 2	DP-9-6-7 <sup>3</sup> 12/14/11 6 - 7	DP-10-0-2 12/14/11 0 - 2	DP-10-13-15 12/14/11 13 - 15	DP-DUPE-1 12/14/11 13 - 15	DP-11-0-2 12/14/11 2 - 2	DP-11-2-3.5 12/14/11 2 - 3.5	MTCA <sup>2</sup> A ULU
Trichlorofluoromethane (CFC-11)	µg/kg	1.2 U	-	0.92 U	-	3.7	-	780	0.9 U	0.97 U	1.1 U	0.89 U	NE
Vinyl Acetate	µg/kg	5.9 U	-	4.6 U	-	6.7 U	-	4.8 U	4.5 U	4.9 U	5.4 U	4.4 U	NE
Vinyl Chloride	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	0.96 U	0.9 U	0.97 U	1.1 U	0.89 U	NE
Xylene, m-,p-	µg/kg	2.3 U	-	1.8 U	-	2.7 U	-	1,200	1.8 U	1.9 U	2.2 U	1.8 U	NE
Xylene, o-	µg/kg	1.2 U	-	0.92 U	-	1.3 U	-	620	0.9 U	0.97 U	1.1 U	0.89 U	NE
Xylene, Total	µg/kg	3.5 U	-	2.7 U	-	4.0 U	-	1,820	2.7 U	2.9 U	3.3 U	2.7 U	9,000
<b>SVOCs by SW8270</b>													
1,2-Diphenylhydrazine	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
1,3-Dinitrobenzene	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,3,4,6-Tetrachlorophenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,3,5,6-Tetrachlorophenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,3-DICHLOROANILINE	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,4,5-Trichlorophenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,4,6-Trichlorophenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,4-Dichlorophenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,4-Dimethylphenol	µg/kg	-	-	-	-	-	-	-	-	-	1900 U	-	NE
2,4-Dinitrophenol	µg/kg	-	-	-	-	-	-	-	-	-	960 U	-	NE
2,4-Dinitrotoluene	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2,6-Dinitrotoluene	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2-Chloronaphthalene	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2-Chlorophenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2-Nitroaniline	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
2-Nitrophenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
3,3'-Dichlorobenzidine	µg/kg	-	-	-	-	-	-	-	-	-	1900 U	-	NE
3-Nitroaniline	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
4,6-Dinitro-2-Methylphenol	µg/kg	-	-	-	-	-	-	-	-	-	960 U	-	NE
4-Bromophenyl phenyl ether	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
4-Chloro-3-Methylphenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
4-Chloroaniline	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
4-Chlorophenyl-Phenylether	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
4-Nitroaniline	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
4-Nitrophenol (p-Nitrophenol)	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Aniline	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Benzene, 1,4-Dinitro-	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Benzidine	µg/kg	-	-	-	-	-	-	-	-	-	1900 U	-	NE
Benzyl Alcohol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Bis(2-Chloroethoxy)Methane	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Bis(2-Chloroethyl)Ether	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Bis(2-chloroisopropyl) ether	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-7-0-2 12/14/11 0 - 2	DP-7-5-6 12/14/11 5 - 6	DP-8-0-2 12/14/11 0 - 2	DP-8-2-3 <sup>3</sup> 12/14/11 2 - 3	DP-9-0-2 12/14/11 0 - 2	DP-9-6-7 <sup>3</sup> 12/14/11 6 - 7	DP-10-0-2 12/14/11 0 - 2	DP-10-13-15 12/14/11 13 - 15	DP-DUPE-1 12/14/11 13 - 15	DP-11-0-2 12/14/11 2 - 2	DP-11-2-3.5 12/14/11 2 - 3.5	MTCA <sup>2</sup> A ULU
Bis(2-Ethylhexyl) Phthalate	µg/kg	-	-	-	-	-	-	-	-	-	46,000	-	NE
Butyl benzyl phthalate	µg/kg	-	-	-	-	-	-	-	-	-	1900 U	-	NE
Carbazole	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Dibutyl phthalate	µg/kg	-	-	-	-	-	-	-	-	-	1900 U	-	NE
Diethyl phthalate	µg/kg	-	-	-	-	-	-	-	-	-	960 U	-	NE
Dimethyl phthalate	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Di-N-Octyl Phthalate	µg/kg	-	-	-	-	-	-	-	-	-	6,700	-	NE
Hexachlorobenzene	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Hexachlorocyclopentadiene	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Hexachloroethane	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Hexanedioic Acid, Bis(2-Ethylhexyl) Ester	µg/kg	-	-	-	-	-	-	-	-	-	960 U	-	NE
Isophorone	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
m,p-Cresol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Naphthalene	µg/kg	-	-	-	-	-	-	-	-	-	26	-	5,000
Nitrobenzene	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
N-Nitrosodimethylamine	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
N-Nitrosodi-n-propylamine	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
N-Nitrosodiphenylamine	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
o-Cresol (2-methylphenol)	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
O-DINITROBENZENE	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Pentachlorophenol	µg/kg	-	-	-	-	-	-	-	-	-	960 U	-	NE
Phenol	µg/kg	-	-	-	-	-	-	-	-	-	190 U	-	NE
Pyridine	µg/kg	-	-	-	-	-	-	-	-	-	1900 U	-	NE

**Notes:**

<sup>1</sup> Chemical analysis performed by OnSite Environmental Inc. of Redmond, Washington.

<sup>2</sup> Model Toxics Control Act (MTCA) Cleanup Regulation Chapter 173-340 WAC. Method A Unrestricted Land Use (ULU) clean-up levels.

<sup>3</sup> Sample holding time was exceeded by one day for follow up analysis.

<sup>4</sup> The published natural background concentration for chromium in the Puget Sound region is 48 mg/kg (*Natural Background Soil Metals Concentrations in Washington State*. Publication #94-115. October 1994.)

<sup>5</sup> cPAH testing and regulatory evaluation is completed for individual carcinogenic compounds as well as the for the summation of the mixture of the seven carcinogenic PAHs (known as Ecology's toxicity equivalency methodology). The summation procedure is completed using toxicity equivalency factors for each individual compound which are then added to produce a total toxicity equivalent concentration (TTEC) which is then compared to the MTCA cleanup level of 0.1 mg/kg (or 100 µg/kg).

PCBs = Polychlorinated biphenyls

PAHs = Polycyclic aromatic hydrocarbons

VOCs = Volatile organic compound

SVOCs = Semivolatile organic compound

µg/kg = microgram per kilogram

mg/kg = milligrams per kilogram

- = Not analyzed

U = The analyte was not detected at a concentration greater than the given RL or MDL.

E = The concentration reported exceeds the quantitation range and is an estimate.

H = The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and may be impacting the sample result

J = The value reported was below the practical quantitation limit. The value is an estimate.

NE = A cleanup level has not been established by Ecology.

**Bold Type** Exceeds MTCA A (Soil - Unrestricted Land Use Cleanup Level)

**TABLE 2C**  
**CHEMICAL ANALYTICAL DATA - SOIL<sup>1</sup>**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-12-0.0-2.0 2/08/12 0 - 2	DP-12-5.0-7.5 2/08/12 5 - 7.5	DP-13-0.0-1.5 2/08/12 0 - 1.5	DP-14-0.0-2.0 2/08/12 0 - 2	DP-15-0.0-2.0 2/08/12 0 - 2	DP-15-5.0-7.0 2/08/12 5 - 7	DP-16-0.0-2.0 2/08/12 0 - 2	DP-17-0.0-2.0 2/08/12 0 - 2	DP-18-0.0-2.0 2/08/12 0 - 2	DP-19-0.0-2.0 2/08/12 0 - 2	DP-19-2.0-3.0 2/08/12 2 - 3	MTCA <sup>2</sup> A ULU
<b>Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx</b>													
Gasoline-range hydrocarbons	mg/kg	5.3 U	4.9 U	6.1 U	5.7 U	7.2 U	6.6 U	5 U	5.2 U	5.4 U	5 U	-	100
Diesel-range hydrocarbons	mg/kg	28 U	30 U	31 U	28 U	30 U	32 U	29 U	29 U	-	-	-	2,000
Lube Oil-range Hydrocarbons	mg/kg	130	59 U	63 U	56 U	60 U	65 U	58 U	58 U	-	-	-	2,000
<b>Metals by EPA6010B/7471A/SW7196A</b>													
Arsenic	mg/kg	11 U	12 U	13 U	11 U	12 U	13 U	12 U	12 U	-	-	-	20
Barium	mg/kg	75	29	97	43	110	62	45	55	-	-	-	NE
Cadmium	mg/kg	0.55 U	0.59 U	0.63 U	0.56 U	0.59 U	0.65 U	0.58 U	0.58 U	-	-	-	2.0
Chromium <sup>4</sup>	mg/kg	34	35	37	31	25	26	28	31	-	-	-	NE
Chromium, Hexavalent	mg/kg	-	-	1.3 U	-	-	-	-	-	-	-	-	19
Lead	mg/kg	19	5.9 U	6.3 U	5.6 U	5.9 U	6.5 U	5.8 U	5.8 U	-	-	-	250
Mercury	mg/kg	0.28 U	0.3 U	0.31 U	0.28 U	0.3 U	0.32 U	0.29 U	0.29 U	-	-	-	2.0
Selenium	mg/kg	11 U	12 U	13 U	11 U	12 U	13 U	12 U	12 U	-	-	-	NE
Silver	mg/kg	0.55 U	0.59 U	0.63 U	0.56 U	0.59 U	0.65 U	0.58 U	0.58 U	-	-	-	NE
<b>PCBs by SW8082</b>													
PCB-aroclor 1016	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	NE
PCB-aroclor 1221	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	NE
PCB-aroclor 1232	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	NE
PCB-aroclor 1242	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	NE
PCB-aroclor 1248	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	NE
PCB-aroclor 1254	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	NE
PCB-aroclor 1260	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	NE
Total Aroclors	µg/kg	55 U	59 U	63 U	56 U	59 U	65 U	58 U	58 U	-	-	-	1,000
<b>PAHs by SW8270</b>													
1-Methylnaphthalene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE
2-Methylnaphthalene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE
Acenaphthene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE
Acenaphthylene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE
Anthracene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE
Benzo(a)anthracene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE
Benzo(a)pyrene	µg/kg	21	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	100
Benzo(b)fluoranthene	µg/kg	11	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE
Benzo(ghi)perylene	µg/kg	46	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	-	-	-	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-12-0.0-2.0 2/08/12 0 - 2	DP-12-5.0-7.5 2/08/12 5 - 7.5	DP-13-0.0-1.5 2/08/12 0 - 1.5	DP-14-0.0-2.0 2/08/12 0 - 2	DP-15-0.0-2.0 2/08/12 0 - 2	DP-15-5.0-7.0 2/08/12 5 - 7	DP-16-0.0-2.0 2/08/12 0 - 2	DP-17-0.0-2.0 2/08/12 0 - 2	DP-18-0.0-2.0 2/08/12 0 - 2	DP-19-0.0-2.0 2/08/12 0 - 2	DP-19-2.0-3.0 2/08/12 2 - 3	MTCA <sup>2</sup> A ULU
Benzo(j,k)fluoranthene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Chrysene	µg/kg	9.6	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Dibenzo(a,h)anthracene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Dibenzofuran	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Fluoranthene	µg/kg	7.4 U	7.9 U	17	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Fluorene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Indeno(1,2,3-cd)pyrene	µg/kg	10	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Naphthalene	µg/kg	7.4 U	7.9 U	8.4 U	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	5,000
Phenanthrene	µg/kg	7.4 U	7.9 U	11	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Pyrene	µg/kg	11	7.9 U	16	7.5 U	7.9 U	8.6 U	7.7 U	7.7 U	--	--	--	NE
Total cPAH <sup>5</sup> TTEC (ND=0.5RL)	µg/kg	24.31	5.96 U	6.34 U	5.66 U	5.96 U	6.49 U	5.81 U	5.81 U	--	--	--	100
<b>VOCs by SW8260</b>													
1,1,1,2-Tetrachloroethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,1,1-Trichloroethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	2,000
1,1,2,2-Tetrachloroethane	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,1,2-Trichloroethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,1-Dichloroethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,1-Dichloroethene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,1-Dichloropropene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,2,3-Trichlorobenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,2,3-Trichloropropane	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,2,4-Trichlorobenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,2,4-Trimethylbenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	21	0.97 U	NE
1,2-Dibromo-3-Chloropropane	µg/kg	4.1 U	4.2 U	310 U	4.1 U	270 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
1,2-dibromoethane (EDB)	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	5.0
1,2-Dichlorobenzene (o-Dichlorobenzene)	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,2-Dichloroethane (EDC)	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,2-Dichloropropane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,3,5-Trimethylbenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	13	0.97 U	NE
1,3-Dichlorobenzene (m-Dichlorobenzene)	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,3-Dichloropropane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
1,4-Dichlorobenzene (p-Dichlorobenzene)	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
2,2-Dichloropropane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
2-Butanone (MEK)	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
2-Chloroethyl vinyl ether	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
2-Chlorotoluene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
2-Hexanone	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-12-0.0-2.0 2/08/12 0 - 2	DP-12-5.0-7.5 2/08/12 5 - 7.5	DP-13-0.0-1.5 2/08/12 0 - 1.5	DP-14-0.0-2.0 2/08/12 0 - 2	DP-15-0.0-2.0 2/08/12 0 - 2	DP-15-5.0-7.0 2/08/12 5 - 7	DP-16-0.0-2.0 2/08/12 0 - 2	DP-17-0.0-2.0 2/08/12 0 - 2	DP-18-0.0-2.0 2/08/12 0 - 2	DP-19-0.0-2.0 2/08/12 0 - 2	DP-19-2.0-3.0 2/08/12 2 - 3	MTCA <sup>2</sup> A ULU
4-Chlorotoluene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
4-Methyl-2-Pentanone (Methyl isobutyl ketone)	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	6.6	4.8 U	NE
Acetone	µg/kg	4.1 U	13	23	4.1 U	100	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
Benzene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	30
Bromobenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Bromochloromethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Bromodichloromethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Bromoform (Tribromomethane)	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Bromomethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Carbon Disulfide	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Carbon Tetrachloride	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Chlorobenzene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Chloroethane	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
Chloroform	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Chloromethane	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
Cis-1,2-Dichloroethene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Cis-1,3-Dichloropropene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Dibromochloromethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Dibromomethane	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Dichlorodifluoromethane (CFC-12)	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Ethylbenzene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	3.5	0.97 U	6,000
Hexachlorobutadiene	µg/kg	4.1 U	4.2 U	310 U	4.1 U	270 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
Isopropylbenzene (Cumene)	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Methyl Iodide (Iodomethane)	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
Methyl t-butyl ether	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	100
Methylene Chloride	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	20
Naphthalene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	2.0	0.97 U	5,000
n-Butylbenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	1.6	0.97 U	NE
n-Propylbenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	2.5	0.97 U	NE
p-Isopropyltoluene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Sec-Butylbenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Styrene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Tert-Butylbenzene	µg/kg	0.83 U	0.83 U	63 U	0.83 U	55 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Tetrachloroethene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	50
Toluene	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	6.6	4.8 U	7,000
Trans-1,2-Dichloroethene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Trans-1,3-Dichloropropene	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Trichloroethene (TCE)	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	30



Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-12-0.0-2.0 2/08/12 0 - 2	DP-12-5.0-7.5 2/08/12 5 - 7.5	DP-13-0.0-1.5 2/08/12 0 - 1.5	DP-14-0.0-2.0 2/08/12 0 - 2	DP-15-0.0-2.0 2/08/12 0 - 2	DP-15-5.0-7.0 2/08/12 5 - 7	DP-16-0.0-2.0 2/08/12 0 - 2	DP-17-0.0-2.0 2/08/12 0 - 2	DP-18-0.0-2.0 2/08/12 0 - 2	DP-19-0.0-2.0 2/08/12 0 - 2	DP-19-2.0-3.0 2/08/12 2 - 3	MTCA <sup>2</sup> A ULU
Trichlorofluoromethane (CFC-11)	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	2.6	0.97 U	NE
Vinyl Acetate	µg/kg	4.1 U	4.2 U	4.8 U	4.1 U	5.4 U	5.4 U	4.4 U	4.7 U	5.3 U	4.4 U	4.8 U	NE
Vinyl Chloride	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	0.88 U	0.97 U	NE
Xylene, m-,p-	µg/kg	1.7 U	1.7 U	1.9 U	1.7 U	2.1 U	2.1 U	1.8 U	1.9 U	2.1 U	21	1.9 U	NE
Xylene, o-	µg/kg	0.83 U	0.83 U	0.97 U	0.83 U	1.1 U	1.1 U	0.88 U	0.94 U	1.1 U	13	0.97 U	NE
Xylene, Total	µg/kg	2.5 U	2.5 U	2.9 U	2.5 U	3.2 U	3.2 U	2.7 U	2.8 U	3.2 U	34	2.9 U	9,000
<b>SVOCs by SW8270</b>													
1,2-Diphenylhydrazine	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
1,3-Dinitrobenzene	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,3,4,6-Tetrachlorophenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,3,5,6-Tetrachlorophenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,3-DICHLOROANILINE	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,4,5-Trichlorophenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,4,6-Trichlorophenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,4-Dichlorophenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,4-Dimethylphenol	µg/kg	370 U	--	--	--	400 U	--	--	--	--	--	--	NE
2,4-Dinitrophenol	µg/kg	180 U	--	--	--	200 U	--	--	--	--	--	--	NE
2,4-Dinitrotoluene	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2,6-Dinitrotoluene	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2-Chloronaphthalene	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2-Chlorophenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2-Nitroaniline	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
2-Nitrophenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
3,3'-Dichlorobenzidine	µg/kg	370 U	--	--	--	400 U	--	--	--	--	--	--	NE
3-Nitroaniline	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
4,6-Dinitro-2-Methylphenol	µg/kg	180 U	--	--	--	200 U	--	--	--	--	--	--	NE
4-Bromophenyl phenyl ether	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
4-Chloro-3-Methylphenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
4-Chloroaniline	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
4-Chlorophenyl-Phenylether	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
4-Nitroaniline	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
4-Nitrophenol (p-Nitrophenol)	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Aniline	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Benzene, 1,4-Dinitro-	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Benzidine	µg/kg	370 U	--	--	--	400 U	--	--	--	--	--	--	NE
Benzyl Alcohol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Bis(2-Chloroethoxy)Methane	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Bis(2-Chloroethyl)Ether	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Bis(2-chloroisopropyl) ether	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE



Analytes	Units	Sample ID, Date and Depth (in ft.)											Screening Levels
		DP-12-0.0-2.0 2/08/12 0 - 2	DP-12-5.0-7.5 2/08/12 5 - 7.5	DP-13-0.0-1.5 2/08/12 0 - 1.5	DP-14-0.0-2.0 2/08/12 0 - 2	DP-15-0.0-2.0 2/08/12 0 - 2	DP-15-5.0-7.0 2/08/12 5 - 7	DP-16-0.0-2.0 2/08/12 0 - 2	DP-17-0.0-2.0 2/08/12 0 - 2	DP-18-0.0-2.0 2/08/12 0 - 2	DP-19-0.0-2.0 2/08/12 0 - 2	DP-19-2.0-3.0 2/08/12 2 - 3	MTCA <sup>2</sup> A ULU
Bis(2-Ethylhexyl) Phthalate	µg/kg	230	--	--	--	200 U	--	--	--	--	--	--	NE
Butyl benzyl phthalate	µg/kg	370 U	--	--	--	400 U	--	--	--	--	--	--	NE
Carbazole	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Dibutyl phthalate	µg/kg	370 U	--	--	--	400 U	--	--	--	--	--	--	NE
Diethyl phthalate	µg/kg	180 U	--	--	--	200 U	--	--	--	--	--	--	NE
Dimethyl phthalate	µg/kg	120	--	--	--	40 U	--	--	--	--	--	--	NE
Di-N-Octyl Phthalate	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Hexachlorobenzene	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Hexachlorocyclopentadiene	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Hexachloroethane	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Hexanedioic Acid, Bis(2-Ethylhexyl) Ester	µg/kg	180 U	--	--	--	200 U	--	--	--	--	--	--	NE
Isophorone	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
m,p-Cresol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Naphthalene	µg/kg	7.4 U	--	--	--	7.9 U	--	--	--	--	--	--	5,000
Nitrobenzene	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
N-Nitrosodimethylamine	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
N-Nitrosodi-n-propylamine	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
N-Nitrosodiphenylamine	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
o-Cresol (2-methylphenol)	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
O-DINITROBENZENE	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Pentachlorophenol	µg/kg	180 U	--	--	--	200 U	--	--	--	--	--	--	NE
Phenol	µg/kg	37 U	--	--	--	40 U	--	--	--	--	--	--	NE
Pyridine	µg/kg	370 U	--	--	--	400 U	--	--	--	--	--	--	NE

**Notes:**

- <sup>1</sup> Chemical analysis performed by OnSite Environmental Inc. of Redmond, Washington.
- <sup>2</sup> Model Toxics Control Act (MTCA) Cleanup Regulation Chapter 173-340 WAC. Method A Unrestricted Land Use (ULU) clean-up levels.
- <sup>3</sup> Sample holding time was exceeded by one day for follow up analysis.
- <sup>4</sup> The published natural background concentration for chromium in the Puget Sound region is 48 mg/kg (*Natural Background Soil Metals Concentrations in Washington State*. Publication #94-115. October 1994.)
- <sup>5</sup> cPAH testing and regulatory evaluation is completed for individual carcinogenic compounds as well as the for the summation of the mixture of the seven carcinogenic PAHs (known as Ecology's toxicity equivalency methodology). The summation procedure is completed using toxicity equivalency factors for each individual compound which are then added to produce a total toxicity equivalent concentration (TEEC) which is then compared to the MTCA cleanup level of 0.1 mg/kg (or 100 µg/kg).

PCBs = Polychlorinated biphenyls  
PAHs = Polycyclic aromatic hydrocarbons  
VOCs = Volatile organic compound  
SVOCs = Semivolatile organic compound  
µg/kg = microgram per kilogram  
mg/kg = milligrams per kilogram  
-- = Not analyzed

U = The analyte was not detected at a concentration greater than the given RL or MDL.  
E = The concentration reported exceeds the quantitation range and is an estimate.  
H = The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and may be impacting the sample result  
J = The value reported was below the practical quantitation limit. The value is an estimate.  
NE = A cleanup level has not been established by Ecology.

**Bold Type** Exceeds MTCA A (Soil - Unrestricted Land Use Cleanup Level)

**TABLE 2D**  
**CHEMICAL ANALYTICAL DATA - SOIL<sup>1</sup>**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Analytes	Units	Sample ID, Date and Depth (in ft.)							Screening Levels
		DP-20-0.0-2.0 2/08/12 0 - 2	DP-21-0.0-2.0 2/08/12 0 - 2	DP-21-2.0-3.0 2/08/12 2 - 3	HA1-0.0-2.0 2/09/12 0 - 2	HA2-0.0-2.0 2/09/12 0 - 2	HA3-0.0-2.0 2/09/12 0 - 2	HA4-0.0-2.0 2/09/12 0 - 2	MTCA <sup>2</sup> A ULU
<b>Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx</b>									
	Units								
Gasoline-range hydrocarbons	mg/kg	5.5 U	--	--	--	--	--	--	100
Diesel-range hydrocarbons	mg/kg	--	100	32 U	31 U	34 U	110 U	--	2,000
Lube Oil-range Hydrocarbons	mg/kg	--	470	63 U	62 U	67 U	660	--	2,000
<b>Metals by EPA6010B/7471A/SW7196A</b>									
Arsenic	mg/kg	--	12 U	13 U	12 U	13 U	12 U	--	20
Barium	mg/kg	--	81	86	120	93	91	--	NE
Cadmium	mg/kg	--	0.83	0.63 U	0.62 U	0.67 U	2.7	--	2.0
Chromium <sup>4</sup>	mg/kg	--	45	44	33	37	58	--	NE
Chromium, Hexavalent	mg/kg	--	--	--	--	--	--	1.3 U	19
Lead	mg/kg	--	48	6.3 U	6.2 U	9.6	150	--	250
Mercury	mg/kg	--	0.29 U	0.32 U	0.31 U	0.34 U	0.3 U	--	2.0
Selenium	mg/kg	--	12 U	13 U	12 U	13 U	12 U	--	NE
Silver	mg/kg	--	0.58 U	0.63 U	0.62 U	0.67 U	0.6 U	--	NE
<b>PCBs by SW8082</b>									
PCB-aroclor 1016	µg/kg	--	--	--	--	--	--	--	NE
PCB-aroclor 1221	µg/kg	--	--	--	--	--	--	--	NE
PCB-aroclor 1232	µg/kg	--	--	--	--	--	--	--	NE
PCB-aroclor 1242	µg/kg	--	--	--	--	--	--	--	NE
PCB-aroclor 1248	µg/kg	--	--	--	--	--	--	--	NE
PCB-aroclor 1254	µg/kg	--	--	--	--	--	--	--	NE
PCB-aroclor 1260	µg/kg	--	--	--	--	--	--	--	NE
Total Aroclors	µg/kg	--	--	--	--	--	--	--	1,000
<b>PAHs by SW8270</b>									
1-Methylnaphthalene	µg/kg	--	--	--	--	--	--	10	NE
2-Methylnaphthalene	µg/kg	--	--	--	--	--	--	19	NE
Acenaphthene	µg/kg	--	--	--	--	--	--	8.4 U	NE
Acenaphthylene	µg/kg	--	--	--	--	--	--	8.4 U	NE
Anthracene	µg/kg	--	--	--	--	--	--	8.4 U	NE
Benzo(a)anthracene	µg/kg	--	--	--	--	--	--	13	NE
Benzo(a)pyrene	µg/kg	--	--	--	--	--	--	13	100
Benzo(b)fluoranthene	µg/kg	--	--	--	--	--	--	18	NE
Benzo(ghi)perylene	µg/kg	--	--	--	--	--	--	19	NE
Benzo(j,k)fluoranthene	µg/kg	--	--	--	--	--	--	8.4 U	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)							Screening Levels
		DP-20-0.0-2.0 2/08/12 0 - 2	DP-21-0.0-2.0 2/08/12 0 - 2	DP-21-2.0-3.0 2/08/12 2 - 3	HA1-0.0-2.0 2/09/12 0 - 2	HA2-0.0-2.0 2/09/12 0 - 2	HA3-0.0-2.0 2/09/12 0 - 2	HA4-0.0-2.0 2/09/12 0 - 2	MTCA <sup>2</sup> A ULU
Chrysene	µg/kg	--	--	--	--	--	--	17	NE
Dibenzo(a,h)anthracene	µg/kg	--	--	--	--	--	--	8.4 U	NE
Dibenzofuran	µg/kg	--	--	--	--	--	--	--	NE
Fluoranthene	µg/kg	--	--	--	--	--	--	36	NE
Fluorene	µg/kg	--	--	--	--	--	--	8.4 U	NE
Indeno(1,2,3-cd)pyrene	µg/kg	--	--	--	--	--	--	13	NE
Naphthalene	µg/kg	--	--	--	--	--	--	21	5,000
Phenanthrene	µg/kg	--	--	--	--	--	--	31	NE
Pyrene	µg/kg	--	--	--	--	--	--	38	NE
Total cPAH <sup>5</sup> TTEC (ND=0.5RL)	µg/kg	--	--	--	--	--	--	17.57 <sup>3</sup>	100
<b>VOCs by SW8260</b>									
1,1,1,2-Tetrachloroethane	µg/kg	1 U	--	--	--	--	--	--	NE
1,1,1-Trichloroethane	µg/kg	1 U	--	--	--	--	--	--	2,000
1,1,2,2-Tetrachloroethane	µg/kg	1 U	--	--	--	--	--	--	NE
1,1,2-Trichloroethane	µg/kg	1 U	--	--	--	--	--	--	NE
1,1-Dichloroethane	µg/kg	1 U	--	--	--	--	--	--	NE
1,1-Dichloroethene	µg/kg	1 U	--	--	--	--	--	--	NE
1,1-Dichloropropene	µg/kg	1 U	--	--	--	--	--	--	NE
1,2,3-Trichlorobenzene	µg/kg	1 U	--	--	--	--	--	--	NE
1,2,3-Trichloropropane	µg/kg	1 U	--	--	--	--	--	--	NE
1,2,4-Trichlorobenzene	µg/kg	1 U	--	--	--	--	--	--	NE
1,2,4-Trimethylbenzene	µg/kg	1 U	--	--	--	--	--	--	NE
1,2-Dibromo-3-Chloropropane	µg/kg	5.1 U	--	--	--	--	--	--	NE
1,2-dibromoethane (EDB)	µg/kg	1 U	--	--	--	--	--	--	5.0
1,2-Dichlorobenzene (o-Dichlorobenzene)	µg/kg	1 U	--	--	--	--	--	--	NE
1,2-Dichloroethane (EDC)	µg/kg	1 U	--	--	--	--	--	--	NE
1,2-Dichloropropane	µg/kg	1 U	--	--	--	--	--	--	NE
1,3,5-Trimethylbenzene	µg/kg	1 U	--	--	--	--	--	--	NE
1,3-Dichlorobenzene (m-Dichlorobenzene)	µg/kg	1 U	--	--	--	--	--	--	NE
1,3-Dichloropropane	µg/kg	1 U	--	--	--	--	--	--	NE
1,4-Dichlorobenzene (p-Dichlorobenzene)	µg/kg	1 U	--	--	--	--	--	--	NE
2,2-Dichloropropane	µg/kg	1 U	--	--	--	--	--	--	NE
2-Butanone (MEK)	µg/kg	5.1 U	--	--	--	--	--	--	NE
2-Chloroethyl vinyl ether	µg/kg	5.1 U	--	--	--	--	--	--	NE
2-Chlorotoluene	µg/kg	1 U	--	--	--	--	--	--	NE
2-Hexanone	µg/kg	5.1 U	--	--	--	--	--	--	NE
4-Chlorotoluene	µg/kg	1 U	--	--	--	--	--	--	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)							Screening Levels
		DP-20-0.0-2.0 2/08/12 0 - 2	DP-21-0.0-2.0 2/08/12 0 - 2	DP-21-2.0-3.0 2/08/12 2 - 3	HA1-0.0-2.0 2/09/12 0 - 2	HA2-0.0-2.0 2/09/12 0 - 2	HA3-0.0-2.0 2/09/12 0 - 2	HA4-0.0-2.0 2/09/12 0 - 2	MTCA <sup>2</sup> A ULU
4-Methyl-2-Pentanone (Methyl isobutyl ketone)	µg/kg	5.1 U	--	--	--	--	--	--	NE
Acetone	µg/kg	5.1 U	--	--	--	--	--	--	NE
Benzene	µg/kg	1 U	--	--	--	--	--	--	30
Bromobenzene	µg/kg	1 U	--	--	--	--	--	--	NE
Bromochloromethane	µg/kg	1 U	--	--	--	--	--	--	NE
Bromodichloromethane	µg/kg	1 U	--	--	--	--	--	--	NE
Bromoform (Tribromomethane)	µg/kg	1 U	--	--	--	--	--	--	NE
Bromomethane	µg/kg	1 U	--	--	--	--	--	--	NE
Carbon Disulfide	µg/kg	1 U	--	--	--	--	--	--	NE
Carbon Tetrachloride	µg/kg	1 U	--	--	--	--	--	--	NE
Chlorobenzene	µg/kg	1 U	--	--	--	--	--	--	NE
Chloroethane	µg/kg	5.1 U	--	--	--	--	--	--	NE
Chloroform	µg/kg	1 U	--	--	--	--	--	--	NE
Chloromethane	µg/kg	5.1 U	--	--	--	--	--	--	NE
Cis-1,2-Dichloroethene	µg/kg	1 U	--	--	--	--	--	--	NE
Cis-1,3-Dichloropropene	µg/kg	1 U	--	--	--	--	--	--	NE
Dibromochloromethane	µg/kg	1 U	--	--	--	--	--	--	NE
Dibromomethane	µg/kg	1 U	--	--	--	--	--	--	NE
Dichlorodifluoromethane (CFC-12)	µg/kg	1 U	--	--	--	--	--	--	NE
Ethylbenzene	µg/kg	1 U	--	--	--	--	--	--	6,000
Hexachlorobutadiene	µg/kg	5.1 U	--	--	--	--	--	--	NE
Isopropylbenzene (Cumene)	µg/kg	1 U	--	--	--	--	--	--	NE
Methyl Iodide (Iodomethane)	µg/kg	5.1 U	--	--	--	--	--	--	NE
Methyl t-butyl ether	µg/kg	1 U	--	--	--	--	--	--	100
Methylene Chloride	µg/kg	5.1 U	--	--	--	--	--	--	20
Naphthalene	µg/kg	1 U	--	--	--	--	--	--	5,000
n-Butylbenzene	µg/kg	1 U	--	--	--	--	--	--	NE
n-Propylbenzene	µg/kg	1 U	--	--	--	--	--	--	NE
p-Isopropyltoluene	µg/kg	1 U	--	--	--	--	--	--	NE
Sec-Butylbenzene	µg/kg	1 U	--	--	--	--	--	--	NE
Styrene	µg/kg	1 U	--	--	--	--	--	--	NE
Tert-Butylbenzene	µg/kg	1 U	--	--	--	--	--	--	NE
Tetrachloroethene	µg/kg	1 U	--	--	--	--	--	--	50
Toluene	µg/kg	5.1 U	--	--	--	--	--	--	7,000
Trans-1,2-Dichloroethene	µg/kg	1 U	--	--	--	--	--	--	NE
Trans-1,3-Dichloropropene	µg/kg	1 U	--	--	--	--	--	--	NE
Trichloroethene (TCE)	µg/kg	1 U	--	--	--	--	--	--	30

Analytes	Units	Sample ID, Date and Depth (in ft.)							Screening Levels
		DP-20-0.0-2.0 2/08/12 0 - 2	DP-21-0.0-2.0 2/08/12 0 - 2	DP-21-2.0-3.0 2/08/12 2 - 3	HA1-0.0-2.0 2/09/12 0 - 2	HA2-0.0-2.0 2/09/12 0 - 2	HA3-0.0-2.0 2/09/12 0 - 2	HA4-0.0-2.0 2/09/12 0 - 2	MTCA <sup>2</sup> A ULU
Trichlorofluoromethane (CFC-11)	µg/kg	1.6	--	--	--	--	--	--	NE
Vinyl Acetate	µg/kg	5.1 U	--	--	--	--	--	--	NE
Vinyl Chloride	µg/kg	1 U	--	--	--	--	--	--	NE
Xylene, m-,p-	µg/kg	2.0 U	--	--	--	--	--	--	NE
Xylene, o-	µg/kg	1.0 U	--	--	--	--	--	--	NE
Xylene, Total	µg/kg	3.0 U	--	--	--	--	--	--	9,000
<b>SVOCs by SW8270</b>									
1,2-Diphenylhydrazine	µg/kg	--	--	--	--	--	--	--	NE
1,3-Dinitrobenzene	µg/kg	--	--	--	--	--	--	--	NE
2,3,4,6-Tetrachlorophenol	µg/kg	--	--	--	--	--	--	--	NE
2,3,5,6-Tetrachlorophenol	µg/kg	--	--	--	--	--	--	--	NE
2,3-DICHLOROANILINE	µg/kg	--	--	--	--	--	--	--	NE
2,4,5-Trichlorophenol	µg/kg	--	--	--	--	--	--	--	NE
2,4,6-Trichlorophenol	µg/kg	--	--	--	--	--	--	--	NE
2,4-Dichlorophenol	µg/kg	--	--	--	--	--	--	--	NE
2,4-Dimethylphenol	µg/kg	--	--	--	--	--	--	--	NE
2,4-Dinitrophenol	µg/kg	--	--	--	--	--	--	--	NE
2,4-Dinitrotoluene	µg/kg	--	--	--	--	--	--	--	NE
2,6-Dinitrotoluene	µg/kg	--	--	--	--	--	--	--	NE
2-Chloronaphthalene	µg/kg	--	--	--	--	--	--	--	NE
2-Chlorophenol	µg/kg	--	--	--	--	--	--	--	NE
2-Nitroaniline	µg/kg	--	--	--	--	--	--	--	NE
2-Nitrophenol	µg/kg	--	--	--	--	--	--	--	NE
3,3'-Dichlorobenzidine	µg/kg	--	--	--	--	--	--	--	NE
3-Nitroaniline	µg/kg	--	--	--	--	--	--	--	NE
4,6-Dinitro-2-Methylphenol	µg/kg	--	--	--	--	--	--	--	NE
4-Bromophenyl phenyl ether	µg/kg	--	--	--	--	--	--	--	NE
4-Chloro-3-Methylphenol	µg/kg	--	--	--	--	--	--	--	NE
4-Chloroaniline	µg/kg	--	--	--	--	--	--	--	NE
4-Chlorophenyl-Phenylether	µg/kg	--	--	--	--	--	--	--	NE
4-Nitroaniline	µg/kg	--	--	--	--	--	--	--	NE
4-Nitrophenol (p-Nitrophenol)	µg/kg	--	--	--	--	--	--	--	NE
Aniline	µg/kg	--	--	--	--	--	--	--	NE
Benzene, 1,4-Dinitro-	µg/kg	--	--	--	--	--	--	--	NE
Benzidine	µg/kg	--	--	--	--	--	--	--	NE
Benzyl Alcohol	µg/kg	--	--	--	--	--	--	--	NE
Bis(2-Chloroethoxy)Methane	µg/kg	--	--	--	--	--	--	--	NE
Bis(2-Chloroethyl)Ether	µg/kg	--	--	--	--	--	--	--	NE

Analytes	Units	Sample ID, Date and Depth (in ft.)							Screening Levels
		DP-20-0.0-2.0 2/08/12 0 - 2	DP-21-0.0-2.0 2/08/12 0 - 2	DP-21-2.0-3.0 2/08/12 2 - 3	HA1-0.0-2.0 2/09/12 0 - 2	HA2-0.0-2.0 2/09/12 0 - 2	HA3-0.0-2.0 2/09/12 0 - 2	HA4-0.0-2.0 2/09/12 0 - 2	MTCA <sup>2</sup> A ULU
Bis(2-chloroisopropyl) ether	µg/kg	--	--	--	--	--	--	--	NE
Bis(2-Ethylhexyl) Phthalate	µg/kg	--	--	--	--	--	--	--	NE
Butyl benzyl phthalate	µg/kg	--	--	--	--	--	--	--	NE
Carbazole	µg/kg	--	--	--	--	--	--	--	NE
Dibutyl phthalate	µg/kg	--	--	--	--	--	--	--	NE
Diethyl phthalate	µg/kg	--	--	--	--	--	--	--	NE
Dimethyl phthalate	µg/kg	--	--	--	--	--	--	--	NE
Di-N-Octyl Phthalate	µg/kg	--	--	--	--	--	--	--	NE
Hexachlorobenzene	µg/kg	--	--	--	--	--	--	--	NE
Hexachlorocyclopentadiene	µg/kg	--	--	--	--	--	--	--	NE
Hexachloroethane	µg/kg	--	--	--	--	--	--	--	NE
Hexanedioic Acid, Bis(2-Ethylhexyl) Ester	µg/kg	--	--	--	--	--	--	--	NE
Isophorone	µg/kg	--	--	--	--	--	--	--	NE
m,p-Cresol	µg/kg	--	--	--	--	--	--	--	NE
Naphthalene	µg/kg	--	--	--	--	--	--	--	<b>5,000</b>
Nitrobenzene	µg/kg	--	--	--	--	--	--	--	NE
N-Nitrosodimethylamine	µg/kg	--	--	--	--	--	--	--	NE
N-Nitrosodi-n-propylamine	µg/kg	--	--	--	--	--	--	--	NE
N-Nitrosodiphenylamine	µg/kg	--	--	--	--	--	--	--	NE
o-Cresol (2-methylphenol)	µg/kg	--	--	--	--	--	--	--	NE
O-DINITROBENZENE	µg/kg	--	--	--	--	--	--	--	NE
Pentachlorophenol	µg/kg	--	--	--	--	--	--	--	NE
Phenol	µg/kg	--	--	--	--	--	--	--	NE
Pyridine	µg/kg	--	--	--	--	--	--	--	NE

**Notes:**

<sup>1</sup> Chemical analysis performed by OnSite Environmental Inc. of Redmond, Washington.

<sup>2</sup> Model Toxics Control Act (MTCA) Cleanup Regulation Chapter 173-340 WAC. Method A Unrestricted Land Use (ULU) clean-up levels.

<sup>3</sup> Sample holding time was exceeded by one day for follow up analysis.

<sup>4</sup> The published natural background concentration for chromium in the Puget Sound region is 48 mg/kg (*Natural Background Soil Metals Concentrations in Washington State*. Publication #94-115. October 1994.)

<sup>5</sup> cPAH testing and regulatory evaluation is completed for individual carcinogenic compounds as well as the for the summation of the mixture of the seven carcinogenic PAHs (known as Ecology's toxicity equivalency methodology). The summation procedure is completed using toxicity equivalency factors for each individual compound which are then added to produce a total toxicity equivalent concentration (TEEC) which is then compared to the MTCA cleanup level of 0.1 mg/kg (or 100 µg/kg).

PCBs = Polychlorinated biphenyls

PAHs = Polycyclic aromatic hydrocarbons

VOCs = Volatile organic compound

SVOCs = Semivolatile organic compound

µg/kg = microgram per kilogram

mg/kg = milligrams per kilogram

-- = Not analyzed

U = The analyte was not detected at a concentration greater than the given RL or MDL.

E = The concentration reported exceeds the quantitation range and is an estimate.

H = The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and may be impacting the sample result

J = The value reported was below the practical quantitation limit. The value is an estimate.

NE = A cleanup level has not been established by Ecology.

**Bold Type** Exceeds MTCA A (Soil - Unrestricted Land Use Cleanup Level)

**TABLE 3**  
**CHEMICAL ANALYTICAL DATA - GROUNDWATER<sup>1</sup>**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Analytes	Units	Sample ID and Date				Screening Levels
		DP1-W 12/13/11	DP12-W 2/08/12	DP15-W 2/08/12	WELL1-W 12/14/11	MTCA <sup>2</sup> A
<b>Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx</b>						
Gasoline-range hydrocarbons	mg/l	0.10 U	0.10 U	0.10 U	0.10 U	1.0
Diesel-range hydrocarbons	mg/l	0.27 U	0.27 U	0.30 U	0.26 U	0.5
Lube Oil-range Hydrocarbons	mg/l	0.43 U	0.43 U	0.48 U	0.41 U	0.5
<b>Metals by EPA200.8/7470A (Total)</b>						
Arsenic	µg/l	18	12	27	3.3 U	5.0
Barium	µg/l	30	300	400	28 U	NE
Cadmium	µg/l	4.4 U	4.4 U	4.4 U	4.4 U	5.0
Chromium	µg/l	26	140	180	11 U	50
Lead	µg/l	1.1 U	12	16	1.1 U	15
Mercury	µg/l	0.5 U	0.5 U	0.5 U	0.5 U	2.0
Selenium	µg/l	5.6 U	5.6 U	6.8	5.6 U	NE
Silver	µg/l	11 U	11 U	11 U	11 U	NE
<b>Metals by EPA200.8/7470A (Dissolved)</b>						
Arsenic	µg/l	21	3.0 U	3.0 U	3.0 U	5.0
Barium	µg/l	25 U	25 U	25 U	25 U	NE
Cadmium	µg/l	4.0 U	4.0 U	4.0 U	4.0 U	5.0
Chromium	µg/l	22	10 U	10 U	10 U	50
Lead	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	15
Mercury	µg/l	0.5 U	0.5 U	0.5 U	0.5 U	2.0
Selenium	µg/l	5.0 U	5.0 U	5.0 U	5.0 U	NE
Silver	µg/l	10 U	10 U	10 U	10 U	NE
<b>PCBs by SW8082</b>						
PCB-aroclor 1016	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
PCB-aroclor 1221	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
PCB-aroclor 1232	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
PCB-aroclor 1242	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
PCB-aroclor 1248	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
PCB-aroclor 1254	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
PCB-aroclor 1260	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
Total Aroclors	µg/l	0.048 U	0.048 U	0.1 U	0.048 U	NE
<b>PAHs by SW8270</b>						
1-Methylnaphthalene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
2-Methylnaphthalene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE

Analytes	Units	Sample ID and Date				Screening Levels
		DP1-W 12/13/11	DP12-W 2/08/12	DP15-W 2/08/12	WELL1-W 12/14/11	MTCA <sup>2</sup> A
Acenaphthene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
Acenaphthylene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
Anthracene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
Benzo(a)anthracene	µg/l	0.0097	0.0098 U	0.0092 U	0.0096 U	NE
Benzo(a)pyrene	µg/l	0.0095 U	0.0098 U	0.0092 U	0.0096 U	0.1
Benzo(b)fluoranthene	µg/l	0.0095 U	0.0098 U	0.0092 U	0.0096 U	NE
Benzo(ghi)perylene	µg/l	0.0095 U	0.0098 U	0.0092 U	0.0096 U	NE
Benzo(j,k)fluoranthene	µg/l	0.0095 U	0.0098 U	0.0092 U	0.0096 U	NE
Chrysene	µg/l	0.0095 U	0.0098 U	0.0092 U	0.0096 U	NE
Dibenzo(a,h)anthracene	µg/l	0.0095 U	0.0098 U	0.0092 U	0.0096 U	NE
Dibenzofuran	µg/l	0.95 U	0.98 U	-	0.96 U	NE
Fluoranthene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
Fluorene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
Indeno(1,2,3-cd)pyrene	µg/l	0.0095 U	0.0098 U	0.0092 U	0.0096 U	NE
Naphthalene	µg/l	0.095 U	0.098 U	0.1	0.19	160
Phenanthrene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
Pyrene	µg/l	0.095 U	0.098 U	0.092 U	0.096 U	NE
Total cPAH <sup>3</sup> TTEC (ND=0.5RL)	µg/l	0.0077	0.0074 U	0.0069 U	0.0072 U	0.1
<b>VOCs by SW8260</b>						
1,1,1,2-Tetrachloroethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,1,1-Trichloroethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	200
1,1,2,2-Tetrachloroethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,1,2-Trichloroethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,1-Dichloroethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,1-Dichloroethene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,1-Dichloropropene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,2,3-Trichlorobenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,2,3-Trichloropropane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,2,4-Trichlorobenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,2,4-Trimethylbenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,2-Dibromo-3-Chloropropane	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	NE
1,2-dibromoethane (EDB)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	0.01
1,2-Dichlorobenzene (o-Dichlorobenzene)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,2-Dichloroethane (EDC)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	5.0
1,2-Dichloropropane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,3,5-Trimethylbenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,3-Dichlorobenzene (m-Dichlorobenzene)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
1,3-Dichloropropane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE



Analytes	Units	Sample ID and Date				Screening Levels
		DP1-W 12/13/11	DP12-W 2/08/12	DP15-W 2/08/12	WELL1-W 12/14/11	MTCA <sup>2</sup> A
1,4-Dichlorobenzene (p-Dichlorobenzene)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
2,2-Dichloropropane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
2-Butanone (MEK)	µg/l	5.0 U	5.0 U	5.0 U	5.0 U	NE
2-Chloroethyl vinyl ether	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	NE
2-Chlorotoluene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
2-Hexanone	µg/l	2.0 U	2.0 U	2.0 U	2.0 U	NE
4-Chlorotoluene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
4-Methyl-2-Pentanone (Methyl isobutyl ketone)	µg/l	2.0 U	2.0 U	2.0 U	2.0 U	NE
Acetone	µg/l	5.0 U	5.0 U	5.0 U	5.0 U	NE
Benzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	5.0
Bromobenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Bromochloromethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Bromodichloromethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Bromoform (Tribromomethane)	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	NE
Bromomethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Carbon Disulfide	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Carbon Tetrachloride	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Chlorobenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Chloroethane	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	NE
Chloroform	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Chloromethane	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	NE
Cis-1,2-Dichloroethene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Cis-1,3-Dichloropropene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Dibromochloromethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Dibromomethane	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Dichlorodifluoromethane (CFC-12)	µg/l	0.52	0.2 U	0.2 U	0.2 U	NE
Ethylbenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	700
Hexachlorobutadiene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Isopropylbenzene (Cumene)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Methyl Iodide (Iodomethane)	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	NE
Methyl t-butyl ether	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	20
Methylene Chloride	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	5.0
Naphthalene	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	160
n-Butylbenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
n-Propylbenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
p-Isopropyltoluene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Sec-Butylbenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Styrene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Tert-Butylbenzene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE

Analytes	Units	Sample ID and Date				Screening Levels
		DP1-W 12/13/11	DP12-W 2/08/12	DP15-W 2/08/12	WELL1-W 12/14/11	MTCA <sup>2</sup> A
Tetrachloroethene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	5.0
Toluene	µg/l	1.0 U	1.0 U	1.0 U	1.0 U	1,000
Trans-1,2-Dichloroethene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Trans-1,3-Dichloropropene	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Trichloroethene (TCE)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	5.0
Trichlorofluoromethane (CFC-11)	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Vinyl Acetate	µg/l	2.0 U	2.0 U	2.0 U	2.0 U	NE
Vinyl Chloride	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	0.2
Xylene, m-,p-	µg/l	0.4 U	0.4 U	0.4 U	0.4 U	NE
Xylene, o-	µg/l	0.2 U	0.2 U	0.2 U	0.2 U	NE
Xylene, Total	µg/l	0.6 U	0.6 U	0.6 U	0.6 U	1,000
<b>SVOCs by SW8270</b>						
1,2-Diphenylhydrazine	µg/l	0.95 U	0.98 U	--	0.96 U	NE
1,3-Dinitrobenzene	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,3,4,6-Tetrachlorophenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,3,5,6-Tetrachlorophenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,3-DICHLOROANILINE	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,4,5-Trichlorophenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,4,6-Trichlorophenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,4-Dichlorophenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,4-Dimethylphenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,4-Dinitrophenol	µg/l	4.8 U	4.9 U	--	4.8 U	NE
2,4-Dinitrotoluene	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2,6-Dinitrotoluene	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2-Chloronaphthalene	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2-Chlorophenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2-Nitroaniline	µg/l	0.95 U	0.98 U	--	0.96 U	NE
2-Nitrophenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
3,3'-Dichlorobenzidine	µg/l	0.95 U	0.98 U	--	0.96 U	NE
3-Nitroaniline	µg/l	0.95 U	0.98 U	--	0.96 U	NE
4,6-Dinitro-2-Methylphenol	µg/l	4.8 U	4.9 U	--	4.8 U	NE
4-Bromophenyl phenyl ether	µg/l	0.95 U	0.98 U	--	0.96 U	NE
4-Chloro-3-Methylphenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
4-Chloroaniline	µg/l	0.95 U	0.98 U	--	0.96 U	NE
4-Chlorophenyl-Phenylether	µg/l	0.95 U	0.98 U	--	0.96 U	NE
4-Nitroaniline	µg/l	0.95 U	0.98 U	--	0.96 U	NE
4-Nitrophenol (p-Nitrophenol)	µg/l	0.95 U	0.98 U	--	0.96 U	NE

Analytes	Units	Sample ID and Date				Screening Levels
		DP1-W 12/13/11	DP12-W 2/08/12	DP15-W 2/08/12	WELL1-W 12/14/11	MTCA <sup>2</sup> A
Aniline	µg/l	4.8 U	4.9 U	--	4.8 U	NE
Benzene, 1,4-Dinitro-	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Benzidine	µg/l	4.8 U	4.9 U	--	4.8 U	NE
Benzyl Alcohol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Bis(2-Chloroethoxy)Methane	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Bis(2-Chloroethyl)Ether	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Bis(2-chloroisopropyl) ether	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Bis(2-Ethylhexyl) Phthalate	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Butyl benzyl phthalate	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Carbazole	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Dibutyl phthalate	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Diethyl phthalate	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Dimethyl phthalate	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Di-N-Octyl Phthalate	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Hexachlorobenzene	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Hexachlorocyclopentadiene	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Hexachloroethane	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Hexanedioic Acid, Bis(2-Ethylhexyl) Ester	µg/l	4.8 U	4.9 U	--	4.8 U	NE
Isophorone	µg/l	0.95 U	0.98 U	--	0.96 U	NE
m,p-Cresol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Naphthalene	µg/l	0.095 U	0.098 U	--	0.19	<b>160</b>
Nitrobenzene	µg/l	0.95 U	0.98 U	--	0.96 U	NE
N-Nitrosodimethylamine	µg/l	0.95 U	0.98 U	--	0.96 U	NE
N-Nitrosodi-n-propylamine	µg/l	0.95 U	0.98 U	--	0.96 U	NE
N-Nitrosodiphenylamine	µg/l	0.95 U	0.98 U	--	0.96 U	NE
o-Cresol (2-methylphenol)	µg/l	0.95 U	0.98 U	--	0.96 U	NE
O-DINITROBENZENE	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Pentachlorophenol	µg/l	4.8 U	4.9 U	--	4.8 U	NE
Phenol	µg/l	0.95 U	0.98 U	--	0.96 U	NE
Pyridine	µg/l	0.95 U	0.98 U	--	0.96 U	NE

**Notes:**

<sup>1</sup> Chemical analysis performed by OnSite Environmental Inc. of Redmond, Washington.

<sup>2</sup> Model Toxics Control Act (MTCA) Cleanup Regulation Chapter 173-340 WAC. Method A clean-up levels.

<sup>3</sup> cPAH testing and regulatory evaluation is completed for individual carcinogenic compounds as well as the for the summation of the mixture of the seven carcinogenic PAHs (known as Ecology's toxicity equivalency methodology). The summation procedure is completed using toxicity equivalency factors for each individual compound which are then added to produce a total toxicity equivalent concentration (TTEC) which is then compared to the MTCA cleanup level of 0.1 µg/l.

PCBs = Polychlorinated biphenyls

PAHs = Polycyclic aromatic hydrocarbons

VOCs = Volatile organic compound

SVOCs = Semivolatile organic compound

E = The concentration reported exceeds the quantitation range and is an estimate.

H = The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and may be impacting the sample result

J = The value reported was below the practical quantitation limit. The value is an estimate.

NE = A cleanup level has not been established by Ecology.

**Bold Type** Exceeds MTCA A (Groundwater - Table Value)

µg/l = microgram per liter

mg/l = milligrams per liter

-- = Not analyzed

U = The analyte was not detected at a concentration greater than the given RL or MDL.

**TABLE 4**  
**CHEMICAL ANALYTICAL DATA - SEDIMENT<sup>1</sup>**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Analytes	Units	Sample ID and Date								Screening Levels
		SED1 12/13/11	SED3 12/13/11	SED4 12/13/11	SED5 12/13/11	SED6 12/13/11	SED7 12/13/11	SED8 12/13/11	SED9 2/09/12	MTCA <sup>2</sup> A ULU
<b>Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx</b>										
Gasoline-range hydrocarbons	mg/kg	7.5 U	9.0 U	9.6 U	7.6 U	9.6 U	5.3 U	8.0 U	-	100
Diesel-range hydrocarbons	mg/kg	35 U	38 U	40 U	120 U	37 U	28 U	33 U	130 U	2,000
Lube Oil-range Hydrocarbons	mg/kg	70 U	77 U	270	950	75 U	57 U	65 U	1,000	2,000
<b>Metals by EPA6010B/7471A/SW7196A</b>										
Arsenic	mg/kg	14 U	15 U	16 U	13 U	15 U	11 U	13 U	-	20
Barium	mg/kg	85	64	88	75	100	41	78	-	NE
Cadmium	mg/kg	0.70 U	0.76 U	0.79 U	0.89	0.9	0.57 U	0.65 U	-	2.0
Chromium <sup>3</sup>	mg/kg	36	41	48	57	37	30	29	-	NE
Lead	mg/kg	17	7.6 U	39	120	13	5.7 U	6.5 U	-	250
Mercury	mg/kg	0.35 U	0.38 U	0.40 U	0.33 U	0.37 U	0.28 U	0.33 U	-	2.0
Selenium	mg/kg	14 U	15 U	16 U	13 U	15 U	11 U	13 U	-	NE
Silver	mg/kg	0.70 U	0.76 U	0.79 U	0.66 U	0.75 U	0.57 U	0.65 U	-	NE
<b>PCBs by SW8082</b>										
PCB-aroclor 1016	µg/kg	70 U	76 U	79 U	66 U	75 U	57 U	65 U	-	NE
PCB-aroclor 1221	µg/kg	70 U	76 U	79 U	67 U	75 U	57 U	65 U	-	NE
PCB-aroclor 1232	µg/kg	70 U	76 U	79 U	68 U	75 U	57 U	65 U	-	NE
PCB-aroclor 1242	µg/kg	70 U	76 U	79 U	69 U	75 U	57 U	65 U	-	NE
PCB-aroclor 1248	µg/kg	70 U	76 U	79 U	70 U	75 U	57 U	65 U	-	NE
PCB-aroclor 1254	µg/kg	130	76 U	79 U	160	75 U	57 U	65 U	-	NE
PCB-aroclor 1260	µg/kg	70 U	76 U	79 U	66 U	75 U	57 U	65 U	-	NE
Total Aroclors	µg/kg	130	76 U	79 U	160	75 U	57 U	65 U	-	1,000
<b>PAHs by SW8270</b>										
1-Methylnaphthalene	µg/kg	14	10 U	11 U	44 U	50 U	7.5 U	8.7 U	17 U	NE
2-Methylnaphthalene	µg/kg	22	10 U	11 U	94	50 U	7.5 U	8.7 U	17 U	NE
Acenaphthene	µg/kg	21	10 U	11 U	44 U	50 U	7.5 U	8.7 U	17 U	NE
Acenaphthylene	µg/kg	9.4 U	10 U	11 U	61	50 U	7.5 U	8.7 U	17 U	NE
Anthracene	µg/kg	9.4 U	10 U	11 U	78	50 U	7.5 U	8.7 U	17 U	NE
Benzo(a)anthracene	µg/kg	11	10 U	11 U	160	50 U	7.5 U	8.7 U	17 U	NE
Benzo(a)pyrene	µg/kg	12	10 U	11 U	270	50 U	7.5 U	8.7 U	17 U	100
Benzo(b)fluoranthene	µg/kg	13	10 U	11 U	330	50 U	7.5 U	8.7 U	17	NE
Benzo(ghi)perylene	µg/kg	9.6	10 U	11 U	230	50 U	7.5 U	8.7 U	17 U	NE
Benzo(j,k)fluoranthene	µg/kg	9.4 U	10 U	11 U	110	50 U	7.5 U	8.7 U	17 U	NE

Analytes	Units	Sample ID and Date								Screening Levels
		SED1 12/13/11	SED3 12/13/11	SED4 12/13/11	SED5 12/13/11	SED6 12/13/11	SED7 12/13/11	SED8 12/13/11	SED9 2/09/12	MTCA <sup>2</sup> A ULU
Chrysene	µg/kg	16	10 U	11 U	200	50 U	7.5 U	8.7 U	17 U	NE
Dibenzo(a,h)anthracene	µg/kg	9.4 U	10 U	11 U	49	50 U	7.5 U	8.7 U	17 U	NE
Dibenzofuran	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Fluoranthene	µg/kg	30	10 U	11 U	300	50 U	7.5 U	8.7 U	17 U	NE
Fluorene	µg/kg	14	10 U	11 U	44 U	50 U	7.5 U	8.7 U	17 U	NE
Indeno(1,2,3-cd)pyrene	µg/kg	9.4 U	10 U	11 U	200	50 U	7.5 U	8.7 U	17 U	NE
Naphthalene	µg/kg	48	10 U	11 U	93	50 U	7.5 U	8.7 U	17 U	5,000
Phenanthrene	µg/kg	59	10 U	11 U	190	50 U	7.5 U	8.7 U	17 U	NE
Pyrene	µg/kg	40	10 U	11 U	290	93	7.5 U	8.7 U	17 U	NE
Total cPAH <sup>4</sup> TTEC (ND=0.5RL)	µg/kg	15.97	7.55 U	8.31 U	356.9	37.8 U	5.66 U	6.57 U	13.69	100
<b>VOCs by SW8260</b>										
1,1,1,2-Tetrachloroethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,1,1-Trichloroethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	2,000
1,1,2,2-Tetrachloroethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,1,2-Trichloroethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,1-Dichloroethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,1-Dichloroethene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,1-Dichloropropene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,2,3-Trichlorobenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,2,3-Trichloropropane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,2,4-Trichlorobenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,2,4-Trimethylbenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,2-Dibromo-3-Chloropropane	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
1,2-dibromoethane (EDB)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	5.0
1,2-Dichlorobenzene (o-Dichlorobenzene)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,2-Dichloroethane (EDC)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,2-Dichloropropane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,3,5-Trimethylbenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,3-Dichlorobenzene (m-Dichlorobenzene)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,3-Dichloropropane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
1,4-Dichlorobenzene (p-Dichlorobenzene)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
2,2-Dichloropropane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
2-Butanone (MEK)	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
2-Chloroethyl vinyl ether	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
2-Chlorotoluene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
2-Hexanone	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
4-Chlorotoluene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
4-Methyl-2-Pentanone (Methyl isobutyl ketone)	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE

Analytes	Units	Sample ID and Date								Screening Levels
		SED1 12/13/11	SED3 12/13/11	SED4 12/13/11	SED5 12/13/11	SED6 12/13/11	SED7 12/13/11	SED8 12/13/11	SED9 2/09/12	MTCA <sup>2</sup> A ULU
Acetone	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
Benzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	30
Bromobenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Bromochloromethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Bromodichloromethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Bromoform (Tribromomethane)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Bromomethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Carbon Disulfide	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Carbon Tetrachloride	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Chlorobenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Chloroethane	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
Chloroform	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Chloromethane	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
Cis-1,2-Dichloroethene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Cis-1,3-Dichloropropene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Dibromochloromethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Dibromomethane	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Dichlorodifluoromethane (CFC-12)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Ethylbenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	6,000
Hexachlorobutadiene	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
Isopropylbenzene (Cumene)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Methyl Iodide (Iodomethane)	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE
Methyl t-butyl ether	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	100
Methylene Chloride	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	20
Naphthalene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	5,000
n-Butylbenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
n-Propylbenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
p-Isopropyltoluene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Sec-Butylbenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Styrene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Tert-Butylbenzene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Tetrachloroethene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	50
Toluene	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	7,000
Trans-1,2-Dichloroethene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Trans-1,3-Dichloropropene	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Trichloroethene (TCE)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	30
Trichlorofluoromethane (CFC-11)	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Vinyl Acetate	µg/kg	6.9 U	6.4 U	8.9 U	6.8 U	6.9 U	4.9 U	5.9 U	--	NE

Analytes	Units	Sample ID and Date								Screening Levels
		SED1 12/13/11	SED3 12/13/11	SED4 12/13/11	SED5 12/13/11	SED6 12/13/11	SED7 12/13/11	SED8 12/13/11	SED9 2/09/12	MTCA <sup>2</sup> A ULU
Vinyl Chloride	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Xylene, m-,p-	µg/kg	2.8 U	2.6 U	3.6 U	2.7 U	2.7 U	1.9 U	2.4 U	--	NE
Xylene, o-	µg/kg	1.4 U	1.3 U	1.8 U	1.4 U	1.4 U	0.97 U	1.2 U	--	NE
Xylene, Total	µg/kg	4.2 U	3.9 U	5.4 U	4.1 U	4.1 U	2.9 U	3.6 U	--	9,000
<b>SVOCs by SW8270</b>										
1,2-Diphenylhydrazine	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
1,3-Dinitrobenzene	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,3,4,6-Tetrachlorophenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,3,5,6-Tetrachlorophenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,3-DICHLOROANILINE	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,4,5-Trichlorophenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,4,6-Trichlorophenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,4-Dichlorophenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,4-Dimethylphenol	µg/kg	--	--	--	2200 U	500 U	--	--	--	NE
2,4-Dinitrophenol	µg/kg	--	--	--	1100 U	250 U	--	--	--	NE
2,4-Dinitrotoluene	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2,6-Dinitrotoluene	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2-Chloronaphthalene	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2-Chlorophenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2-Nitroaniline	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
2-Nitrophenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
3,3'-Dichlorobenzidine	µg/kg	--	--	--	2200 U	500 U	--	--	--	NE
3-Nitroaniline	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
4,6-Dinitro-2-Methylphenol	µg/kg	--	--	--	1100 U	250 U	--	--	--	NE
4-Bromophenyl phenyl ether	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
4-Chloro-3-Methylphenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
4-Chloroaniline	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
4-Chlorophenyl-Phenylether	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
4-Nitroaniline	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
4-Nitrophenol (p-Nitrophenol)	µg/kg	--	--	--	220 U	130	--	--	--	NE
Aniline	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Benzene, 1,4-Dinitro-	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Benzidine	µg/kg	--	--	--	2200 U	500 U	--	--	--	NE
Benzyl Alcohol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Bis(2-Chloroethoxy)Methane	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Bis(2-Chloroethyl)Ether	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Bis(2-chloroisopropyl) ether	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Bis(2-Ethylhexyl) Phthalate	µg/kg	--	--	--	1100 U	250 U	--	--	--	NE

Analytes	Units	Sample ID and Date								Screening Levels
		SED1 12/13/11	SED3 12/13/11	SED4 12/13/11	SED5 12/13/11	SED6 12/13/11	SED7 12/13/11	SED8 12/13/11	SED9 2/09/12	MTCA <sup>2</sup> A ULU
Butyl benzyl phthalate	µg/kg	--	--	--	2200 U	500 U	--	--	--	NE
Carbazole	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Dibutyl phthalate	µg/kg	--	--	--	2200 U	500 U	--	--	--	NE
Diethyl phthalate	µg/kg	--	--	--	1100 U	250 U	--	--	--	NE
Dimethyl phthalate	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Di-N-Octyl Phthalate	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Hexachlorobenzene	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Hexachlorocyclopentadiene	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Hexachloroethane	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Hexanedioic Acid, Bis(2-Ethylhexyl) Ester	µg/kg	--	--	--	1100 U	250 U	--	--	--	NE
Isophorone	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
m,p-Cresol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Naphthalene	µg/kg	48	10 U	11 U	93	50 U	7.5 U	8.7 U	17 U	<b>5,000</b>
Nitrobenzene	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
N-Nitrosodimethylamine	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
N-Nitrosodi-n-propylamine	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
N-Nitrosodiphenylamine	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
o-Cresol (2-methylphenol)	µg/kg	--	--	--	220 U	64	--	--	--	NE
O-DINITROBENZENE	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Pentachlorophenol	µg/kg	--	--	--	1100 U	250 U	--	--	--	NE
Phenol	µg/kg	--	--	--	220 U	50 U	--	--	--	NE
Pyridine	µg/kg	--	--	--	2200 U	500 U	--	--	--	NE

**Notes:**

<sup>1</sup> Chemical analysis performed by OnSite Environmental Inc. of Redmond, Washington.

<sup>2</sup> Model Toxics Control Act (MTCA) Cleanup Regulation Chapter 173-340 WAC. Method A Unrestricted Land Use (ULU) clean-up levels.

<sup>3</sup> The published natural background concentration for chromium in the Puget Sound region is 48 mg/kg (*Natural Background Soil Metals Concentrations in Washington State*. Publication #94-115. October 1994.)

<sup>4</sup> cPAH testing and regulatory evaluation is completed for individual carcinogenic compounds as well as the for the summation of the mixture of the seven carcinogenic PAHs (known as Ecology's toxicity equivalency methodology). The summation procedure is completed using toxicity equivalency factors for each individual compound which are then added to produce a total toxicity equivalent concentration (TTEC) which is then compared to the MTCA cleanup level of 0.1 mg/kg (or 100 µg/kg).

PCBs = Polychlorinated biphenyls  
PAHs = Polycyclic aromatic hydrocarbons  
VOCs = Volatile organic compound  
SVOCs = Semivolatile organic compound  
µg/kg = microgram per kilogram  
mg/kg = milligrams per kilogram  
-- = Not analyzed  
U = The analyte was not detected at a concentration greater than the given RL or MDL.  
E = The concentration reported exceeds the quantitation range and is an estimate.  
H = The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and may be impacting the sample result  
J = The value reported was below the practical quantitation limit. The value is an estimate.  
NE = A cleanup level has not been established by Ecology.

**Bold Type** Exceeds MTCA A (Soil - Unrestricted Land Use Cleanup Level)



**TABLE 5**  
**CHEMICAL ANALYTICAL DATA - SURFACE WATER<sup>1</sup>**  
**WSDOT MIDWAY METALS**  
**SEQUIM, WASHINGTON**

Analytes	Units	Sample ID and Date			Screening Levels			
		SW1 12/13/2011	SW2 12/13/2011	SW9 2/08/2012	Surface Water MTCA B - Carcinogenic Table Value	Surface Water ARAR - Aquatic Life Fresh/Chronic - Ch. 173-201A WAC	Surface Water ARAR - Aquatic Life Fresh/Chronic - Clean Water Act §304	Surface Water ARAR - Aquatic Life Fresh/Chronic - National Toxics Rule, 40 CFR 131
<b>Total Petroleum Hydrocarbons by NWTPH-Gx and NWTPH-Dx</b>								
Gasoline-range hydrocarbons	mg/l	0.1 U	0.1 U	0.1 U	NE	NE	NE	NE
Diesel-range hydrocarbons	mg/l	0.29 U	0.29 U	0.27 U	NE	NE	NE	NE
Lube Oil-range Hydrocarbons	mg/l	0.47 U	0.46 U	0.43 U	NE	NE	NE	NE
<b>Metals by EPA200.8/7470A (Total)</b>								
Arsenic	µg/l	3.3 U	3.3 U	3.3 U	0.058	190	150	190
Barium	µg/l	41	31	28 U	3,200	NE	NE	NE
Cadmium	µg/l	4.4 U	4.4 U	4.4 U	16	0.37	0.25	1.0
Chromium	µg/l	11 U	11 U	11 U	NE	NE	NE	NE
Lead	µg/l	7.4	4.0	8.3	NE	0.54	2.5	2.5
Mercury	µg/l	0.5 U	0.5 U	0.5 U	NE	0.012	0.77	0.012
Selenium	µg/l	5.6 U	5.6 U	5.6 U	80	5.0	5.0	5.0
Silver	µg/l	11 U	11 U	11 U	80	NE	NE	NE
<b>Metals by EPA200.8/7470A (Dissolved)</b>								
Arsenic	µg/l	3.0 U	3.0 U	3.0 U	0.058	190	150	190
Barium	µg/l	31	25 U	25 U	3,200	NE	NE	NE
Cadmium	µg/l	4.0 U	4.0 U	4.0 U	16	0.37	0.25	1.0
Chromium	µg/l	10 U	10 U	10 U	NE	NE	NE	NE
Lead	µg/l	1.4	1.4	1.6	NE	0.54	2.5	2.5
Mercury	µg/l	0.5 U	0.5 U	0.5 U	NE	0.012	0.77	0.012
Selenium	µg/l	5.0 U	5.0 U	5.0 U	80	5.0	5.0	5.0
Silver	µg/l	10 U	10 U	10 U	80	NE	NE	NE
<b>PCBs by SW8082</b>								
PCB-aroclor 1016	µg/l	0.049 U	0.048 U	0.053 U	1.1	NE	NE	0.014
PCB-aroclor 1221	µg/l	0.049 U	0.048 U	0.053 U	NE	NE	NE	NE
PCB-aroclor 1232	µg/l	0.049 U	0.048 U	0.053 U	NE	NE	NE	NE
PCB-aroclor 1242	µg/l	0.074	0.048 U	0.053 U	NE	NE	NE	NE
PCB-aroclor 1248	µg/l	0.049 U	0.048 U	0.053 U	NE	NE	NE	NE
PCB-aroclor 1254	µg/l	0.049 U	0.048 U	0.053 U	0.32	NE	NE	0.014
PCB-aroclor 1260	µg/l	0.049 U	0.048 U	0.053 U	NE	NE	NE	0.014
Total Aroclors	µg/l	0.074 T	0.048 UT	0.053 UT	NE	0.014	0.014	0.14

Analytes	Units	Sample ID and Date			Screening Levels			
		SW1 12/13/2011	SW2 12/13/2011	SW9 2/08/2012	Surface Water MTCA B - Carcinogenic Table Value	Surface Water ARAR - Aquatic Life Fresh/Chronic - Ch. 173-201A WAC	Surface Water ARAR - Aquatic Life Fresh/Chronic - Clean Water Act §304	Surface Water ARAR - Aquatic Life Fresh/Chronic - National Toxics Rule, 40 CFR 131
<b>PAHs by SW8270</b>								
1-Methylnaphthalene	µg/l	0.098 U	0.095 U	0.1 U	2.4	NE	NE	NE
2-Methylnaphthalene	µg/l	0.098 U	0.095 U	0.1 U	32	NE	NE	NE
Acenaphthene	µg/l	0.098 U	0.095 U	0.1 U	960	NE	NE	NE
Acenaphthylene	µg/l	0.098 U	0.095 U	0.1 U	NE	NE	NE	NE
Anthracene	µg/l	0.098 U	0.095 U	0.1 U	4,800	NE	NE	NE
Benzo(a)anthracene	µg/l	0.01	0.0095 U	0.01 U	NE	NE	NE	NE
Benzo(a)pyrene	µg/l	0.0098 U	0.0095 U	0.01 U	0.012	NE	NE	NE
Benzo(b)fluoranthene	µg/l	0.0098 U	0.0095 U	0.01 U	NE	NE	NE	NE
Benzo(ghi)perylene	µg/l	0.0098 U	0.0095 U	0.01 U	NE	NE	NE	NE
Benzo(j,k)fluoranthene	µg/l	0.0098 U	0.0095 U	0.01 U	NE	NE	NE	NE
Chrysene	µg/l	0.0098 U	0.0095 U	0.01 U	NE	NE	NE	NE
Dibenzo(a,h)anthracene	µg/l	0.0098 U	0.0095 U	0.01 U	NE	NE	NE	NE
Dibenzofuran	µg/l	-	0.95 U	-	16	NE	NE	NE
Fluoranthene	µg/l	0.098 U	0.095 U	0.1 U	640	NE	NE	NE
Fluorene	µg/l	0.098 U	0.095 U	0.1 U	640	NE	NE	NE
Indeno(1,2,3-cd)pyrene	µg/l	0.0098 U	0.0095 U	0.01 U	NE	NE	NE	NE
Naphthalene	µg/l	0.12	0.095 U	0.1 U	160	NE	NE	NE
Phenanthrene	µg/l	0.098 U	0.095 U	0.1 U	NE	NE	NE	NE
Pyrene	µg/l	0.098 U	0.095 U	0.1 U	480	NE	NE	NE
Total cPAH TEQ (ND=0.5RL)	µg/l	0.007909 T	0.007172 UT	0.00755 UT	NE	NE	NE	NE
<b>VOCs by SW8260</b>								
1,1,1,2-Tetrachloroethane	µg/l	0.2 U	0.2 U	0.2 U	1.7	NE	NE	NE
1,1,1-Trichloroethane	µg/l	0.2 U	0.2 U	0.2 U	7,200	NE	NE	NE
1,1,1,2-Tetrachloroethane	µg/l	0.2 U	0.2 U	0.2 U	0.22	NE	NE	NE
1,1,2-Trichloroethane	µg/l	0.2 U	0.2 U	0.2 U	0.77	NE	NE	NE
1,1-Dichloroethane	µg/l	0.2 U	0.2 U	0.2 U	800	NE	NE	NE
1,1-Dichloroethene	µg/l	0.2 U	0.2 U	0.2 U	400	NE	NE	NE
1,1-Dichloropropene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
1,2,3-Trichlorobenzene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
1,2,3-Trichloropropane	µg/l	0.2 U	0.2 U	0.2 U	0.00146	NE	NE	NE
1,2,4-Trichlorobenzene	µg/l	0.2 U	0.2 U	0.2 U	80	NE	NE	NE
1,2,4-Trimethylbenzene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
1,2-Dibromo-3-Chloropropane	µg/l	1 U	1 U	1 U	0.0547	NE	NE	NE
1,2-dibromoethane (EDB)	µg/l	0.2 U	0.2 U	0.2 U	0.00051	NE	NE	NE
1,2-Dichlorobenzene (o-Dichlorobenzene)	µg/l	0.2 U	0.2 U	0.2 U	720	NE	NE	NE
1,2-Dichloroethane (EDC)	µg/l	0.2 U	0.2 U	0.2 U	0.48	NE	NE	NE

Analytes	Units	Sample ID and Date			Screening Levels			
		SW1 12/13/2011	SW2 12/13/2011	SW9 2/08/2012	Surface Water MTCA B - Carcinogenic Table Value	Surface Water ARAR - Aquatic Life Fresh/Chronic - Ch. 173-201A WAC	Surface Water ARAR - Aquatic Life Fresh/Chronic - Clean Water Act §304	Surface Water ARAR - Aquatic Life Fresh/Chronic - National Toxics Rule, 40 CFR 131
1,2-Dichloropropane	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
1,3,5-Trimethylbenzene	µg/l	0.2 U	0.2 U	0.2 U	80	NE	NE	NE
1,3-Dichlorobenzene (m-Dichlorobenzene)	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
1,3-Dichloropropane	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
1,4-Dichlorobenzene (p-Dichlorobenzene)	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
2,2-Dichloropropane	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
2-Butanone (MEK)	µg/l	5 U	5 U	5 U	4,800	NE	NE	NE
2-Chloroethyl vinyl ether	µg/l	1 U	1 U	1 U	NE	NE	NE	NE
2-Chlorotoluene	µg/l	0.2 U	0.2 U	0.2 U	160	NE	NE	NE
2-Hexanone	µg/l	2 U	2 U	2 U	NE	NE	NE	NE
4-Chlorotoluene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
4-Methyl-2-Pentanone (Methyl isobutyl ketone)	µg/l	2 U	2 U	2 U	640	NE	NE	NE
Acetone	µg/l	5 U	5 U	5 U	7,200	NE	NE	NE
Benzene	µg/l	0.2 U	0.2 U	0.2 U	0.8	NE	NE	NE
Bromobenzene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
Bromochloromethane	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
Bromodichloromethane	µg/l	0.2 U	0.2 U	0.2 U	0.71	NE	NE	NE
Bromoform (Tribromomethane)	µg/l	1 U	1 U	1 U	5.5	NE	NE	NE
Bromomethane	µg/l	0.2 U	0.2 U	0.2 U	11	NE	NE	NE
Carbon Disulfide	µg/l	0.2 U	0.2 U	0.2 U	800	NE	NE	NE
Carbon Tetrachloride	µg/l	0.2 U	0.2 U	0.2 U	0.625	NE	NE	NE
Chlorobenzene	µg/l	0.2 U	0.2 U	0.2 U	160	NE	NE	NE
Chloroethane	µg/l	1 U	1 U	1 U	NE	NE	NE	NE
Chloroform	µg/l	0.2 U	0.2 U	0.2 U	80	NE	NE	NE
Chloromethane	µg/l	1 U	1 U	1 U	NE	NE	NE	NE
Cis-1,2-Dichloroethene	µg/l	0.2 U	0.2 U	0.2 U	16	NE	NE	NE
Cis-1,3-Dichloropropene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
Dibromochloromethane	µg/l	0.2 U	0.2 U	0.2 U	0.52	NE	NE	NE
Dibromomethane	µg/l	0.2 U	0.2 U	0.2 U	80	NE	NE	NE
Dichlorodifluoromethane (CFC-12)	µg/l	0.25	0.2 U	0.2 U	1,600	NE	NE	NE
Ethylbenzene	µg/l	0.2 U	0.2 U	0.2 U	800	NE	NE	NE
Hexachlorobutadiene	µg/l	0.2 U	0.2 U	0.2 U	0.56	NE	NE	NE
Isopropylbenzene (Cumene)	µg/l	0.2 U	0.2 U	0.2 U	800	NE	NE	NE
Methyl Iodide (Iodomethane)	µg/l	1 U	1 U	1 U	NE	NE	NE	NE
Methyl t-butyl ether	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
Methylene Chloride	µg/l	1 U	1 U	1 U	5.8	NE	NE	NE
Naphthalene	µg/l	1 U	1 U	1 U	160	NE	NE	NE

Analytes	Units	Sample ID and Date			Screening Levels			
		SW1 12/13/2011	SW2 12/13/2011	SW9 2/08/2012	Surface Water MTCA B - Carcinogenic Table Value	Surface Water ARAR - Aquatic Life Fresh/Chronic - Ch. 173-201A WAC	Surface Water ARAR - Aquatic Life Fresh/Chronic - Clean Water Act §304	Surface Water ARAR - Aquatic Life Fresh/Chronic - National Toxics Rule, 40 CFR 131
n-Butylbenzene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
n-Propylbenzene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
p-Isopropyltoluene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
Sec-Butylbenzene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
Styrene	µg/l	0.2 U	0.2 U	0.2 U	1,600	NE	NE	NE
Tert-Butylbenzene	µg/l	0.2 U	0.2 U	0.2 U		NE	NE	NE
Tetrachloroethene	µg/l	0.2 U	0.2 U	0.2 U	0.081	NE	NE	NE
Toluene	µg/l	1 U	1 U	1 U	640	NE	NE	NE
Trans-1,2-Dichloroethene	µg/l	0.2 U	0.2 U	0.2 U	160	NE	NE	NE
Trans-1,3-Dichloropropene	µg/l	0.2 U	0.2 U	0.2 U	NE	NE	NE	NE
Trichloroethene (TCE)	µg/l	0.2 U	0.2 U	0.2 U	0.11	NE	NE	NE
Trichlorofluoromethane (CFC-11)	µg/l	0.62	0.2 U	0.2 U	2,400	NE	NE	NE
Vinyl Acetate	µg/l	2 U	2 U	2 U	8,000	NE	NE	NE
Vinyl Chloride	µg/l	0.2 U	0.2 U	0.2 U	0.0608	NE	NE	NE
Xylene, m-,p-	µg/l	0.4 U	0.4 U	0.4 U	NE	NE	NE	NE
Xylene, o-	µg/l	0.2 U	0.2 U	0.2 U	1,600	NE	NE	NE
Xylene, Total	µg/l	0.6 U	0.6 U	0.6 U	NE	NE	NE	NE
<b>SVOCs by SW8270</b>								
1,2-Diphenylhydrazine	µg/l	--	0.95 U	--	0.11	NE	NE	NE
1,3-Dinitrobenzene	µg/l	--	0.95 U	--	1.6	NE	NE	NE
2,3,4,6-Tetrachlorophenol	µg/l	--	0.95 U	--	480	NE	NE	NE
2,3,5,6-Tetrachlorophenol	µg/l	--	0.95 U	--	NE	NE	NE	NE
2,3-Dichloroaniline	µg/l	--	0.95 U	--	NE	NE	NE	NE
2,4,5-Trichlorophenol	µg/l	--	0.95 U	--	800	NE	NE	NE
2,4,6-Trichlorophenol	µg/l	--	0.95 U	--	4	NE	NE	NE
2,4-Dichlorophenol	µg/l	--	0.95 U	--	24	NE	NE	NE
2,4-Dimethylphenol	µg/l	--	0.95 U	--	160	NE	NE	NE
2,4-Dinitrophenol	µg/l	--	4.8 U	--	32	NE	NE	NE
2,4-Dinitrotoluene	µg/l	--	0.95 U	--	32	NE	NE	NE
2,6-Dinitrotoluene	µg/l	--	0.95 U	--	16	NE	NE	NE
2-Chloronaphthalene	µg/l	--	0.95 U	--	640	NE	NE	NE
2-Chlorophenol	µg/l	--	0.95 U	--	40	NE	NE	NE
2-Nitroaniline	µg/l	--	0.95 U	--	NE	NE	NE	NE
2-Nitrophenol	µg/l	--	0.95 U	--	NE	NE	NE	NE
3,3'-Dichlorobenzidine	µg/l	--	0.95 U	--	0.19	NE	NE	NE
3-Nitroaniline	µg/l	--	0.95 U	--	NE	NE	NE	NE

Analytes	Units	Sample ID and Date			Screening Levels			
		SW1 12/13/2011	SW2 12/13/2011	SW9 2/08/2012	Surface Water MTCA B - Carcinogenic Table Value	Surface Water ARAR - Aquatic Life Fresh/Chronic - Ch. 173-201A WAC	Surface Water ARAR - Aquatic Life Fresh/Chronic - Clean Water Act §304	Surface Water ARAR - Aquatic Life Fresh/Chronic - National Toxics Rule, 40 CFR 131
4,6-Dinitro-2-Methylphenol	µg/l	--	4.8 U	--	NE	NE	NE	NE
4-Bromophenyl phenyl ether	µg/l	--	0.95 U	--	NE	NE	NE	NE
4-Chloro-3-Methylphenol	µg/l	--	0.95 U	--	NE	NE	NE	NE
4-Chloroaniline	µg/l	--	0.95 U	--	32	NE	NE	NE
4-Chlorophenyl-Phenylether	µg/l	--	0.95 U	--	NE	NE	NE	NE
4-Nitroaniline	µg/l	--	0.95 U	--	NE	NE	NE	NE
4-Nitrophenol (p-Nitrophenol)	µg/l	--	0.95 U	--	NE	NE	NE	NE
Aniline	µg/l	--	4.8 U	--	7.7	NE	NE	NE
Benzene, 1,4-Dinitro-	µg/l	--	0.95 U	--	1.6	NE	NE	NE
Benzidine	µg/l	--	4.8 U	--	0.00038	NE	NE	NE
Benzyl Alcohol	µg/l	--	0.95 U	--	800	NE	NE	NE
Bis(2-Chloroethoxy)Methane	µg/l	--	0.95 U	--	NE	NE	NE	NE
Bis(2-Chloroethyl)Ether	µg/l	--	0.95 U	--	0.04	NE	NE	NE
Bis(2-chloroisopropyl) ether	µg/l	--	0.95 U	--	NE	NE	NE	NE
Bis(2-Ethylhexyl) Phthalate	µg/l	--	0.95 U	--	6.3	NE	NE	NE
Butyl benzyl phthalate	µg/l	--	0.95 U	--	3,200	NE	NE	NE
Carbazole	µg/l	--	0.95 U	--	NE	NE	NE	NE
Dibutyl phthalate	µg/l	--	0.95 U	--	1,600	NE	NE	NE
Diethyl phthalate	µg/l	--	0.95 U	--	13,000	NE	NE	NE
Dimethyl phthalate	µg/l	--	0.95 U	--	NE	NE	NE	NE
Di-N-Octyl Phthalate	µg/l	--	0.95 U	--	NE	NE	NE	NE
Hexachlorobenzene	µg/l	--	0.95 U	--	0.055	NE	NE	NE
Hexachlorocyclopentadiene	µg/l	--	0.95 U	--	48	NE	NE	NE
Hexachloroethane	µg/l	--	0.95 U	--	3.1	NE	NE	NE
Hexanedioic Acid, Bis(2-Ethylhexyl) Ester	µg/l	--	4.8 U	--	73	NE	NE	NE
Isophorone	µg/l	--	0.95 U	--	46	NE	NE	NE
m,p-Cresol	µg/l	--	0.95 U	--	NE	NE	NE	NE
Naphthalene	µg/l	0.12	0.095 U	0.1 U	160	NE	NE	NE
Nitrobenzene	µg/l	--	0.95 U	--	16	NE	NE	NE
N-Nitrosodimethylamine	µg/l	--	0.95 U	--	0.00086	NE	NE	NE
N-Nitrosodi-n-propylamine	µg/l	--	0.95 U	--	NE	NE	NE	NE
N-Nitrosodiphenylamine	µg/l	--	0.95 U	--	NE	NE	NE	NE
o-Cresol (2-methylphenol)	µg/l	--	0.95 U	--	400	NE	NE	NE
o-Dinitrobenzene	µg/l	--	0.95 U	--	1.6	NE	NE	NE
Pentachlorophenol	µg/l	--	4.8 U	--	0.219	12.79	15	13

Analytes	Units	Sample ID and Date			Screening Levels			
		SW1 12/13/2011	SW2 12/13/2011	SW9 2/08/2012	Surface Water MTCA B - Carcinogenic Table Value	Surface Water ARAR - Aquatic Life Fresh/Chronic - Ch. 173-201A WAC	Surface Water ARAR - Aquatic Life Fresh/Chronic - Clean Water Act §304	Surface Water ARAR - Aquatic Life Fresh/Chronic - National Toxics Rule, 40 CFR 131
Phenol	µg/l	-	0.95 U	-	<b>2,400</b>	NE	NE	NE
Pyridine	µg/l	-	0.95 U	-	<b>8</b>	NE	NE	NE

**Notes:**

<sup>1</sup> Chemical analysis performed by OnSite Environmental Inc. of Redmond, Washington.

<sup>2</sup> Model Toxics Control Act (MTCA) Cleanup Regulation Chapter 173-340 WAC. Method A clean-up levels.

<sup>3</sup> cPAH testing and regulatory evaluation is completed for individual carcinogenic compounds as well as the for the summation of the mixture of the seven carcinogenic PAHs (known as Ecology's toxicity equivalency methodology). The summation procedure is completed using toxicity equivalency factors for each individual compound which are then added to produce a total toxicity equivalent concentration (TTEC) which is then compared to the MTCA cleanup level of 0.1 µg/l.

ARAR = Applicable or Relevant and Appropriate Requirements

PCBs = Polychlorinated biphenyls

PAHs = Polycyclic aromatic hydrocarbons

VOCs = Volatile organic compound

SVOCs = Semivolatile organic compound

µg/l = microgram per liter

mg/l = milligrams per liter

- = Not analyzed

U = The analyte was not detected at a concentration greater than the given RL or MDL.

E = The concentration reported exceeds the quantitation range and is an estimate.

H = The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and may be impacting the sample result

J = The value reported was below the practical quantitation limit. The value is an estimate.

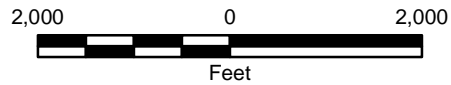
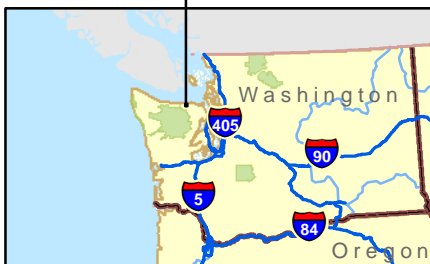
NE = A cleanup level has not been established by Ecology.

**Bold Type** Exceeds One or More MTCA B (Surface Water - Table Value) or ARAR Screening Level

Map Revised: 29 November 2011 syi

Path: P:\0180292\GIS\018029200\_F1.mxd

Office: TACO



Notes:

1. The locations of all features shown are approximate.
2. This drawing is for information purposes. It is intended to assist in showing features discussed in an attached document. GeoEngineers, Inc. can not guarantee the accuracy and content of electronic files. The master file is stored by GeoEngineers, Inc. and will serve as the official record of this communication.
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Data Sources: ESRI Data & Maps, Street Maps 2005  
 Transverse Mercator, State Plane South, North American Datum 1983  
 North arrow oriented to grid north

**Vicinity Map**

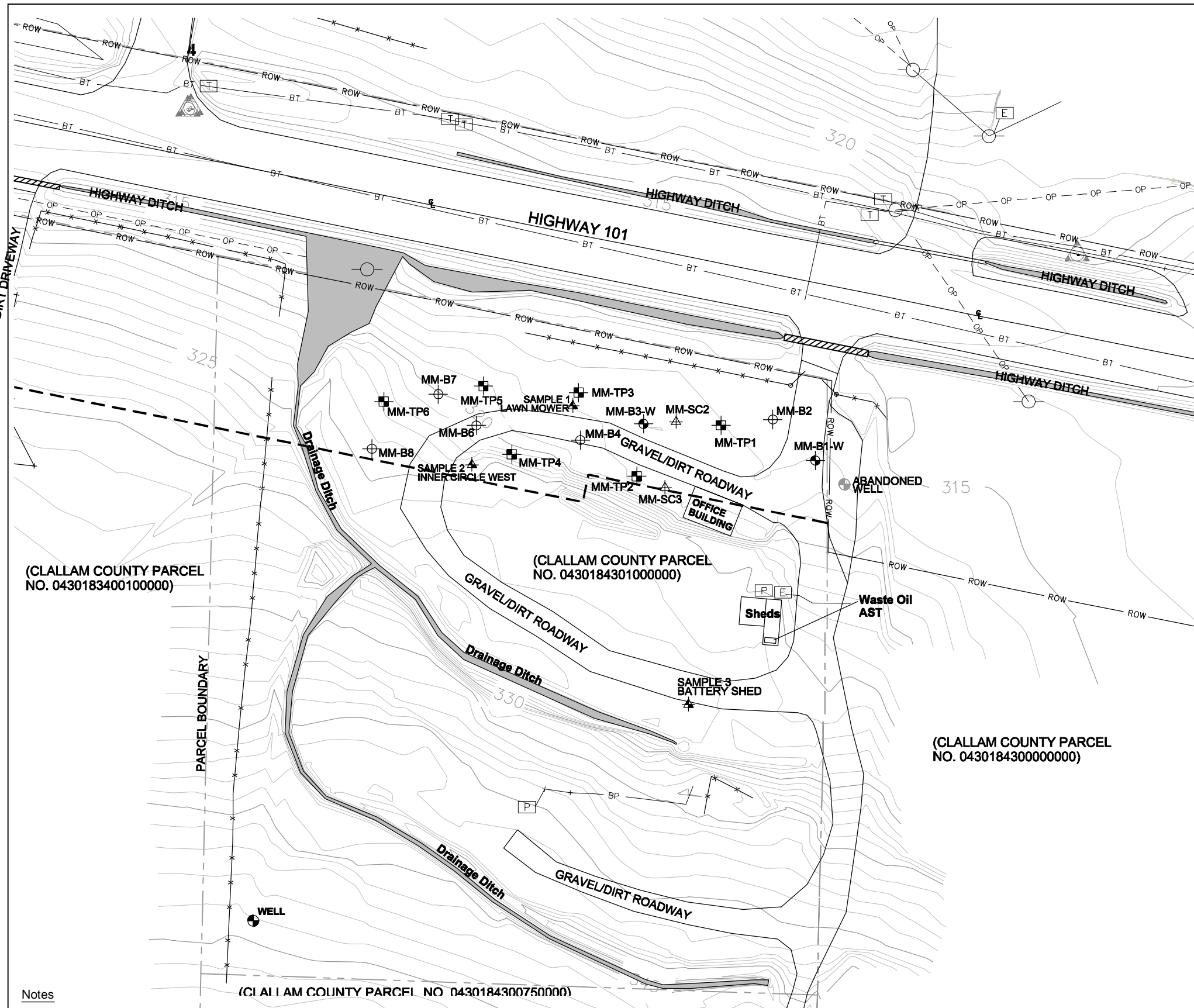
US 101 Midway Metals  
 Clallam County, Washington



**Figure 1**

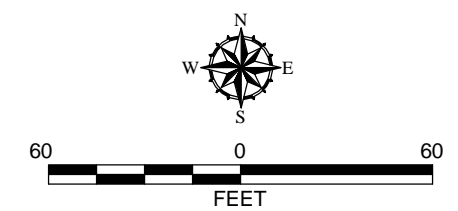


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**LEGEND**

- Approximate Location of Clallam Co. Surface Soil Sample
- Approximate Location of WSDOT Surface Soil Sample
- Approximate Location of WSDOT Direct Push Soil Boring
- Approximate Location of WSDOT Direct Push Soil/Groundwater Boring
- Approximate Location of WSDOT Test Pit
- Location of Existing Groundwater Well
- Location of Abandoned Groundwater Well (Plate Welded On Top of Casing)
- Existing WSDOT Right-of-Way
- Proposed WSDOT Right-of-Way
- Property Boundary
- Fence line
- Elevation Contours (1-foot interval)
- Buried Telephone
- Buried Power
- Overhead Power
- Water Meter
- Power Meter Box
- Utility Pole
- Telephone Vault
- Pad Mount Transformer
- Culvert



**Notes**

1. The locations of all features shown are approximate.
2. This drawing is for information purposes. It is intended to assist in showing features discussed in an attached document. GeoEngineers, Inc. cannot guarantee the accuracy and content of electronic files. The master file is stored by GeoEngineers, Inc. and will serve as the official record of this communication.

Reference: Base files provided by WSDOT Survey Support on January 04, 2012.

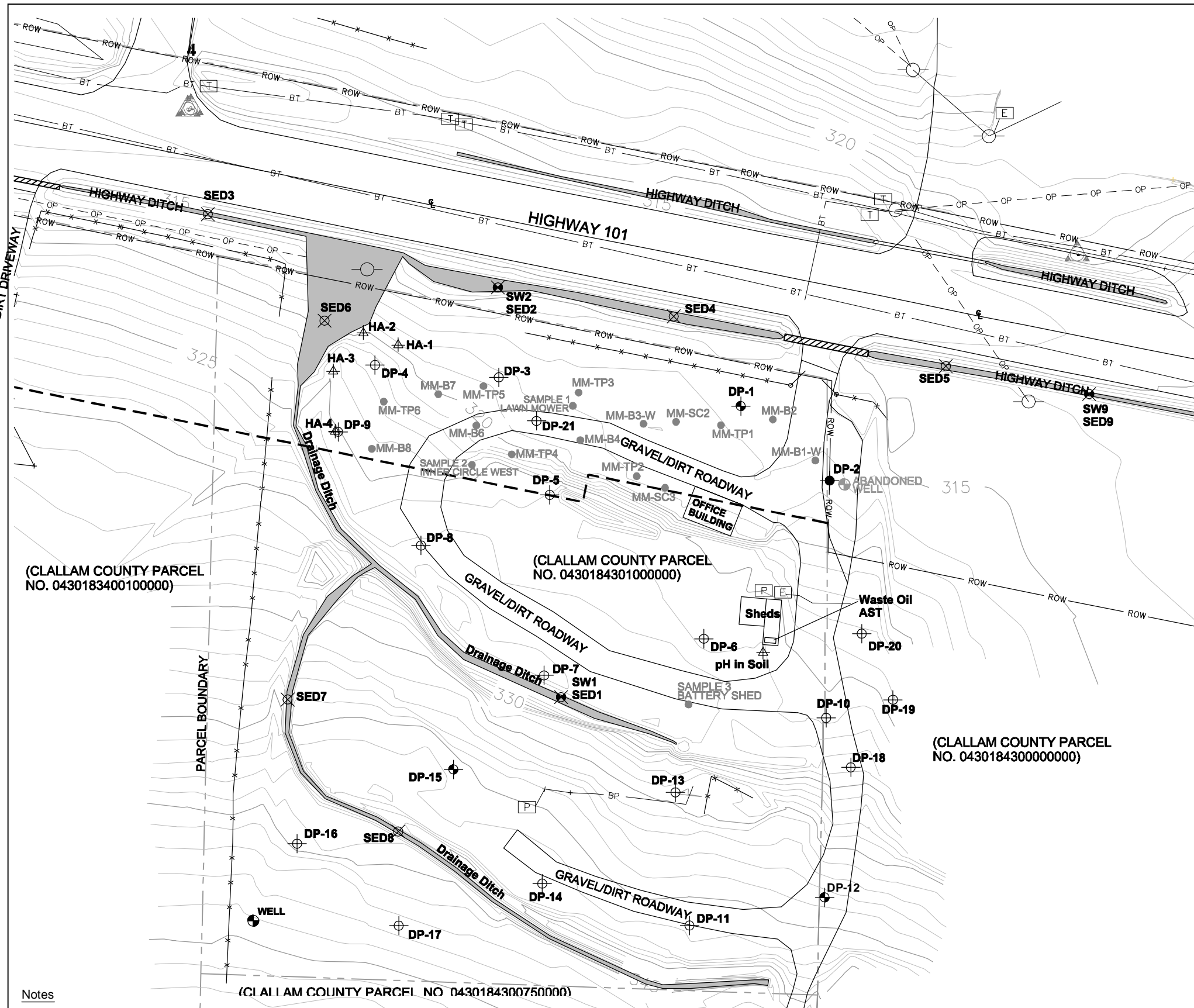
**Site Plan with Previous Sample Locations**

US 101 Midway Metals  
Clallam County, Washington

**Figure 2**

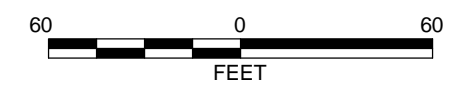


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**LEGEND**

- Approximate Location of Previous Investigations Samples
- ⊗ Location of GeoEngineers Sediment Sample
- ⊗ Location of GeoEngineers Sediment/Surface Water Grab Sample
- ⊕ Location of GeoEngineers Surface Soil Sample (0' to 2')
- ⊕ Location of GeoEngineers Direct Push Soil Boring
- ⊕ Location of GeoEngineers Direct Push Soil/Groundwater Boring
- ⊕ Location of GeoEngineers Direct Push Groundwater Boring
- ⊕ Location of Existing Groundwater Well
- ⊕ Location of Abandoned Groundwater Well (Plate Welded On Top of Casing)
- ROW — Existing WSDOT Right-of-Way
- - - Proposed WSDOT Right-of-Way
- - - Property Boundary
- x x x x Fence line
- 213 Elevation Contours (1-foot interval)
- BT Buried Telephone
- BP Buried Power
- OP OP Overhead Power
- ⊕ Water Meter
- ⊕ Power Meter Box
- ⊕ Utility Pole
- ⊕ Telephone Vault
- ⊕ Pad Mount Transformer
- ▨ Culvert



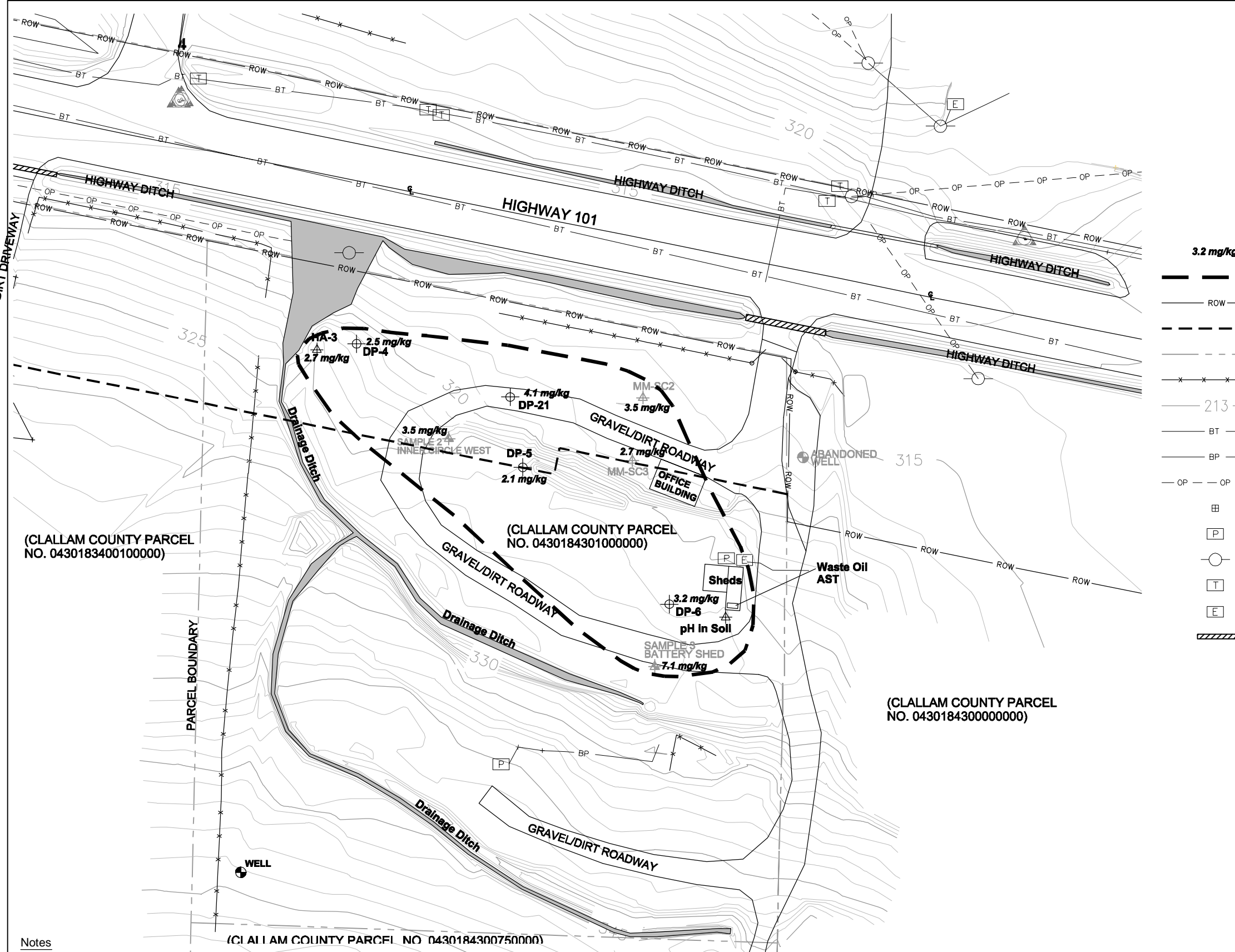
**Notes**

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Reference: Base files provided by WSDOT Survey Support on January 04, 2012.

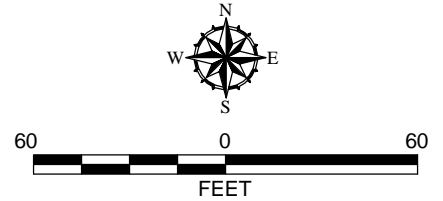
<b>Remedial Investigation Plan</b>	
US 101 Midway Metals Clallam County, Washington	
<b>GEOENGINEERS</b>	<b>Figure 3</b>

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**LEGEND**

- Approximate Location of Clallam Co. Surface Soil Sample
- Approximate Location of WSDOT Surface Soil Sample
- Location of GeoEngineers Surface Soil Sample (0' to 2')
- Location of GeoEngineers Direct Push Soil Boring
- Location of Existing Groundwater Well
- Location of Abandoned Groundwater Well (Plate Welded On Top of Casing)
- 3.2 mg/kg** Cadmium concentration in surface soil (<2' bgs)
- Approximate Extent of Shallow Contamination
- Existing WSDOT Right-of-Way
- Proposed WSDOT Right-of-Way
- Property Boundary
- Fence line
- Elevation Contours (1-foot interval)
- Buried Telephone
- Buried Power
- Overhead Power
- Water Meter
- Power Meter Box
- Utility Pole
- Telephone Vault
- Pad Mount Transformer
- Culvert



**Notes**

1. The locations of all features shown are approximate.
2. This drawing is for information purposes. It is intended to assist in showing features discussed in an attached document. GeoEngineers, Inc. cannot guarantee the accuracy and content of electronic files. The master file is stored by GeoEngineers, Inc. and will serve as the official record of this communication.

Reference: Base files provided by WSDOT Survey Support on January 04, 2012.

<b>Contaminant Boundary</b>	
US 101 Midway Metals Clallam County, Washington	
<b>GEOENGINEERS</b>	<b>Figure 4</b>

**APPENDIX A**  
**Laboratory Analytical Data**





14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

December 23, 2011

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1112-098

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on December 14, 2011.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098  
Project: 0180-292-00

### Case Narrative

Samples were collected on December 13 and 14, 2011 and received by the laboratory on December 14, 2011. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### NWTPH Gx Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

#### Volatiles EPA 8260B Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

The value reported for Acetone in sample DP-1-0-2 exceeds the calibration range and is therefore an estimate. The sample was re-analyzed at the lowest possible dilution allowed by Method 5035A with non-detect results for Acetone.

Internal Standard 1,4-Dichlorobenzene-d4 does not meet acceptance criteria for sample DP-5-0-2 due to sample matrix effects. The sample was re-analyzed with similar results. All results, including Practical Quantitation Limits, from Bromobenzene onward should be considered estimates.

**Please note that any other QA/QC issues associated with these extractions and analyses will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.**

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098  
Project: 0180-292-00

**ANALYTICAL REPORT FOR SAMPLES**

<b>Client ID</b>	<b>Laboratory ID</b>	<b>Matrix</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Notes</b>
DP-1-0-2	12-098-01	Soil	12-13-11	12-14-11	
DP-1-13-15	12-098-03	Soil	12-13-11	12-14-11	
DP-3-0-2	12-098-04	Soil	12-13-11	12-14-11	
DP-4-0-2	12-098-07	Soil	12-13-11	12-14-11	
DP-5-0-2	12-098-09	Soil	12-13-11	12-14-11	
DP-6-0-2	12-098-12	Soil	12-13-11	12-14-11	
DP-6-11.5-12.5	12-098-14	Soil	12-13-11	12-14-11	
DP-7-0-2	12-098-15	Soil	12-14-11	12-14-11	
DP-8-0-2	12-098-17	Soil	12-14-11	12-14-11	
DP-9-0-2	12-098-20	Soil	12-14-11	12-14-11	
DP-10-0-2	12-098-22	Soil	12-14-11	12-14-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

### NWTPH-Gx

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-0-2</b>					
Laboratory ID:	12-098-01					
Gasoline	<b>ND</b>	6.2	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	92	68-124				
<b>Client ID:</b>	<b>DP-1-13-15</b>					
Laboratory ID:	12-098-03					
Gasoline	<b>ND</b>	5.9	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	98	68-124				
<b>Client ID:</b>	<b>DP-3-0-2</b>					
Laboratory ID:	12-098-04					
Gasoline	<b>ND</b>	6.3	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	97	68-124				
<b>Client ID:</b>	<b>DP-4-0-2</b>					
Laboratory ID:	12-098-07					
Gasoline	<b>ND</b>	7.6	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	100	68-124				
<b>Client ID:</b>	<b>DP-5-0-2</b>					
Laboratory ID:	12-098-09					
Gasoline	<b>ND</b>	5.7	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	99	68-124				
<b>Client ID:</b>	<b>DP-6-0-2</b>					
Laboratory ID:	12-098-12					
Gasoline	<b>ND</b>	5.4	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	96	68-124				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

### NWTPH-Gx

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-11.5-12.5</b>					
Laboratory ID:	12-098-14					
Gasoline	<b>ND</b>	5.7	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	94	68-124				
<b>Client ID:</b>	<b>DP-7-0-2</b>					
Laboratory ID:	12-098-15					
Gasoline	<b>ND</b>	7.2	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	99	68-124				
<b>Client ID:</b>	<b>DP-8-0-2</b>					
Laboratory ID:	12-098-17					
Gasoline	<b>18</b>	5.2	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	95	68-124				
<b>Client ID:</b>	<b>DP-9-0-2</b>					
Laboratory ID:	12-098-20					
Gasoline	<b>ND</b>	8.4	NWTPH-Gx	12-16-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	99	68-124				
<b>Client ID:</b>	<b>DP-10-0-2</b>					
Laboratory ID:	12-098-22					
Gasoline	<b>42</b>	5.0	NWTPH-Gx	12-16-11	12-19-11	O
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	100	68-124				



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**NWTPH-Dx**  
**(with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-0-2</b>					
Laboratory ID:	12-098-01					
Diesel Range Organics	<b>ND</b>	29	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	58	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	106	50-150				
<b>Client ID:</b>	<b>DP-1-13-15</b>					
Laboratory ID:	12-098-03					
Diesel Range Organics	<b>ND</b>	30	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	59	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	102	50-150				
<b>Client ID:</b>	<b>DP-3-0-2</b>					
Laboratory ID:	12-098-04					
Diesel Range Organics	<b>ND</b>	180	NWTPH-Dx	12-21-11	12-21-11	U1
Lube Oil	<b>880</b>	59	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	102	50-150				
<b>Client ID:</b>	<b>DP-4-0-2</b>					
Laboratory ID:	12-098-07					
Diesel Range Organics	<b>ND</b>	100	NWTPH-Dx	12-21-11	12-22-11	U1
Lube Oil	<b>580</b>	67	NWTPH-Dx	12-21-11	12-22-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	88	50-150				
<b>Client ID:</b>	<b>DP-5-0-2</b>					
Laboratory ID:	12-098-09					
Diesel Range Organics	<b>ND</b>	200	NWTPH-Dx	12-21-11	12-21-11	U1
Lube Oil	<b>1100</b>	58	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				
<b>Client ID:</b>	<b>DP-6-0-2</b>					
Laboratory ID:	12-098-12					
Diesel Range Organics	<b>280</b>	28	NWTPH-Dx	12-21-11	12-21-11	N
Lube Oil	<b>1100</b>	56	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	129	50-150				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**NWTPH-Dx**  
**(with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-11.5-12.5</b>					
Laboratory ID:	12-098-14					
Diesel Range Organics	<b>ND</b>	30	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	60	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	108	50-150				
<b>Client ID:</b>	<b>DP-7-0-2</b>					
Laboratory ID:	12-098-15					
Diesel Range Organics	<b>34</b>	32	NWTPH-Dx	12-21-11	12-21-11	N
Lube Oil	<b>130</b>	63	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	107	50-150				
<b>Client ID:</b>	<b>DP-8-0-2</b>					
Laboratory ID:	12-098-17					
Diesel Range Organics	<b>ND</b>	30	NWTPH-Dx	12-21-11	12-21-11	U1
Lube Oil	<b>150</b>	55	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	113	50-150				
<b>Client ID:</b>	<b>DP-9-0-2</b>					
Laboratory ID:	12-098-20					
Diesel Range Organics	<b>ND</b>	73	NWTPH-Dx	12-21-11	12-21-11	U1
Lube Oil	<b>360</b>	70	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	108	50-150				
<b>Client ID:</b>	<b>DP-10-0-2</b>					
Laboratory ID:	12-098-22					
Diesel Range Organics	<b>ND</b>	34	NWTPH-Dx	12-21-11	12-21-11	U1
Lube Oil	<b>120</b>	58	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	97	50-150				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
 Page 1 of 2

Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-0-2</b>					
Laboratory ID:	12-098-01					
Dichlorodifluoromethane	0.0022	0.00094	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0047	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0047	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Acetone	0.20	0.0047	EPA 8260	12-16-11	12-16-11	E
Iodomethane	ND	0.0047	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0047	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0047	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
2-Butanone	0.035	0.0047	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Benzene	0.0014	0.00094	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0047	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0047	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0047	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.00094	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-0-2</b>					
Laboratory ID:	12-098-01					
1,1,2-Trichloroethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0047	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0019	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.00094	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	0.0043	0.00094	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0047	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0047	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.00094	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>95</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>96</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>82</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-1-13-15</b>					
<b>Laboratory ID:</b>	<b>12-098-03</b>					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0053	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0053	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Acetone	0.046	0.0053	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0053	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0053	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0053	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Butanone	ND	0.0053	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0053	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0053	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0053	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-13-15</b>					
Laboratory ID:	12-098-03					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0053	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0021	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0053	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0053	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>91</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>98</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>93</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-3-0-2</b>					
Laboratory ID:	12-098-04					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0055	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0055	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	0.64	0.064	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Acetone	0.52	0.32	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0055	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0055	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0055	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Butanone	0.10	0.0055	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0055	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0055	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0055	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-3-0-2</b>					
Laboratory ID:	12-098-04					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0055	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0022	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0055	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0055	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>92</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>95</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>80</i>	<i>55-121</i>				



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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-4-0-2</b>					
<b>Laboratory ID:</b>	<b>12-098-07</b>					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0056	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0056	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	3.6	0.075	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Acetone	ND	0.0056	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0056	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0056	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0056	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Butanone	ND	0.0056	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0056	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0056	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0056	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-4-0-2</b>					
Laboratory ID:	12-098-07					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0056	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0022	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.0011	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.075	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.075	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.38	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.38	EPA 8260	12-19-11	12-19-11	
Naphthalene	ND	0.075	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.075	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>92</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>95</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>77</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-5-0-2</b>					
Laboratory ID:	12-098-09					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloromethane	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Vinyl Chloride	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromomethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloroethane	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Trichlorofluoromethane	0.011	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Acetone	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Carbon Disulfide	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methylene Chloride	ND	0.0052	EPA 8260	12-19-11	12-19-11	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Vinyl Acetate	ND	0.0052	EPA 8260	12-19-11	12-19-11	
2,2-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Butanone	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Bromochloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloroform	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Carbon Tetrachloride	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Benzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Trichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Dibromomethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromodichloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Chloroethyl Vinyl Ether	ND	0.0052	EPA 8260	12-19-11	12-19-11	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methyl Isobutyl Ketone	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Toluene	ND	0.0052	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-5-0-2</b>					
Laboratory ID:	12-098-09					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Tetrachloroethene	0.0040	0.0010	EPA 8260	12-19-11	12-19-11	
1,3-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Hexanone	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Dibromochloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dibromoethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Ethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	ND	0.0021	EPA 8260	12-19-11	12-19-11	
o-Xylene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Styrene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromoform	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Isopropylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.0052	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.0052	EPA 8260	12-19-11	12-19-11	
Naphthalene	0.0042	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>91</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>94</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>79</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-0-2</b>					
Laboratory ID:	12-098-12					
Dichlorodifluoromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	0.0090	0.00095	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Acetone	0.053	0.0048	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0048	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0048	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Butanone	0.0060	0.0048	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0048	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0048	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.00095	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-0-2</b>					
Laboratory ID:	12-098-12					
1,1,2-Trichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	0.0016	0.00095	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0019	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>91</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>95</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>83</i>	<i>55-121</i>				

Date of Report: December 23, 2011  
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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-11.5-12.5</b>					
Laboratory ID:	12-098-14					
Dichlorodifluoromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Acetone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0048	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0048	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Butanone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0048	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0048	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.00095	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-11.5-12.5</b>					
Laboratory ID:	12-098-14					
1,1,2-Trichloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0019	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.00095	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.00095	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>90</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>98</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>93</i>	<i>55-121</i>				



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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-7-0-2</b>					
Laboratory ID:	12-098-15					
Dichlorodifluoromethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Acetone	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0059	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0059	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
2-Butanone	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0059	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0059	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-7-0-2</b>					
Laboratory ID:	12-098-15					
1,1,2-Trichloroethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0023	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.0012	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0059	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0059	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.0012	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>93</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>97</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>89</i>	<i>55-121</i>				

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**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-8-0-2</b>					
Laboratory ID:	12-098-17					
Dichlorodifluoromethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Acetone	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0046	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0046	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
2-Butanone	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0046	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0046	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.00092	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-8-0-2</b>					
Laboratory ID:	12-098-17					
1,1,2-Trichloroethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0018	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.00092	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0046	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0046	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.00092	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>92</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>98</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>89</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-9-0-2</b>					
Laboratory ID:	12-098-20					
Dichlorodifluoromethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0067	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0067	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	0.0037	0.0013	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Acetone	0.097	0.0067	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0067	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0067	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0067	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
2-Butanone	0.0091	0.0067	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0067	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0067	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0067	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-9-0-2</b>					
Laboratory ID:	12-098-20					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0067	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0027	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.0013	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0067	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0067	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.0013	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>95</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>99</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>81</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-10-0-2</b>					
Laboratory ID:	12-098-22					
Dichlorodifluoromethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	0.78	0.059	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Acetone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	0.0095	0.0048	EPA 8260	12-16-11	12-16-11	H
(trans) 1,2-Dichloroethene	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.00096	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0048	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.00096	EPA 8260	12-16-11	12-16-11	
2-Butanone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.00096	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.00096	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Benzene	0.013	0.00096	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.00096	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0048	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Toluene	0.41	0.29	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.00096	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-10-0-2</b>					
Laboratory ID:	12-098-22					
1,1,2-Trichloroethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	0.019	0.00096	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0048	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.00096	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	0.23	0.059	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	1.2	0.12	EPA 8260	12-19-11	12-19-11	
o-Xylene	0.62	0.059	EPA 8260	12-19-11	12-19-11	
Styrene	0.0055	0.00096	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.00096	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	0.040	0.00096	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.059	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.059	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	0.27	0.059	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.059	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.059	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	0.84	0.059	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	2.5	0.059	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	0.098	0.059	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.059	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.29	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.29	EPA 8260	12-19-11	12-19-11	
Naphthalene	0.66	0.059	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.059	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>93</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>99</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>74</i>	<i>55-121</i>				



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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-0-2</b>					
<b>Laboratory ID:</b>	<b>12-098-01</b>					
n-Nitrosodimethylamine	ND	0.039	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	0.39	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.039	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.039	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.039	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.039	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	<b>0.044</b>	0.039	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.039	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.039	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.039	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	0.39	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
Naphthalene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
4-Chloroaniline	ND	0.039	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.039	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Hexachlorocyclopentadiene	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.039	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
3-Nitroaniline	ND	0.039	EPA 8270	12-19-11	12-20-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-0-2</b>					
Laboratory ID:	12-098-01					
2,4-Dinitrophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
Acenaphthene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
4-Nitrophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.039	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.039	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	0.19	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.039	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.039	EPA 8270	12-19-11	12-20-11	
Fluorene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
4,6-Dinitro-2-methylphenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.039	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.039	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.039	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.039	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
Phenanthrene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Carbazole	ND	0.039	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	0.39	EPA 8270	12-19-11	12-20-11	
Fluoranthene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Benzidine	ND	0.39	EPA 8270	12-19-11	12-20-11	
Pyrene	<b>0.013</b>	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Butylbenzylphthalate	ND	0.39	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	0.19	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.39	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
bis(2-Ethylhexyl)phthalate	ND	0.19	EPA 8270	12-19-11	12-20-11	
Di-n-octylphthalate	ND	0.039	EPA 8270	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Indeno[1,2,3-cd]pyrene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0078	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	73	30 - 97				
Phenol-d6	70	40 - 104				
Nitrobenzene-d5	68	35 - 102				
2-Fluorobiphenyl	89	44 - 97				
2,4,6-Tribromophenol	110	41 - 110				
Terphenyl-d14	71	53 - 107				

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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-0-2</b>					
<b>Laboratory ID:</b>	<b>12-098-12</b>					
n-Nitrosodimethylamine	ND	0.19	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	1.9	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.19	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.19	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.19	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.19	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.19	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	1.9	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Naphthalene	<b>0.038</b>	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
4-Chloroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.19	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.045</b>	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>0.016</b>	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Hexachlorocyclopentadiene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	<b>0.0099</b>	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
3-Nitroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-0-2</b>					
Laboratory ID:	12-098-12					
2,4-Dinitrophenol	ND	0.94	EPA 8270	12-19-11	12-20-11	
Acenaphthene	0.0085	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
4-Nitrophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	0.94	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.19	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
Fluorene	0.013	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
4,6-Dinitro-2-methylphenol	ND	0.94	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.19	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.19	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	0.94	EPA 8270	12-19-11	12-20-11	
Phenanthrene	0.040	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	0.015	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Carbazole	ND	0.19	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	1.9	EPA 8270	12-19-11	12-20-11	
Fluoranthene	0.045	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzidine	ND	1.9	EPA 8270	12-19-11	12-20-11	
Pyrene	0.063	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Butylbenzylphthalate	ND	1.9	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	0.94	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	1.9	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	0.026	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	0.035	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
bis(2-Ethylhexyl)phthalate	0.97	0.94	EPA 8270	12-19-11	12-20-11	
Di-n-octylphthalate	ND	0.19	EPA 8270	12-19-11	12-20-11	
Benzo[b]fluoranthene	0.047	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	0.016	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	0.033	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Indeno[1,2,3-cd]pyrene	0.030	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	0.038	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	46	30 - 97				
Phenol-d6	47	40 - 104				
Nitrobenzene-d5	39	35 - 102				
2-Fluorobiphenyl	77	44 - 97				
2,4,6-Tribromophenol	101	41 - 110				
Terphenyl-d14	66	53 - 107				

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 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-13-15</b>					
Laboratory ID:	12-098-03					
Naphthalene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0079	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>50</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>55</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>51</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-3-0-2</b>					
Laboratory ID:	12-098-04					
Naphthalene	<b>0.021</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.020</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>0.010</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>0.011</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.034</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>0.011</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.045</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.047</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>0.024</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>0.036</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>0.039</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>0.011</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>0.028</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>0.020</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>0.025</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>57</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>64</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>78</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
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**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-4-0-2</b>					
Laboratory ID:	12-098-07					
Naphthalene	<b>0.015</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.018</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.017</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.017</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.020</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>0.014</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>0.015</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>0.0095</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>0.012</b>	0.0089	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>57</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>62</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>64</i>	<i>33 - 119</i>				

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**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-5-0-2</b>					
Laboratory ID:	12-098-09					
Naphthalene	<b>0.033</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.032</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>0.015</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>0.0082</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.049</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>0.014</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.063</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.068</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>0.036</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>0.050</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>0.062</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>0.017</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>0.046</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>0.038</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>0.0092</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>0.045</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>69</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>65</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>81</i>	<i>33 - 119</i>				



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-11.5-12.5</b>					
Laboratory ID:	12-098-14					
Naphthalene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	ND	0.0080	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>57</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>66</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>60</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-7-0-2</b>					
Laboratory ID:	12-098-15					
Naphthalene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.0088</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.012</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.016</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.015</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>0.027</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>0.0097</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0085	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>54</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>59</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>56</i>	<i>33 - 119</i>				

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 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-8-0-2</b>					
Laboratory ID:	12-098-17					
Naphthalene	<b>0.0081</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.0082</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.0085</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.010</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>0.0091</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0073	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>59</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>64</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>67</i>	<i>33 - 119</i>				

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**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-9-0-2</b>					
Laboratory ID:	12-098-20					
Naphthalene	<b>0.016</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.013</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>ND</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>0.037</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>0.019</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>0.028</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.22</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>0.072</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.35</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.28</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>0.16</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>0.15</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>0.17</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>0.057</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>0.14</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>0.082</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>0.020</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>0.076</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>50</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>60</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>61</i>	<i>33 - 119</i>				

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**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-10-0-2</b>					
Laboratory ID:	12-098-22					
Naphthalene	<b>0.029</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.037</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>0.020</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.013</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.010</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.011</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0078	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>58</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>64</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>62</i>	<i>33 - 119</i>				

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### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-1-0-2</b>					
Laboratory ID:	12-098-01					
Aroclor 1016	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.058	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	81	42-123				
<b>Client ID:</b>	<b>DP-1-13-15</b>					
Laboratory ID:	12-098-03					
Aroclor 1016	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.059	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	79	42-123				
<b>Client ID:</b>	<b>DP-3-0-2</b>					
Laboratory ID:	12-098-04					
Aroclor 1016	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.059	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.059	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	70	42-123				

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 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-4-0-2</b>					
Laboratory ID:	12-098-07					
Aroclor 1016	ND	0.067	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.067	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.067	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	0.42	0.067	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.067	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.067	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.067	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	65	42-123				
<b>Client ID:</b>	<b>DP-5-0-2</b>					
Laboratory ID:	12-098-09					
Aroclor 1016	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	0.089	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	0.14	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.058	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	77	42-123				
<b>Client ID:</b>	<b>DP-6-0-2</b>					
Laboratory ID:	12-098-12					
Aroclor 1016	ND	0.056	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.056	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.056	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	0.14	0.056	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.056	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.056	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	0.13	0.056	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	79	42-123				

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### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-11.5-12.5</b>					
Laboratory ID:	12-098-14					
Aroclor 1016	ND	0.060	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.060	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.060	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	ND	0.060	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.060	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.060	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.060	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	82	42-123				
<b>Client ID:</b>	<b>DP-7-0-2</b>					
Laboratory ID:	12-098-15					
Aroclor 1016	ND	0.063	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.063	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.063	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	0.12	0.063	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.063	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.063	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	0.079	0.063	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	78	42-123				
<b>Client ID:</b>	<b>DP-8-0-2</b>					
Laboratory ID:	12-098-17					
Aroclor 1016	ND	0.055	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.055	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.055	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	ND	0.055	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.055	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	0.058	0.055	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.055	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	83	42-123				



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### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-9-0-2</b>					
Laboratory ID:	12-098-20					
Aroclor 1016	ND	0.070	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.070	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.070	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	ND	0.070	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.070	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.070	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.070	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	81	42-123				
<b>Client ID:</b>	<b>DP-10-0-2</b>					
Laboratory ID:	12-098-22					
Aroclor 1016	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	ND	0.058	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	ND	0.058	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	81	42-123				

Date of Report: December 23, 2011  
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**TOTAL METALS**  
**EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-01					
<b>Client ID:</b>	<b>DP-1-0-2</b>					
Arsenic	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Barium	<b>70</b>	2.9	6010B	12-19-11	12-19-11	
Cadmium	<b>ND</b>	0.58	6010B	12-19-11	12-19-11	
Chromium	<b>34</b>	0.58	6010B	12-19-11	12-19-11	
Lead	<b>ND</b>	5.8	6010B	12-19-11	12-19-11	
Mercury	<b>ND</b>	0.29	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.58	6010B	12-19-11	12-19-11	

Lab ID:	12-098-03					
<b>Client ID:</b>	<b>DP-1-13-15</b>					
Arsenic	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Barium	<b>30</b>	3.0	6010B	12-19-11	12-19-11	
Cadmium	<b>ND</b>	0.59	6010B	12-19-11	12-19-11	
Chromium	<b>29</b>	0.59	6010B	12-19-11	12-19-11	
Lead	<b>ND</b>	5.9	6010B	12-19-11	12-19-11	
Mercury	<b>ND</b>	0.30	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.59	6010B	12-19-11	12-19-11	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-04					
<b>Client ID:</b>	<b>DP-3-0-2</b>					
Arsenic	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Barium	<b>89</b>	2.9	6010B	12-19-11	12-19-11	
Cadmium	<b>0.80</b>	0.59	6010B	12-19-11	12-19-11	
Chromium	<b>36</b>	0.59	6010B	12-19-11	12-19-11	
Lead	<b>40</b>	5.9	6010B	12-19-11	12-19-11	
Mercury	<b>ND</b>	0.29	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.59	6010B	12-19-11	12-19-11	

Lab ID:	12-098-07					
<b>Client ID:</b>	<b>DP-4-0-2</b>					
Arsenic	<b>ND</b>	13	6010B	12-19-11	12-19-11	
Barium	<b>99</b>	3.3	6010B	12-19-11	12-19-11	
Cadmium	<b>2.5</b>	0.67	6010B	12-19-11	12-19-11	
Chromium	<b>42</b>	0.67	6010B	12-19-11	12-19-11	
Lead	<b>150</b>	6.7	6010B	12-19-11	12-19-11	
Mercury	<b>0.67</b>	0.33	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	13	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.67	6010B	12-19-11	12-19-11	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-09					
<b>Client ID:</b>	<b>DP-5-0-2</b>					
Arsenic	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Barium	<b>120</b>	2.9	6010B	12-19-11	12-19-11	
Cadmium	<b>2.1</b>	0.58	6010B	12-19-11	12-19-11	
Chromium	<b>35</b>	0.58	6010B	12-19-11	12-19-11	
Lead	<b>130</b>	5.8	6010B	12-19-11	12-19-11	
Mercury	<b>0.32</b>	0.29	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	12	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.58	6010B	12-19-11	12-19-11	

Lab ID:	12-098-12					
<b>Client ID:</b>	<b>DP-6-0-2</b>					
Arsenic	<b>ND</b>	11	6010B	12-19-11	12-19-11	
Barium	<b>72</b>	2.8	6010B	12-19-11	12-19-11	
Cadmium	<b>3.2</b>	0.56	6010B	12-19-11	12-19-11	
Chromium	<b>41</b>	0.56	6010B	12-19-11	12-19-11	
Lead	<b>160</b>	5.6	6010B	12-19-11	12-19-11	
Mercury	<b>ND</b>	0.28	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	11	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.56	6010B	12-19-11	12-19-11	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-14					
<b>Client ID:</b>	<b>DP-6-11.5-12.5</b>					
Arsenic	ND	12	6010B	12-19-11	12-19-11	
Barium	42	3.0	6010B	12-19-11	12-19-11	
Cadmium	ND	0.60	6010B	12-19-11	12-19-11	
Chromium	26	0.60	6010B	12-19-11	12-19-11	
Lead	ND	6.0	6010B	12-19-11	12-19-11	
Mercury	ND	0.30	7471A	12-19-11	12-19-11	
Selenium	ND	12	6010B	12-19-11	12-19-11	
Silver	ND	0.60	6010B	12-19-11	12-19-11	

Lab ID:	12-098-15					
<b>Client ID:</b>	<b>DP-7-0-2</b>					
Arsenic	ND	13	6010B	12-19-11	12-19-11	
Barium	88	3.2	6010B	12-19-11	12-19-11	
Cadmium	0.81	0.63	6010B	12-19-11	12-19-11	
Chromium	31	0.63	6010B	12-19-11	12-19-11	
Lead	93	6.3	6010B	12-19-11	12-19-11	
Mercury	ND	0.32	7471A	12-19-11	12-19-11	
Selenium	ND	13	6010B	12-19-11	12-19-11	
Silver	ND	0.63	6010B	12-19-11	12-19-11	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-17					
<b>Client ID:</b>	<b>DP-8-0-2</b>					
Arsenic	ND	11	6010B	12-19-11	12-19-11	
Barium	46	2.7	6010B	12-19-11	12-19-11	
Cadmium	ND	0.55	6010B	12-19-11	12-19-11	
Chromium	29	0.55	6010B	12-19-11	12-19-11	
Lead	24	5.5	6010B	12-19-11	12-19-11	
Mercury	ND	0.27	7471A	12-19-11	12-19-11	
Selenium	ND	11	6010B	12-19-11	12-19-11	
Silver	ND	0.55	6010B	12-19-11	12-19-11	

Lab ID:	12-098-20					
<b>Client ID:</b>	<b>DP-9-0-2</b>					
Arsenic	ND	14	6010B	12-19-11	12-19-11	
Barium	84	3.5	6010B	12-19-11	12-19-11	
Cadmium	1.6	0.70	6010B	12-19-11	12-19-11	
Chromium	46	0.70	6010B	12-19-11	12-19-11	
Lead	85	7.0	6010B	12-19-11	12-19-11	
Mercury	ND	0.35	7471A	12-19-11	12-19-11	
Selenium	ND	14	6010B	12-19-11	12-19-11	
Silver	ND	0.70	6010B	12-19-11	12-19-11	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-22					
Client ID:	DP-10-0-2					
Arsenic	ND	12	6010B	12-19-11	12-19-11	
Barium	70	2.9	6010B	12-19-11	12-19-11	
Cadmium	ND	0.58	6010B	12-19-11	12-19-11	
Chromium	34	0.58	6010B	12-19-11	12-19-11	
Lead	8.3	5.8	6010B	12-19-11	12-19-11	
Mercury	ND	0.29	7471A	12-19-11	12-19-11	
Selenium	ND	12	6010B	12-19-11	12-19-11	
Silver	ND	0.58	6010B	12-19-11	12-19-11	

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**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1216S1					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	93	68-124				
Laboratory ID:	MB1216S2					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	87	68-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-099-08							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	30	
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				100	103	68-124		
Laboratory ID:	12-100-03							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	30	
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				102	106	68-124		



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**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1221S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>111</i>	<i>50-150</i>				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-098-14					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			108 109	50-150		
Laboratory ID:	12-098-17					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	U1
Lube Oil	<b>137</b>	<b>117</b>		16	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			113 101	50-150		

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Chloromethane	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Vinyl Chloride	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Bromomethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Chloroethane	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Trichlorofluoromethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Acetone	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Iodomethane	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Carbon Disulfide	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Methylene Chloride	ND	0.0050	EPA 8260	12-16-11	12-16-11	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,1-Dichloroethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Vinyl Acetate	ND	0.0050	EPA 8260	12-16-11	12-16-11	
2,2-Dichloropropane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
2-Butanone	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Bromochloromethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Chloroform	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Carbon Tetrachloride	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,1-Dichloropropene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Benzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2-Dichloroethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Trichloroethene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2-Dichloropropane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Dibromomethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Bromodichloromethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	12-16-11	12-16-11	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Toluene	ND	0.0050	EPA 8260	12-16-11	12-16-11	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-16-11	12-16-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Tetrachloroethene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,3-Dichloropropane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
2-Hexanone	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Dibromochloromethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2-Dibromoethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Chlorobenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Ethylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
m,p-Xylene	ND	0.0020	EPA 8260	12-16-11	12-16-11	
o-Xylene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Styrene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Bromoform	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Isopropylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Bromobenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	12-16-11	12-16-11	
n-Propylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
2-Chlorotoluene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
4-Chlorotoluene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
tert-Butylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
sec-Butylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
p-Isopropyltoluene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
n-Butylbenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	12-16-11	12-16-11	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
Hexachlorobutadiene	ND	0.0050	EPA 8260	12-16-11	12-16-11	
Naphthalene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>93</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>101</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>95</i>	<i>55-121</i>				

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**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloromethane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Vinyl Chloride	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromomethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloroethane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Trichlorofluoromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Acetone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Carbon Disulfide	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methylene Chloride	ND	0.0050	EPA 8260	12-19-11	12-19-11	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Vinyl Acetate	ND	0.0050	EPA 8260	12-19-11	12-19-11	
2,2-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Butanone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Bromochloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloroform	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Carbon Tetrachloride	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Benzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Trichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Dibromomethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromodichloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	12-19-11	12-19-11	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Toluene	ND	0.0050	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Tetrachloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Hexanone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Dibromochloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dibromoethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Ethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	ND	0.0020	EPA 8260	12-19-11	12-19-11	
o-Xylene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Styrene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromoform	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Isopropylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Naphthalene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>95</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>99</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>99</i>	<i>55-121</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID: SB1216S1										
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0554</b>	<b>0.0572</b>	0.0500	0.0500	111	114	70-130	3	19	
Benzene	<b>0.0505</b>	<b>0.0515</b>	0.0500	0.0500	101	103	70-125	2	15	
Trichloroethene	<b>0.0482</b>	<b>0.0477</b>	0.0500	0.0500	96	95	70-122	1	14	
Toluene	<b>0.0494</b>	<b>0.0489</b>	0.0500	0.0500	99	98	73-120	1	16	
Chlorobenzene	<b>0.0465</b>	<b>0.0446</b>	0.0500	0.0500	93	89	74-109	4	12	
<i>Surrogate:</i>										
					85	85	63-127			
					89	91	65-129			
					85	86	55-121			
Laboratory ID: SB1219S1										
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0559</b>	<b>0.0550</b>	0.0500	0.0500	112	110	70-130	2	19	
Benzene	<b>0.0514</b>	<b>0.0506</b>	0.0500	0.0500	103	101	70-125	2	15	
Trichloroethene	<b>0.0471</b>	<b>0.0470</b>	0.0500	0.0500	94	94	70-122	0	14	
Toluene	<b>0.0493</b>	<b>0.0484</b>	0.0500	0.0500	99	97	73-120	2	16	
Chlorobenzene	<b>0.0455</b>	<b>0.0432</b>	0.0500	0.0500	91	86	74-109	5	12	
<i>Surrogate:</i>										
					89	84	63-127			
					91	92	65-129			
					87	88	55-121			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S2					
n-Nitrosodimethylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.033	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	0.33	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Naphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4-Chloroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
3-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S2					
2,4-Dinitrophenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4-Nitrophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	0.17	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Carbazole	ND	0.033	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	0.33	EPA 8270	12-19-11	12-20-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzidine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Butylbenzylphthalate	ND	0.33	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	0.17	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
bis(2-Ethylhexyl)phthalate	ND	0.17	EPA 8270	12-19-11	12-20-11	
Di-n-octylphthalate	ND	0.033	EPA 8270	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	85	30 - 97				
Phenol-d6	80	40 - 104				
Nitrobenzene-d5	74	35 - 102				
2-Fluorobiphenyl	92	44 - 97				
2,4,6-Tribromophenol	103	41 - 110				
Terphenyl-d14	79	53 - 107				



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
	SB	SBD	SB	SBD	SB	SBD				
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1219S2									
Phenol	<b>0.925</b>	<b>0.950</b>	1.33	1.33	70	71	31 - 111	3	34	
2-Chlorophenol	<b>1.04</b>	<b>1.06</b>	1.33	1.33	78	80	29 - 112	2	37	
1,4-Dichlorobenzene	<b>0.522</b>	<b>0.528</b>	0.667	0.667	78	79	24 - 100	1	37	
n-Nitroso-di-n-propylamine	<b>0.396</b>	<b>0.405</b>	0.667	0.667	59	61	35 - 104	2	32	
1,2,4-Trichlorobenzene	<b>0.685</b>	<b>0.713</b>	0.667	0.667	103	107	29 - 110	4	35	
4-Chloro-3-methylphenol	<b>1.13</b>	<b>1.19</b>	1.33	1.33	85	89	53 - 104	5	25	
Acenaphthene	<b>0.544</b>	<b>0.563</b>	0.667	0.667	82	84	50 - 95	3	23	
4-Nitrophenol	<b>1.22</b>	<b>1.30</b>	1.33	1.33	92	98	42 - 126	6	30	
2,4-Dinitrotoluene	<b>0.722</b>	<b>0.749</b>	0.667	0.667	108	112	53 - 115	4	31	
Pentachlorophenol	<b>0.939</b>	<b>1.03</b>	1.33	1.33	71	77	50 - 116	9	30	
Pyrene	<b>0.446</b>	<b>0.468</b>	0.667	0.667	67	70	57 - 120	5	27	
<i>Surrogate:</i>										
2-Fluorophenol					83	82	30 - 97			
Phenol-d6					76	78	40 - 104			
Nitrobenzene-d5					72	74	35 - 102			
2-Fluorobiphenyl					91	91	44 - 97			
2,4,6-Tribromophenol					105	111	41 - 110			
Terphenyl-d14					78	82	53 - 107			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
Naphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>75</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>85</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>77</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent		Recovery	RPD	RPD	Flags
	MS	MSD	MS	MSD	Result	Recovery	Limits	RPD	Limit		
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-110-03										
	MS	MSD	MS	MSD		MS	MSD				
Naphthalene	<b>0.0546</b>	<b>0.0593</b>	0.0833	0.0833	ND	66	71	39 - 110	8	21	
Acenaphthylene	<b>0.0530</b>	<b>0.0593</b>	0.0833	0.0833	ND	64	71	47 - 124	11	21	
Acenaphthene	<b>0.0489</b>	<b>0.0560</b>	0.0833	0.0833	ND	59	67	50 - 120	14	20	
Fluorene	<b>0.0487</b>	<b>0.0568</b>	0.0833	0.0833	ND	58	68	52 - 126	15	21	
Phenanthrene	<b>0.0478</b>	<b>0.0551</b>	0.0833	0.0833	ND	57	66	41 - 130	14	22	
Anthracene	<b>0.0473</b>	<b>0.0554</b>	0.0833	0.0833	ND	57	67	48 - 124	16	23	
Fluoranthene	<b>0.0495</b>	<b>0.0581</b>	0.0833	0.0833	ND	59	70	40 - 137	16	23	
Pyrene	<b>0.0487</b>	<b>0.0579</b>	0.0833	0.0833	ND	58	70	36 - 139	17	23	
Benzo[a]anthracene	<b>0.0496</b>	<b>0.0593</b>	0.0833	0.0833	ND	60	71	43 - 127	18	21	
Chrysene	<b>0.0476</b>	<b>0.0570</b>	0.0833	0.0833	ND	57	68	41 - 133	18	19	
Benzo[b]fluoranthene	<b>0.0452</b>	<b>0.0555</b>	0.0833	0.0833	ND	54	67	40 - 132	20	25	
Benzo(j,k)fluoranthene	<b>0.0436</b>	<b>0.0526</b>	0.0833	0.0833	ND	52	63	35 - 132	19	25	
Benzo[a]pyrene	<b>0.0471</b>	<b>0.0563</b>	0.0833	0.0833	ND	57	68	37 - 131	18	26	
Indeno(1,2,3-c,d)pyrene	<b>0.0444</b>	<b>0.0541</b>	0.0833	0.0833	ND	53	65	39 - 134	20	23	
Dibenz[a,h]anthracene	<b>0.0457</b>	<b>0.0553</b>	0.0833	0.0833	ND	55	66	40 - 137	19	21	
Benzo[g,h,i]perylene	<b>0.0456</b>	<b>0.0550</b>	0.0833	0.0833	ND	55	66	35 - 135	19	22	
<i>Surrogate:</i>											
<i>2-Fluorobiphenyl</i>						<i>54</i>	<i>60</i>	<i>43 - 109</i>			
<i>Pyrene-d10</i>						<i>56</i>	<i>67</i>	<i>38 - 128</i>			
<i>Terphenyl-d14</i>						<i>53</i>	<i>63</i>	<i>33 - 119</i>			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1218S1					
Aroclor 1016	<b>ND</b>	0.050	EPA 8082	12-18-11	12-19-11	
Aroclor 1221	<b>ND</b>	0.050	EPA 8082	12-18-11	12-19-11	
Aroclor 1232	<b>ND</b>	0.050	EPA 8082	12-18-11	12-19-11	
Aroclor 1242	<b>ND</b>	0.050	EPA 8082	12-18-11	12-19-11	
Aroclor 1248	<b>ND</b>	0.050	EPA 8082	12-18-11	12-19-11	
Aroclor 1254	<b>ND</b>	0.050	EPA 8082	12-18-11	12-19-11	
Aroclor 1260	<b>ND</b>	0.050	EPA 8082	12-18-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	89		42-123			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-098-03										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	<b>0.396</b>	<b>0.443</b>	0.500	0.500	ND	<b>79</b>	<b>89</b>	44-125	11	15	
<i>Surrogate:</i>											
DCB						74	84	42-123			

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098  
Project: 0180-292-00

**TOTAL METALS  
EPA 6010B/7471A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-19-11  
Date Analyzed: 12-19-11  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB1219SM1&MB1219S1

Analyte	Method	Result	PQL
Arsenic	6010B	ND	10
Barium	6010B	ND	2.5
Cadmium	6010B	ND	0.50
Chromium	6010B	ND	0.50
Lead	6010B	ND	5.0
Mercury	7471A	ND	0.25
Selenium	6010B	ND	10
Silver	6010B	ND	0.50

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-19-11  
 Date Analyzed: 12-19-11  
 Matrix: Soil  
 Units: mg/kg (ppm)  
 Lab ID: 12-098-17

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	42.3	47.4	11	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	26.3	27.9	6	0.50	
Lead	21.6	22.6	5	5.0	
Mercury	ND	ND	NA	0.25	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-19-11

Date Analyzed: 12-19-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 12-098-17

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>88.7</b>	89	<b>89.0</b>	89	0	
Barium	100	<b>142</b>	100	<b>140</b>	97	2	
Cadmium	50.0	<b>47.0</b>	94	<b>47.2</b>	94	0	
Chromium	100	<b>119</b>	92	<b>121</b>	94	2	
Lead	250	<b>253</b>	93	<b>241</b>	88	5	
Mercury	0.500	<b>0.512</b>	102	<b>0.490</b>	98	4	
Selenium	100	<b>91.6</b>	92	<b>92.3</b>	92	1	
Silver	25.0	<b>21.0</b>	84	<b>21.1</b>	84	0	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098  
Project: 0180-292-00

**% MOISTURE**

Date Analyzed: 12-16-11

Client ID	Lab ID	% Moisture
DP-1-0-2	12-098-01	14
DP-1-13-15	12-098-03	16
DP-3-0-2	12-098-04	15
DP-4-0-2	12-098-07	25
DP-5-0-2	12-098-09	13
DP-6-0-2	12-098-12	11
DP-6-11.5-12.5	12-098-14	17
DP-7-0-2	12-098-15	21
DP-8-0-2	12-098-17	9
DP-9-0-2	12-098-20	29
DP-10-0-2	12-098-22	14





### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference



14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

December 27, 2011

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1112-110

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on December 15, 2011.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: December 27, 2011  
Samples Submitted: December 15, 2011  
Laboratory Reference: 1112-110  
Project: 0180-292-00

### Case Narrative

Samples were collected on December 14, 2011 and received by the laboratory on December 15, 2011. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### NWTPH Gx and Volatiles EPA 8260B Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

#### PAHs EPA 8270D/SIM Analysis

Sample DP-DUPE-1 had one surrogate recovery out of control limits. This is within allowance of our standard operating procedure as long as the recovery is above 10%.

**Please note that any other QA/QC issues associated with these extractions and analyses will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.**

Date of Report: December 27, 2011  
Samples Submitted: December 15, 2011  
Laboratory Reference: 1112-110  
Project: 0180-292-00

**ANALYTICAL REPORT FOR SAMPLES**

<b>Client ID</b>	<b>Laboratory ID</b>	<b>Matrix</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Notes</b>
DP-10-13-15	12-110-01	Soil	12-14-11	12-15-11	
DP-11-0-2	12-110-02	Soil	12-14-11	12-15-11	
DP-11-2-3.5	12-110-03	Soil	12-14-11	12-15-11	
DP-DUPE-1	12-110-05	Soil	12-14-11	12-15-11	

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**NWTPH-Gx**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-10-13-15</b>					
Laboratory ID:	12-110-01					
Gasoline	<b>ND</b>	5.1	NWTPH-Gx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	92	68-124				
<b>Client ID:</b>	<b>DP-11-0-2</b>					
Laboratory ID:	12-110-02					
Gasoline	<b>ND</b>	5.8	NWTPH-Gx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	100	68-124				
<b>Client ID:</b>	<b>DP-11-2-3.5</b>					
Laboratory ID:	12-110-03					
Gasoline	<b>ND</b>	5.1	NWTPH-Gx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	102	68-124				
<b>Client ID:</b>	<b>DP-DUPE-1</b>					
Laboratory ID:	12-110-05					
Gasoline	<b>ND</b>	5.7	NWTPH-Gx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	99	68-124				

Date of Report: December 27, 2011  
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**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-10-13-15</b>					
Laboratory ID:	12-110-01					
Diesel Range Organics	<b>ND</b>	28	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	56	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	109	50-150				

<b>Client ID:</b>	<b>DP-11-0-2</b>					
Laboratory ID:	12-110-02					
Diesel Range Organics	<b>ND</b>	91	NWTPH-Dx	12-19-11	12-19-11	U1
Lube Oil	<b>480</b>	58	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	103	50-150				

<b>Client ID:</b>	<b>DP-11-2-3.5</b>					
Laboratory ID:	12-110-03					
Diesel Range Organics	<b>ND</b>	29	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	58	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	115	50-150				

<b>Client ID:</b>	<b>DP-DUPE-1</b>					
Laboratory ID:	12-110-05					
Diesel Range Organics	<b>ND</b>	28	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	56	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	104	50-150				

Date of Report: December 27, 2011  
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**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-10-13-15</b>					
Laboratory ID:	12-110-01					
Dichlorodifluoromethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Chloromethane	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Vinyl Chloride	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Bromomethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Chloroethane	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Trichlorofluoromethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Acetone	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Carbon Disulfide	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Methylene Chloride	ND	0.0045	EPA 8260	12-19-11	12-19-11	
(trans) 1,2-Dichloroethene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Methyl t-Butyl Ether	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Vinyl Acetate	ND	0.0045	EPA 8260	12-19-11	12-19-11	
2,2-Dichloropropane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
(cis) 1,2-Dichloroethene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
2-Butanone	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Bromochloromethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Chloroform	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,1,1-Trichloroethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Carbon Tetrachloride	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,1-Dichloropropene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Benzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2-Dichloroethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Trichloroethene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2-Dichloropropane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Dibromomethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Bromodichloromethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
2-Chloroethyl Vinyl Ether	ND	0.0045	EPA 8260	12-19-11	12-19-11	
(cis) 1,3-Dichloropropene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Methyl Isobutyl Ketone	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Toluene	ND	0.0045	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.00090	EPA 8260	12-19-11	12-19-11	

Date of Report: December 27, 2011  
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**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-10-13-15</b>					
Laboratory ID:	12-110-01					
1,1,2-Trichloroethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Tetrachloroethene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,3-Dichloropropane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
2-Hexanone	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Dibromochloromethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2-Dibromoethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Chlorobenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,1,1,2-Tetrachloroethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Ethylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	ND	0.0018	EPA 8260	12-19-11	12-19-11	
o-Xylene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Styrene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Bromoform	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Isopropylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Bromobenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.00090	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.0045	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.0045	EPA 8260	12-19-11	12-19-11	
Naphthalene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.00090	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>91</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>99</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>93</i>	<i>55-121</i>				



Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-11-0-2</b>					
Laboratory ID:	12-110-02					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Chloromethane	ND	0.0054	EPA 8260	12-19-11	12-19-11	
Vinyl Chloride	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Bromomethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Chloroethane	ND	0.0054	EPA 8260	12-19-11	12-19-11	
Trichlorofluoromethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Acetone	0.049	0.0054	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0054	EPA 8260	12-19-11	12-19-11	
Carbon Disulfide	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Methylene Chloride	ND	0.0054	EPA 8260	12-19-11	12-19-11	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Vinyl Acetate	ND	0.0054	EPA 8260	12-19-11	12-19-11	
2,2-Dichloropropane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
2-Butanone	ND	0.0054	EPA 8260	12-19-11	12-19-11	
Bromochloromethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Chloroform	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Carbon Tetrachloride	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,1-Dichloropropene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Benzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2-Dichloroethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Trichloroethene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2-Dichloropropane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Dibromomethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Bromodichloromethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
2-Chloroethyl Vinyl Ether	ND	0.0054	EPA 8260	12-19-11	12-19-11	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Methyl Isobutyl Ketone	ND	0.0054	EPA 8260	12-19-11	12-19-11	
Toluene	ND	0.0054	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-19-11	12-19-11	

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-11-0-2</b>					
Laboratory ID:	12-110-02					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Tetrachloroethene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,3-Dichloropropane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
2-Hexanone	ND	0.0054	EPA 8260	12-19-11	12-19-11	
Dibromochloromethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2-Dibromoethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Chlorobenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Ethylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	ND	0.0022	EPA 8260	12-19-11	12-19-11	
o-Xylene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Styrene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Bromoform	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Isopropylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Bromobenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.0054	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.0054	EPA 8260	12-19-11	12-19-11	
Naphthalene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>90</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>96</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>80</i>	<i>55-121</i>				

Date of Report: December 27, 2011  
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 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-11-2-3.5</b>					
Laboratory ID:	12-110-03					
Dichlorodifluoromethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Chloromethane	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Vinyl Chloride	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Bromomethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Chloroethane	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Trichlorofluoromethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Acetone	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Carbon Disulfide	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Methylene Chloride	ND	0.0044	EPA 8260	12-19-11	12-19-11	
(trans) 1,2-Dichloroethene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Methyl t-Butyl Ether	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Vinyl Acetate	ND	0.0044	EPA 8260	12-19-11	12-19-11	
2,2-Dichloropropane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
(cis) 1,2-Dichloroethene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
2-Butanone	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Bromochloromethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Chloroform	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,1,1-Trichloroethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Carbon Tetrachloride	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,1-Dichloropropene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Benzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2-Dichloroethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Trichloroethene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2-Dichloropropane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Dibromomethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Bromodichloromethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
2-Chloroethyl Vinyl Ether	ND	0.0044	EPA 8260	12-19-11	12-19-11	
(cis) 1,3-Dichloropropene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Methyl Isobutyl Ketone	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Toluene	ND	0.0044	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.00089	EPA 8260	12-19-11	12-19-11	

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 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-11-2-3.5</b>					
Laboratory ID:	12-110-03					
1,1,2-Trichloroethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Tetrachloroethene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,3-Dichloropropane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
2-Hexanone	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Dibromochloromethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2-Dibromoethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Chlorobenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,1,1,2-Tetrachloroethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Ethylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	ND	0.0018	EPA 8260	12-19-11	12-19-11	
o-Xylene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Styrene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Bromoform	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Isopropylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Bromobenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.00089	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.0044	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.0044	EPA 8260	12-19-11	12-19-11	
Naphthalene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.00089	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>94</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>100</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>95</i>	<i>55-121</i>				

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-DUPE-1</b>					
Laboratory ID:	12-110-05					
Dichlorodifluoromethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Chloromethane	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Vinyl Chloride	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Bromomethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Chloroethane	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Trichlorofluoromethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Acetone	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Carbon Disulfide	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Methylene Chloride	ND	0.0049	EPA 8260	12-19-11	12-19-11	
(trans) 1,2-Dichloroethene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Methyl t-Butyl Ether	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Vinyl Acetate	ND	0.0049	EPA 8260	12-19-11	12-19-11	
2,2-Dichloropropane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
(cis) 1,2-Dichloroethene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
2-Butanone	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Bromochloromethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Chloroform	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,1,1-Trichloroethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Carbon Tetrachloride	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,1-Dichloropropene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Benzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2-Dichloroethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Trichloroethene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2-Dichloropropane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Dibromomethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Bromodichloromethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
2-Chloroethyl Vinyl Ether	ND	0.0049	EPA 8260	12-19-11	12-19-11	
(cis) 1,3-Dichloropropene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Methyl Isobutyl Ketone	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Toluene	ND	0.0049	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.00097	EPA 8260	12-19-11	12-19-11	

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 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-DUPE-1</b>					
Laboratory ID:	12-110-05					
1,1,2-Trichloroethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Tetrachloroethene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,3-Dichloropropane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
2-Hexanone	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Dibromochloromethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2-Dibromoethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Chlorobenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,1,1,2-Tetrachloroethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Ethylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	ND	0.0019	EPA 8260	12-19-11	12-19-11	
o-Xylene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Styrene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Bromoform	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Isopropylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Bromobenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.00097	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.0049	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.0049	EPA 8260	12-19-11	12-19-11	
Naphthalene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.00097	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>90</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>97</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>94</i>	<i>55-121</i>				

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Client ID:	DP-11-0-2					
Laboratory ID:	12-110-02					
n-Nitrosodimethylamine	ND	0.19	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	1.9	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.19	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.19	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.19	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.19	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.19	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	1.9	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Naphthalene	0.026	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
4-Chloroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.19	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Hexachlorocyclopentadiene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
3-Nitroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	

Date of Report: December 27, 2011  
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**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-11-0-2</b>					
Laboratory ID:	12-110-02					
2,4-Dinitrophenol	ND	0.96	EPA 8270	12-19-11	12-20-11	
Acenaphthene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
4-Nitrophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.19	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	0.96	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.19	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.19	EPA 8270	12-19-11	12-20-11	
Fluorene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
4,6-Dinitro-2-methylphenol	ND	0.96	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.19	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.19	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.19	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.19	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	0.96	EPA 8270	12-19-11	12-20-11	
Phenanthrene	<b>0.010</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Carbazole	ND	0.19	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	1.9	EPA 8270	12-19-11	12-20-11	
Fluoranthene	<b>0.015</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzidine	ND	1.9	EPA 8270	12-19-11	12-20-11	
Pyrene	<b>0.022</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Butylbenzylphthalate	ND	1.9	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	0.96	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	1.9	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>0.011</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
bis(2-Ethylhexyl)phthalate	<b>46</b>	3.9	EPA 8270	12-19-11	12-21-11	
Di-n-octylphthalate	<b>6.7</b>	0.77	EPA 8270	12-19-11	12-21-11	
Benzo[b]fluoranthene	<b>0.0099</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>0.010</b>	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Indeno[1,2,3-cd]pyrene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	ND	0.0077	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	52	30 - 97				
Phenol-d6	54	40 - 104				
Nitrobenzene-d5	52	35 - 102				
2-Fluorobiphenyl	81	44 - 97				
2,4,6-Tribromophenol	49	41 - 110				
Terphenyl-d14	79	53 - 107				



Date of Report: December 27, 2011  
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 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-10-13-15</b>					
Laboratory ID:	12-110-01					
Naphthalene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>48</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>59</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>55</i>	<i>33 - 119</i>				

Date of Report: December 27, 2011  
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 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-11-2-3.5</b>					
Laboratory ID:	12-110-03					
Naphthalene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0077	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>49</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>52</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>46</i>	<i>33 - 119</i>				

Date of Report: December 27, 2011  
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 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-DUPE-1</b>					
Laboratory ID:	12-110-05					
Naphthalene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0075	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>40</i>	<i>43 - 109</i>				<i>Q</i>
<i>Pyrene-d10</i>	<i>47</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>45</i>	<i>33 - 119</i>				

Date of Report: December 27, 2011  
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 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-10-13-15</b>					
Laboratory ID:	12-110-01					
Aroclor 1016	ND	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1221	ND	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1232	ND	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1242	ND	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1248	ND	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1254	ND	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1260	ND	0.056	EPA 8082	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	74	42-123				
<b>Client ID:</b>	<b>DP-11-0-2</b>					
Laboratory ID:	12-110-02					
Aroclor 1016	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1221	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1232	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1242	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1248	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1254	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1260	ND	0.058	EPA 8082	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	116	42-123				
<b>Client ID:</b>	<b>DP-11-2-3.5</b>					
Laboratory ID:	12-110-03					
Aroclor 1016	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1221	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1232	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1242	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1248	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1254	ND	0.058	EPA 8082	12-21-11	12-21-11	
Aroclor 1260	ND	0.058	EPA 8082	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	59	42-123				

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**PCBs by EPA 8082**

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-DUPE-1</b>					
Laboratory ID:	12-110-05					
Aroclor 1016	<b>ND</b>	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1221	<b>ND</b>	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1232	<b>ND</b>	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1242	<b>ND</b>	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1248	<b>ND</b>	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1254	<b>ND</b>	0.056	EPA 8082	12-21-11	12-21-11	
Aroclor 1260	<b>ND</b>	0.056	EPA 8082	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>110</i>	<i>42-123</i>				

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-110-01					
<b>Client ID:</b>	<b>DP-10-13-15</b>					
Arsenic	<b>ND</b>	11	6010B	12-20-11	12-20-11	
Barium	<b>58</b>	2.8	6010B	12-20-11	12-20-11	
Cadmium	<b>ND</b>	0.56	6010B	12-20-11	12-20-11	
Chromium	<b>38</b>	0.56	6010B	12-20-11	12-20-11	
Lead	<b>ND</b>	5.6	6010B	12-20-11	12-20-11	
Mercury	<b>ND</b>	0.28	7471A	12-20-11	12-20-11	
Selenium	<b>ND</b>	11	6010B	12-20-11	12-20-11	
Silver	<b>ND</b>	0.56	6010B	12-20-11	12-20-11	

Lab ID:	12-110-02					
<b>Client ID:</b>	<b>DP-11-0-2</b>					
Arsenic	<b>ND</b>	12	6010B	12-20-11	12-20-11	
Barium	<b>82</b>	2.9	6010B	12-20-11	12-20-11	
Cadmium	<b>0.89</b>	0.58	6010B	12-20-11	12-20-11	
Chromium	<b>37</b>	0.58	6010B	12-20-11	12-20-11	
Lead	<b>26</b>	5.8	6010B	12-20-11	12-20-11	
Mercury	<b>ND</b>	0.29	7471A	12-20-11	12-20-11	
Selenium	<b>ND</b>	12	6010B	12-20-11	12-20-11	
Silver	<b>ND</b>	0.58	6010B	12-20-11	12-20-11	

Date of Report: December 27, 2011  
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 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-110-03					
Client ID:	DP-11-2-3.5					
Arsenic	ND	12	6010B	12-20-11	12-20-11	
Barium	74	2.9	6010B	12-20-11	12-20-11	
Cadmium	ND	0.58	6010B	12-20-11	12-20-11	
Chromium	45	0.58	6010B	12-20-11	12-20-11	
Lead	ND	5.8	6010B	12-20-11	12-20-11	
Mercury	ND	0.29	7471A	12-20-11	12-20-11	
Selenium	ND	12	6010B	12-20-11	12-20-11	
Silver	ND	0.58	6010B	12-20-11	12-20-11	

Lab ID:	12-110-05					
Client ID:	DP-DUPE-1					
Arsenic	ND	11	6010B	12-20-11	12-20-11	
Barium	56	2.8	6010B	12-20-11	12-20-11	
Cadmium	ND	0.56	6010B	12-20-11	12-20-11	
Chromium	33	0.56	6010B	12-20-11	12-20-11	
Lead	ND	5.6	6010B	12-20-11	12-20-11	
Mercury	ND	0.28	7471A	12-20-11	12-20-11	
Selenium	ND	11	6010B	12-20-11	12-20-11	
Silver	ND	0.56	6010B	12-20-11	12-20-11	

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1219S1					
Gasoline	ND	5.0	NWTPH-Gx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	94	68-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-111-02							
	ORIG	DUP						
Gasoline	ND	ND	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>			98	92	68-124			



Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1219S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	120	50-150				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-110-01					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			109 114	50-150		

Date of Report: December 27, 2011  
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 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloromethane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Vinyl Chloride	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromomethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloroethane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Trichlorofluoromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Acetone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Iodomethane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Carbon Disulfide	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methylene Chloride	ND	0.0050	EPA 8260	12-19-11	12-19-11	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Vinyl Acetate	ND	0.0050	EPA 8260	12-19-11	12-19-11	
2,2-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Butanone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Bromochloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chloroform	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Carbon Tetrachloride	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Benzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Trichloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Dibromomethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromodichloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	12-19-11	12-19-11	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Toluene	ND	0.0050	EPA 8260	12-19-11	12-19-11	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-19-11	12-19-11	

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Tetrachloroethene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3-Dichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Hexanone	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Dibromochloromethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dibromoethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Chlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Ethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
m,p-Xylene	ND	0.0020	EPA 8260	12-19-11	12-19-11	
o-Xylene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Styrene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromoform	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Isopropylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Bromobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	12-19-11	12-19-11	
n-Propylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
2-Chlorotoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
4-Chlorotoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
tert-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
sec-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
p-Isopropyltoluene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
n-Butylbenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	12-19-11	12-19-11	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
Hexachlorobutadiene	ND	0.0050	EPA 8260	12-19-11	12-19-11	
Naphthalene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>95</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>99</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>99</i>	<i>55-121</i>				

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1219S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0559</b>	<b>0.0550</b>	0.0500	0.0500	112	110	70-130	2	19	
Benzene	<b>0.0514</b>	<b>0.0506</b>	0.0500	0.0500	103	101	70-125	2	15	
Trichloroethene	<b>0.0471</b>	<b>0.0470</b>	0.0500	0.0500	94	94	70-122	0	14	
Toluene	<b>0.0493</b>	<b>0.0484</b>	0.0500	0.0500	99	97	73-120	2	16	
Chlorobenzene	<b>0.0455</b>	<b>0.0432</b>	0.0500	0.0500	91	86	74-109	5	12	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					89	84	63-127			
<i>Toluene-d8</i>					91	92	65-129			
<i>4-Bromofluorobenzene</i>					87	88	55-121			

Date of Report: December 27, 2011  
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**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S2					
n-Nitrosodimethylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.033	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	0.33	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Naphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4-Chloroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
3-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	

Date of Report: December 27, 2011  
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**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S2					
2,4-Dinitrophenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4-Nitrophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	0.17	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Carbazole	ND	0.033	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	0.33	EPA 8270	12-19-11	12-20-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzidine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Butylbenzylphthalate	ND	0.33	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	0.17	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
bis(2-Ethylhexyl)phthalate	ND	0.17	EPA 8270	12-19-11	12-20-11	
Di-n-octylphthalate	ND	0.033	EPA 8270	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>85</i>	<i>30 - 97</i>				
<i>Phenol-d6</i>	<i>80</i>	<i>40 - 104</i>				
<i>Nitrobenzene-d5</i>	<i>74</i>	<i>35 - 102</i>				
<i>2-Fluorobiphenyl</i>	<i>92</i>	<i>44 - 97</i>				
<i>2,4,6-Tribromophenol</i>	<i>103</i>	<i>41 - 110</i>				
<i>Terphenyl-d14</i>	<i>79</i>	<i>53 - 107</i>				

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					Recovery	Limits	Limit			
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1219S2									
	SB	SBD	SB	SBD	SB	SBD				
Phenol	<b>0.925</b>	<b>0.950</b>	1.33	1.33	70	71	31 - 111	3	34	
2-Chlorophenol	<b>1.04</b>	<b>1.06</b>	1.33	1.33	78	80	29 - 112	2	37	
1,4-Dichlorobenzene	<b>0.522</b>	<b>0.528</b>	0.667	0.667	78	79	24 - 100	1	37	
n-Nitroso-di-n-propylamine	<b>0.396</b>	<b>0.405</b>	0.667	0.667	59	61	35 - 104	2	32	
1,2,4-Trichlorobenzene	<b>0.685</b>	<b>0.713</b>	0.667	0.667	103	107	29 - 110	4	35	
4-Chloro-3-methylphenol	<b>1.13</b>	<b>1.19</b>	1.33	1.33	85	89	53 - 104	5	25	
Acenaphthene	<b>0.544</b>	<b>0.563</b>	0.667	0.667	82	84	50 - 95	3	23	
4-Nitrophenol	<b>1.22</b>	<b>1.30</b>	1.33	1.33	92	98	42 - 126	6	30	
2,4-Dinitrotoluene	<b>0.722</b>	<b>0.749</b>	0.667	0.667	108	112	53 - 115	4	31	
Pentachlorophenol	<b>0.939</b>	<b>1.03</b>	1.33	1.33	71	77	50 - 116	9	30	
Pyrene	<b>0.446</b>	<b>0.468</b>	0.667	0.667	67	70	57 - 120	5	27	
<i>Surrogate:</i>										
2-Fluorophenol					83	82	30 - 97			
Phenol-d6					76	78	40 - 104			
Nitrobenzene-d5					72	74	35 - 102			
2-Fluorobiphenyl					91	91	44 - 97			
2,4,6-Tribromophenol					105	111	41 - 110			
Terphenyl-d14					78	82	53 - 107			

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
Naphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>75</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>85</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>77</i>	<i>33 - 119</i>				



Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent		Recovery	RPD	RPD	Flags
	MS	MSD	MS	MSD	Result	Recovery	Limits	RPD	Limit		
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-110-03										
	MS	MSD	MS	MSD		MS	MSD				
Naphthalene	<b>0.0546</b>	<b>0.0593</b>	0.0833	0.0833	ND	66	71	39 - 110	8	21	
Acenaphthylene	<b>0.0530</b>	<b>0.0593</b>	0.0833	0.0833	ND	64	71	47 - 124	11	21	
Acenaphthene	<b>0.0489</b>	<b>0.0560</b>	0.0833	0.0833	ND	59	67	50 - 120	14	20	
Fluorene	<b>0.0487</b>	<b>0.0568</b>	0.0833	0.0833	ND	58	68	52 - 126	15	21	
Phenanthrene	<b>0.0478</b>	<b>0.0551</b>	0.0833	0.0833	ND	57	66	41 - 130	14	22	
Anthracene	<b>0.0473</b>	<b>0.0554</b>	0.0833	0.0833	ND	57	67	48 - 124	16	23	
Fluoranthene	<b>0.0495</b>	<b>0.0581</b>	0.0833	0.0833	ND	59	70	40 - 137	16	23	
Pyrene	<b>0.0487</b>	<b>0.0579</b>	0.0833	0.0833	ND	58	70	36 - 139	17	23	
Benzo[a]anthracene	<b>0.0496</b>	<b>0.0593</b>	0.0833	0.0833	ND	60	71	43 - 127	18	21	
Chrysene	<b>0.0476</b>	<b>0.0570</b>	0.0833	0.0833	ND	57	68	41 - 133	18	19	
Benzo[b]fluoranthene	<b>0.0452</b>	<b>0.0555</b>	0.0833	0.0833	ND	54	67	40 - 132	20	25	
Benzo(j,k)fluoranthene	<b>0.0436</b>	<b>0.0526</b>	0.0833	0.0833	ND	52	63	35 - 132	19	25	
Benzo[a]pyrene	<b>0.0471</b>	<b>0.0563</b>	0.0833	0.0833	ND	57	68	37 - 131	18	26	
Indeno(1,2,3-c,d)pyrene	<b>0.0444</b>	<b>0.0541</b>	0.0833	0.0833	ND	53	65	39 - 134	20	23	
Dibenz[a,h]anthracene	<b>0.0457</b>	<b>0.0553</b>	0.0833	0.0833	ND	55	66	40 - 137	19	21	
Benzo[g,h,i]perylene	<b>0.0456</b>	<b>0.0550</b>	0.0833	0.0833	ND	55	66	35 - 135	19	22	
<i>Surrogate:</i>											
<i>2-Fluorobiphenyl</i>						<i>54</i>	<i>60</i>	<i>43 - 109</i>			
<i>Pyrene-d10</i>						<i>56</i>	<i>67</i>	<i>38 - 128</i>			
<i>Terphenyl-d14</i>						<i>53</i>	<i>63</i>	<i>33 - 119</i>			

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1221S1					
Aroclor 1016	ND	0.050	EPA 8082	12-21-11	12-21-11	
Aroclor 1221	ND	0.050	EPA 8082	12-21-11	12-21-11	
Aroclor 1232	ND	0.050	EPA 8082	12-21-11	12-21-11	
Aroclor 1242	ND	0.050	EPA 8082	12-21-11	12-21-11	
Aroclor 1248	ND	0.050	EPA 8082	12-21-11	12-21-11	
Aroclor 1254	ND	0.050	EPA 8082	12-21-11	12-21-11	
Aroclor 1260	ND	0.050	EPA 8082	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	88		42-123			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-110-05										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.237	0.281	0.500	0.500	ND	47	56	44-125	17	20	
<i>Surrogate:</i>											
DCB						47	53	42-123			

Date of Report: December 27, 2011  
Samples Submitted: December 15, 2011  
Laboratory Reference: 1112-110  
Project: 0180-292-00

**TOTAL METALS  
EPA 6010B/7471A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-20-11  
Date Analyzed: 12-20-11  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB1220SM1&MB1220S1

Analyte	Method	Result	PQL
Arsenic	6010B	ND	10
Barium	6010B	ND	2.5
Cadmium	6010B	ND	0.50
Chromium	6010B	ND	0.50
Lead	6010B	ND	5.0
Mercury	7471A	ND	0.25
Selenium	6010B	ND	10
Silver	6010B	ND	0.50

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-20-11  
 Date Analyzed: 12-20-11  
  
 Matrix: Soil  
 Units: mg/kg (ppm)  
  
 Lab ID: 12-110-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	52.1	53.9	4	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	34.4	31.0	10	0.50	
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	

Date of Report: December 27, 2011  
 Samples Submitted: December 15, 2011  
 Laboratory Reference: 1112-110  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-20-11

Date Analyzed: 12-20-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 12-110-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>88.6</b>	89	<b>87.2</b>	87	2	
Barium	100	<b>149</b>	97	<b>149</b>	97	0	
Cadmium	50.0	<b>46.7</b>	93	<b>46.0</b>	92	1	
Chromium	100	<b>127</b>	93	<b>123</b>	89	3	
Lead	250	<b>227</b>	91	<b>225</b>	90	1	
Mercury	0.500	<b>0.492</b>	98	<b>0.492</b>	98	0	
Selenium	100	<b>92.3</b>	92	<b>91.3</b>	91	1	
Silver	25.0	<b>20.4</b>	82	<b>20.1</b>	80	2	

Date of Report: December 27, 2011  
Samples Submitted: December 15, 2011  
Laboratory Reference: 1112-110  
Project: 0180-292-00

**% MOISTURE**

Date Analyzed: 12-16-11

Client ID	Lab ID	% Moisture
DP-10-13-15	12-110-01	11
DP-11-0-2	12-110-02	14
DP-11-2-3.5	12-110-03	14
DP-DUPE-1	12-110-05	11



### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference



# Onsite Environmental Inc.

Analytical Laboratory Testing Services  
14648 NE 95th Street • Redmond, WA 98052  
Phone: (425) 883-3881 • www.onsite-env.com

## Chain of Custody

12-110

### Turnaround Request (in working days)

(Check One)

Same Day  1 Day

2 Days  3 Days

Standard (7 Days) (TPH analysis 5 Days)

\_\_\_\_\_ (other)

### Laboratory Number:

Company: Geo Engineers  
 Project Number: 0180-292-00  
 Project Name: WSDOT Midway Metals  
 Project Manager: Aaron Weigman  
 Sampled by: [Signature]

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	No. of Cont.
1	DP-10-13-15	12/14/11	1400	Soil	5
2	DP-11-0-2		1600		
3	DP-11-2-3,5		1610		
4	DP-11-6-8		1630		
5	DP-DUPR-1		0800		

Method	1	2	3	4	5
NWTPH-HCID					
NWTPH-Gx/BTEX					
NWTPH-Gx	X	X	X	X	X
NWTPH-Dx	X	X	X	X	X
Volatiles 8260B	X	X	X	X	X
Halogenated Volatiles 8260B					
Semivolatiles 8270D/SIM (with low-level PAHs)	X	X	X	X	X
PAHs 8270D/SIM (low-level)	X	X	X	X	X
PCBs 8082	X	X	X	X	X
Organochlorine Pesticides 8081A					
Organophosphorus Pesticides 8270D/SIM					
Chlorinated Acid Herbicides 8151A					
Total RCRA Metals	X	X	X	X	X
Total MTCA Metals					
TCLP Metals					
HEM (oil and grease) 1664					
% Moisture	X	X	X	X	X

Signature	Company	Date	Time	Comments/Special Instructions
[Signature]	Geo Engineers	12/15/11	1120	
[Signature]	Speery Hsing	12-15-11	1120	
[Signature]	"	"	1414	
[Signature]	OBE	12/15/11	1414	

Relinquished  
Received  
Relinquished  
Received  
Relinquished  
Received  
Reviewed/Date

Reviewed/Date

Chromatograms with final report





14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

February 16, 2012

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1202-085

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on February 9, 2012.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister  
Project Manager

Enclosures

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

### **Case Narrative**

Samples were collected on February 8 and 9, 2012 and received by the laboratory on February 9, 2012. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### NWTPH Gx and Volatiles EPA 8260B Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**ANALYTICAL REPORT FOR SAMPLES**

<b>Client ID</b>	<b>Laboratory ID</b>	<b>Matrix</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Notes</b>
DP12-0.0-2.0	02-085-01	Soil	2-8-12	2-9-12	
DP13-0.0-1.5	02-085-02	Soil	2-8-12	2-9-12	
DP14-0.0-2.0	02-085-03	Soil	2-8-12	2-9-12	
DP15-0.0-2.0	02-085-04	Soil	2-8-12	2-9-12	
DP16-0.0-2.0	02-085-05	Soil	2-8-12	2-9-12	
DP17-0.0-2.0	02-085-06	Soil	2-8-12	2-9-12	
DP18-0.0-2.0	02-085-07	Soil	2-8-12	2-9-12	
DP19-0.0-2.0	02-085-08	Soil	2-8-12	2-9-12	
DP20-0.0-2.0	02-085-09	Soil	2-8-12	2-9-12	
DP21-0.0-2.0	02-085-10	Soil	2-8-12	2-9-12	
DP12-5.0-7.5	02-085-12	Soil	2-8-12	2-9-12	
DP15-5.0-7.0	02-085-19	Soil	2-8-12	2-9-12	
DP21-2.0-3.0	02-085-27	Soil	2-8-12	2-9-12	
HA1-0.0-2.0	02-085-28	Soil	2-9-12	2-9-12	
HA2-0.0-2.0	02-085-29	Soil	2-9-12	2-9-12	
HA3-0.0-2.0	02-085-30	Soil	2-9-12	2-9-12	
HA4-0.0-2.0	02-085-31	Soil	2-9-12	2-9-12	
SED9	02-085-32	Soil	2-9-12	2-9-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
Laboratory ID:	02-085-01					
Diesel Range Organics	<b>ND</b>	28	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil	<b>130</b>	55	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	102	50-150				
<b>Client ID:</b>	<b>DP13-0.0-1.5</b>					
Laboratory ID:	02-085-02					
Diesel Range Organics	<b>ND</b>	31	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	63	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	101	50-150				
<b>Client ID:</b>	<b>DP14-0.0-2.0</b>					
Laboratory ID:	02-085-03					
Diesel Range Organics	<b>ND</b>	28	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	56	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Laboratory ID:	02-085-04					
Diesel Range Organics	<b>ND</b>	30	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	60	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	92	50-150				
<b>Client ID:</b>	<b>DP16-0.0-2.0</b>					
Laboratory ID:	02-085-05					
Diesel Range Organics	<b>ND</b>	29	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	58	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	105	50-150				
<b>Client ID:</b>	<b>DP17-0.0-2.0</b>					
Laboratory ID:	02-085-06					
Diesel Range Organics	<b>ND</b>	29	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	58	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP21-0.0-2.0</b>					
Laboratory ID:	02-085-10					
Diesel Fuel #2	<b>100</b>	29	NWTPH-Dx	2-10-12	2-10-12	N
Lube Oil	<b>470</b>	59	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	106	50-150				
<b>Client ID:</b>	<b>DP12-5.0-7.5</b>					
Laboratory ID:	02-085-12					
Diesel Range Organics	<b>ND</b>	30	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	59	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	104	50-150				
<b>Client ID:</b>	<b>DP15-5.0-7.0</b>					
Laboratory ID:	02-085-19					
Diesel Range Organics	<b>ND</b>	32	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	65	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	97	50-150				
<b>Client ID:</b>	<b>DP21-2.0-3.0</b>					
Laboratory ID:	02-085-27					
Diesel Range Organics	<b>ND</b>	32	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	63	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	100	50-150				
<b>Client ID:</b>	<b>HA1-0.0-2.0</b>					
Laboratory ID:	02-085-28					
Diesel Range Organics	<b>ND</b>	31	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	62	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	102	50-150				
<b>Client ID:</b>	<b>HA2-0.0-2.0</b>					
Laboratory ID:	02-085-29					
Diesel Range Organics	<b>ND</b>	34	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	67	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	95	50-150				

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**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>HA3-0.0-2.0</b>					
Laboratory ID:	02-085-30					
Diesel Range Organics	<b>ND</b>	110	NWTPH-Dx	2-10-12	2-10-12	U1
Lube Oil	<b>660</b>	60	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				
<b>Client ID:</b>	<b>SED9</b>					
Laboratory ID:	02-085-32					
Diesel Range Organics	<b>ND</b>	130	NWTPH-Dx	2-10-12	2-10-12	U1
Lube Oil	<b>1000</b>	130	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	97	50-150				

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### NWTPH-Gx

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
Laboratory ID:	02-085-01					
Gasoline	<b>ND</b>	5.3	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	68-124				
<b>Client ID:</b>	<b>DP13-0.0-1.5</b>					
Laboratory ID:	02-085-02					
Gasoline	<b>ND</b>	6.1	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	101	68-124				
<b>Client ID:</b>	<b>DP14-0.0-2.0</b>					
Laboratory ID:	02-085-03					
Gasoline	<b>ND</b>	5.7	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	68-124				
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Laboratory ID:	02-085-04					
Gasoline	<b>ND</b>	7.2	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	87	68-124				
<b>Client ID:</b>	<b>DP16-0.0-2.0</b>					
Laboratory ID:	02-085-05					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	95	68-124				
<b>Client ID:</b>	<b>DP17-0.0-2.0</b>					
Laboratory ID:	02-085-06					
Gasoline	<b>ND</b>	5.2	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	97	68-124				

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### NWTPH-Gx

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP18-0.0-2.0</b>					
Laboratory ID:	02-085-07					
Gasoline	<b>ND</b>	5.4	NWTPH-Gx	2-13-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	95	68-124				
<b>Client ID:</b>	<b>DP19-0.0-2.0</b>					
Laboratory ID:	02-085-08					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	2-13-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	68-124				
<b>Client ID:</b>	<b>DP20-0.0-2.0</b>					
Laboratory ID:	02-085-09					
Gasoline	<b>ND</b>	5.5	NWTPH-Gx	2-13-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	94	68-124				
<b>Client ID:</b>	<b>DP12-5.0-7.5</b>					
Laboratory ID:	02-085-12					
Gasoline	<b>ND</b>	4.9	NWTPH-Gx	2-13-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	88	68-124				
<b>Client ID:</b>	<b>DP15-5.0-7.0</b>					
Laboratory ID:	02-085-19					
Gasoline	<b>ND</b>	6.6	NWTPH-Gx	2-13-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	108	68-124				



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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
Laboratory ID:	02-085-01					
Dichlorodifluoromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0041	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0041	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0041	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0041	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
Laboratory ID:	02-085-01					
1,1,2-Trichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0017	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>112</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>107</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>83</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP13-0.0-1.5</b>					
Laboratory ID:	02-085-02					
Dichlorodifluoromethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0048	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0048	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Acetone	0.023	0.0048	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0048	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0048	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0048	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0048	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0048	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0048	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0048	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.00097	EPA 8260	2-13-12	2-13-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP13-0.0-1.5</b>					
Laboratory ID:	02-085-02					
1,1,2-Trichloroethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0048	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0019	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.00097	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.063	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.063	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.31	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.31	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.063	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.063	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>110</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>108</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>80</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP14-0.0-2.0</b>					
<b>Laboratory ID:</b>	02-085-03					
Dichlorodifluoromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0041	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0041	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0041	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0041	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP14-0.0-2.0</b>					
Laboratory ID:	02-085-03					
1,1,2-Trichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0017	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0041	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0041	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>113</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>112</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>89</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Laboratory ID:	02-085-04					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Acetone	0.10	0.0054	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0054	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0054	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0054	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0054	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	2-13-12	2-13-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Laboratory ID:	02-085-04					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0021	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.055	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.055	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.27	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.27	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.055	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.055	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>108</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>105</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>75</i>	<i>55-121</i>				



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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP16-0.0-2.0</b>					
Laboratory ID:	02-085-05					
Dichlorodifluoromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0044	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0044	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0044	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0044	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.00088	EPA 8260	2-13-12	2-13-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP16-0.0-2.0</b>					
Laboratory ID:	02-085-05					
1,1,2-Trichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0018	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>108</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>106</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>89</i>	<i>55-121</i>				

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**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP17-0.0-2.0</b>					
Laboratory ID:	02-085-06					
Dichlorodifluoromethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0047	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0047	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0047	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0047	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.00094	EPA 8260	2-13-12	2-13-12	

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 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP17-0.0-2.0</b>					
Laboratory ID:	02-085-06					
1,1,2-Trichloroethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0019	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.00094	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0047	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0047	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.00094	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>111</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>110</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>89</i>	<i>55-121</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
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 Project: 0180-292-00

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP18-0.0-2.0</b>					
Laboratory ID:	02-085-07					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Chloromethane	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Vinyl Chloride	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Bromomethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Chloroethane	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Trichlorofluoromethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,1-Dichloroethene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Acetone	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Iodomethane	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Carbon Disulfide	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Methylene Chloride	ND	0.0053	EPA 8260	2-14-12	2-14-12	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,1-Dichloroethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Vinyl Acetate	ND	0.0053	EPA 8260	2-14-12	2-14-12	
2,2-Dichloropropane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
2-Butanone	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Bromochloromethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Chloroform	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Carbon Tetrachloride	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,1-Dichloropropene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Benzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2-Dichloroethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Trichloroethene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2-Dichloropropane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Dibromomethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Bromodichloromethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
2-Chloroethyl Vinyl Ether	ND	0.0053	EPA 8260	2-14-12	2-14-12	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Methyl Isobutyl Ketone	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Toluene	ND	0.0053	EPA 8260	2-14-12	2-14-12	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	2-14-12	2-14-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP18-0.0-2.0</b>					
Laboratory ID:	02-085-07					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Tetrachloroethene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,3-Dichloropropane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
2-Hexanone	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Dibromochloromethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2-Dibromoethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Chlorobenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Ethylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
m,p-Xylene	ND	0.0021	EPA 8260	2-14-12	2-14-12	
o-Xylene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Styrene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Bromoform	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Isopropylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Bromobenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260	2-14-12	2-14-12	
n-Propylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
2-Chlorotoluene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
4-Chlorotoluene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
tert-Butylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
sec-Butylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
p-Isopropyltoluene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
n-Butylbenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2-Dibromo-3-chloropropane	ND	0.0053	EPA 8260	2-14-12	2-14-12	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
Hexachlorobutadiene	ND	0.0053	EPA 8260	2-14-12	2-14-12	
Naphthalene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>109</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>108</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>87</i>	<i>55-121</i>				

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 Project: 0180-292-00

**VOLATILES by EPA 8260B**

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP19-0.0-2.0</b>					
<b>Laboratory ID:</b>	02-085-08					
Dichlorodifluoromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	0.0026	0.00088	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0044	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0044	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0044	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	0.0066	0.0044	EPA 8260	2-13-12	2-13-12	
Toluene	0.0066	0.0044	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.00088	EPA 8260	2-13-12	2-13-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP19-0.0-2.0</b>					
Laboratory ID:	02-085-08					
1,1,2-Trichloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	0.0035	0.00088	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	0.021	0.0018	EPA 8260	2-13-12	2-13-12	
o-Xylene	0.013	0.00088	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.00088	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	0.0025	0.00088	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	0.013	0.00088	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	0.021	0.00088	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	0.0016	0.00088	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0044	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0044	EPA 8260	2-13-12	2-13-12	
Naphthalene	0.0020	0.00088	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.00088	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>115</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>109</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>89</i>	<i>55-121</i>				



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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP20-0.0-2.0</b>					
Laboratory ID:	02-085-09					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	0.0016	0.0010	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0051	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0051	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0051	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0051	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	2-13-12	2-13-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP20-0.0-2.0</b>					
Laboratory ID:	02-085-09					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0020	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0051	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0051	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>108</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>108</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>87</i>	<i>55-121</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP12-5.0-7.5</b>					
Laboratory ID:	02-085-12					
Dichlorodifluoromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0042	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0042	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Acetone	0.013	0.0042	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0042	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0042	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0042	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0042	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0042	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0042	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0042	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.00083	EPA 8260	2-13-12	2-13-12	

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 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-5.0-7.5</b>					
Laboratory ID:	02-085-12					
1,1,2-Trichloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0042	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0017	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.00083	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0042	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0042	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.00083	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>105</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>105</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>89</i>	<i>55-121</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP15-5.0-7.0</b>					
Laboratory ID:	02-085-19					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0054	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0054	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0054	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0054	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	2-13-12	2-13-12	

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 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP15-5.0-7.0</b>					
Laboratory ID:	02-085-19					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0021	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0054	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0054	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>110</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>113</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>90</i>	<i>55-121</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
<b>Laboratory ID:</b>	<b>02-085-01</b>					
n-Nitrosodimethylamine	ND	0.037	EPA 8270	2-15-12	2-15-12	
Pyridine	ND	0.37	EPA 8270	2-15-12	2-15-12	
Phenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
Aniline	ND	0.037	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroethyl)ether	ND	0.037	EPA 8270	2-15-12	2-15-12	
2-Chlorophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
1,3-Dichlorobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
1,4-Dichlorobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
Benzyl alcohol	ND	0.037	EPA 8270	2-15-12	2-15-12	
1,2-Dichlorobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
2-Methylphenol (o-Cresol)	ND	0.037	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroisopropyl)ether	ND	0.037	EPA 8270	2-15-12	2-15-12	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.037	EPA 8270	2-15-12	2-15-12	
n-Nitroso-di-n-propylamine	ND	0.037	EPA 8270	2-15-12	2-15-12	
Hexachloroethane	ND	0.037	EPA 8270	2-15-12	2-15-12	
Nitrobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
Isophorone	ND	0.037	EPA 8270	2-15-12	2-15-12	
2-Nitrophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,4-Dimethylphenol	ND	0.37	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroethoxy)methane	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,4-Dichlorophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
1,2,4-Trichlorobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
Naphthalene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
4-Chloroaniline	ND	0.037	EPA 8270	2-15-12	2-15-12	
Hexachlorobutadiene	ND	0.037	EPA 8270	2-15-12	2-15-12	
4-Chloro-3-methylphenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
2-Methylnaphthalene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
1-Methylnaphthalene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Hexachlorocyclopentadiene	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,4,6-Trichlorophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,3-Dichloroaniline	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,4,5-Trichlorophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
2-Chloronaphthalene	ND	0.037	EPA 8270	2-15-12	2-15-12	
2-Nitroaniline	ND	0.037	EPA 8270	2-15-12	2-15-12	
1,4-Dinitrobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
Dimethylphthalate	0.12	0.037	EPA 8270	2-15-12	2-15-12	
1,3-Dinitrobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,6-Dinitrotoluene	ND	0.037	EPA 8270	2-15-12	2-15-12	
1,2-Dinitrobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
Acenaphthylene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
3-Nitroaniline	ND	0.037	EPA 8270	2-15-12	2-15-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
Laboratory ID:	02-085-01					
2,4-Dinitrophenol	ND	0.18	EPA 8270	2-15-12	2-15-12	
Acenaphthene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
4-Nitrophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,4-Dinitrotoluene	ND	0.037	EPA 8270	2-15-12	2-15-12	
Dibenzofuran	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,3,5,6-Tetrachlorophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
2,3,4,6-Tetrachlorophenol	ND	0.037	EPA 8270	2-15-12	2-15-12	
Diethylphthalate	ND	0.18	EPA 8270	2-15-12	2-15-12	
4-Chlorophenyl-phenylether	ND	0.037	EPA 8270	2-15-12	2-15-12	
4-Nitroaniline	ND	0.037	EPA 8270	2-15-12	2-15-12	
Fluorene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
4,6-Dinitro-2-methylphenol	ND	0.18	EPA 8270	2-15-12	2-15-12	
n-Nitrosodiphenylamine	ND	0.037	EPA 8270	2-15-12	2-15-12	
1,2-Diphenylhydrazine	ND	0.037	EPA 8270	2-15-12	2-15-12	
4-Bromophenyl-phenylether	ND	0.037	EPA 8270	2-15-12	2-15-12	
Hexachlorobenzene	ND	0.037	EPA 8270	2-15-12	2-15-12	
Pentachlorophenol	ND	0.18	EPA 8270	2-15-12	2-15-12	
Phenanthrene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Anthracene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Carbazole	ND	0.037	EPA 8270	2-15-12	2-15-12	
Di-n-butylphthalate	ND	0.37	EPA 8270	2-15-12	2-15-12	
Fluoranthene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Benzidine	ND	0.37	EPA 8270	2-15-12	2-15-12	
Pyrene	0.011	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Butylbenzylphthalate	ND	0.37	EPA 8270	2-15-12	2-15-12	
bis-2-Ethylhexyladipate	ND	0.18	EPA 8270	2-15-12	2-15-12	
3,3'-Dichlorobenzidine	ND	0.37	EPA 8270	2-15-12	2-15-12	
Benzo[a]anthracene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Chrysene	0.0096	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
bis(2-Ethylhexyl)phthalate	0.23	0.18	EPA 8270	2-15-12	2-15-12	
Di-n-octylphthalate	ND	0.037	EPA 8270	2-15-12	2-15-12	
Benzo[b]fluoranthene	0.011	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Benzo(j,k)fluoranthene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Benzo[a]pyrene	0.021	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Indeno[1,2,3-cd]pyrene	0.010	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Dibenz[a,h]anthracene	ND	0.0074	EPA 8270/SIM	2-15-12	2-15-12	
Benzo[g,h,i]perylene	0.046	0.037	EPA 8270	2-15-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	56	30 - 97				
Phenol-d6	68	40 - 104				
Nitrobenzene-d5	63	35 - 102				
2-Fluorobiphenyl	81	44 - 97				
2,4,6-Tribromophenol	96	41 - 110				
Terphenyl-d14	82	53 - 107				



Date of Report: February 16, 2012  
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 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Laboratory ID:	02-085-04					
n-Nitrosodimethylamine	ND	0.040	EPA 8270	2-15-12	2-15-12	
Pyridine	ND	0.40	EPA 8270	2-15-12	2-15-12	
Phenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
Aniline	ND	0.040	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroethyl)ether	ND	0.040	EPA 8270	2-15-12	2-15-12	
2-Chlorophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,3-Dichlorobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,4-Dichlorobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
Benzyl alcohol	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,2-Dichlorobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
2-Methylphenol (o-Cresol)	ND	0.040	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroisopropyl)ether	ND	0.040	EPA 8270	2-15-12	2-15-12	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.040	EPA 8270	2-15-12	2-15-12	
n-Nitroso-di-n-propylamine	ND	0.040	EPA 8270	2-15-12	2-15-12	
Hexachloroethane	ND	0.040	EPA 8270	2-15-12	2-15-12	
Nitrobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
Isophorone	ND	0.040	EPA 8270	2-15-12	2-15-12	
2-Nitrophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,4-Dimethylphenol	ND	0.40	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroethoxy)methane	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,4-Dichlorophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,2,4-Trichlorobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
Naphthalene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
4-Chloroaniline	ND	0.040	EPA 8270	2-15-12	2-15-12	
Hexachlorobutadiene	ND	0.040	EPA 8270	2-15-12	2-15-12	
4-Chloro-3-methylphenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
2-Methylnaphthalene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
1-Methylnaphthalene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Hexachlorocyclopentadiene	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,4,6-Trichlorophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,3-Dichloroaniline	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,4,5-Trichlorophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
2-Chloronaphthalene	ND	0.040	EPA 8270	2-15-12	2-15-12	
2-Nitroaniline	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,4-Dinitrobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
Dimethylphthalate	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,3-Dinitrobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,6-Dinitrotoluene	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,2-Dinitrobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
Acenaphthylene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
3-Nitroaniline	ND	0.040	EPA 8270	2-15-12	2-15-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Laboratory ID:	02-085-04					
2,4-Dinitrophenol	ND	0.20	EPA 8270	2-15-12	2-15-12	
Acenaphthene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
4-Nitrophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,4-Dinitrotoluene	ND	0.040	EPA 8270	2-15-12	2-15-12	
Dibenzofuran	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,3,5,6-Tetrachlorophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
2,3,4,6-Tetrachlorophenol	ND	0.040	EPA 8270	2-15-12	2-15-12	
Diethylphthalate	ND	0.20	EPA 8270	2-15-12	2-15-12	
4-Chlorophenyl-phenylether	ND	0.040	EPA 8270	2-15-12	2-15-12	
4-Nitroaniline	ND	0.040	EPA 8270	2-15-12	2-15-12	
Fluorene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
4,6-Dinitro-2-methylphenol	ND	0.20	EPA 8270	2-15-12	2-15-12	
n-Nitrosodiphenylamine	ND	0.040	EPA 8270	2-15-12	2-15-12	
1,2-Diphenylhydrazine	ND	0.040	EPA 8270	2-15-12	2-15-12	
4-Bromophenyl-phenylether	ND	0.040	EPA 8270	2-15-12	2-15-12	
Hexachlorobenzene	ND	0.040	EPA 8270	2-15-12	2-15-12	
Pentachlorophenol	ND	0.20	EPA 8270	2-15-12	2-15-12	
Phenanthrene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Anthracene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Carbazole	ND	0.040	EPA 8270	2-15-12	2-15-12	
Di-n-butylphthalate	ND	0.40	EPA 8270	2-15-12	2-15-12	
Fluoranthene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Benzidine	ND	0.40	EPA 8270	2-15-12	2-15-12	
Pyrene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Butylbenzylphthalate	ND	0.40	EPA 8270	2-15-12	2-15-12	
bis-2-Ethylhexyladipate	ND	0.20	EPA 8270	2-15-12	2-15-12	
3,3'-Dichlorobenzidine	ND	0.40	EPA 8270	2-15-12	2-15-12	
Benzo[a]anthracene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Chrysene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
bis(2-Ethylhexyl)phthalate	ND	0.20	EPA 8270	2-15-12	2-15-12	
Di-n-octylphthalate	ND	0.040	EPA 8270	2-15-12	2-15-12	
Benzo[b]fluoranthene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Benzo(j,k)fluoranthene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Benzo[a]pyrene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Indeno[1,2,3-cd]pyrene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Dibenz[a,h]anthracene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
Benzo[g,h,i]perylene	ND	0.0079	EPA 8270/SIM	2-15-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	55	30 - 97				
Phenol-d6	65	40 - 104				
Nitrobenzene-d5	60	35 - 102				
2-Fluorobiphenyl	74	44 - 97				
2,4,6-Tribromophenol	81	41 - 110				
Terphenyl-d14	72	53 - 107				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP13-0.0-1.5</b>					
Laboratory ID:	02-085-02					
Naphthalene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	<b>0.011</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	<b>0.017</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	<b>0.016</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	<b>ND</b>	0.0084	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>71</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>73</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>68</i>	<i>33 - 119</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP14-0.0-2.0</b>					
Laboratory ID:	02-085-03					
Naphthalene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	ND	0.0075	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>71</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>85</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>80</i>	<i>33 - 119</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP16-0.0-2.0</b>					
Laboratory ID:	02-085-05					
Naphthalene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	69	43 - 109				
<i>Pyrene-d10</i>	73	38 - 128				
<i>Terphenyl-d14</i>	70	33 - 119				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP17-0.0-2.0</b>					
Laboratory ID:	02-085-06					
Naphthalene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	ND	0.0077	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	63	43 - 109				
<i>Pyrene-d10</i>	65	38 - 128				
<i>Terphenyl-d14</i>	60	33 - 119				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP12-5.0-7.5</b>					
Laboratory ID:	02-085-12					
Naphthalene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	<b>ND</b>	0.0079	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>77</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>80</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>77</i>	<i>33 - 119</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP15-5.0-7.0</b>					
Laboratory ID:	02-085-19					
Naphthalene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	<b>ND</b>	0.0086	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>58</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>67</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>61</i>	<i>33 - 119</i>				



Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED9</b>					
Laboratory ID:	02-085-32					
Naphthalene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	<b>0.017</b>	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	ND	0.017	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>54</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>67</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>58</i>	<i>33 - 119</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
Laboratory ID:	02-085-01					
Aroclor 1016	ND	0.055	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	ND	0.055	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	ND	0.055	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	ND	0.055	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	ND	0.055	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	ND	0.055	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	ND	0.055	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	86	42-123				
<b>Client ID:</b>	<b>DP13-0.0-1.5</b>					
Laboratory ID:	02-085-02					
Aroclor 1016	ND	0.063	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	ND	0.063	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	ND	0.063	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	ND	0.063	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	ND	0.063	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	ND	0.063	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	ND	0.063	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	86	42-123				
<b>Client ID:</b>	<b>DP14-0.0-2.0</b>					
Laboratory ID:	02-085-03					
Aroclor 1016	ND	0.056	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	ND	0.056	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	ND	0.056	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	ND	0.056	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	ND	0.056	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	ND	0.056	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	ND	0.056	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	84	42-123				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
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 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Laboratory ID:	02-085-04					
Aroclor 1016	ND	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	ND	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	ND	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	ND	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	ND	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	ND	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	ND	0.059	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	89	42-123				
<b>Client ID:</b>	<b>DP16-0.0-2.0</b>					
Laboratory ID:	02-085-05					
Aroclor 1016	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	ND	0.058	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	91	42-123				
<b>Client ID:</b>	<b>DP17-0.0-2.0</b>					
Laboratory ID:	02-085-06					
Aroclor 1016	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	ND	0.058	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	ND	0.058	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	78	42-123				

Date of Report: February 16, 2012  
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 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP12-5.0-7.5</b>					
Laboratory ID:	02-085-12					
Aroclor 1016	<b>ND</b>	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	<b>ND</b>	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	<b>ND</b>	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	<b>ND</b>	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	<b>ND</b>	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	<b>ND</b>	0.059	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	<b>ND</b>	0.059	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	79	42-123				
<b>Client ID:</b>	<b>DP15-5.0-7.0</b>					
Laboratory ID:	02-085-19					
Aroclor 1016	<b>ND</b>	0.065	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	<b>ND</b>	0.065	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	<b>ND</b>	0.065	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	<b>ND</b>	0.065	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	<b>ND</b>	0.065	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	<b>ND</b>	0.065	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	<b>ND</b>	0.065	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	74	42-123				

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 Project: 0180-292-00

**TOTAL METALS**  
**EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-01					
<b>Client ID:</b>	<b>DP12-0.0-2.0</b>					
Arsenic	ND	11	6010B	2-14-12	2-14-12	
Barium	75	2.8	6010B	2-14-12	2-14-12	
Cadmium	ND	0.55	6010B	2-14-12	2-14-12	
Chromium	34	0.55	6010B	2-14-12	2-14-12	
Lead	19	5.5	6010B	2-14-12	2-14-12	
Mercury	ND	0.28	7471A	2-14-12	2-14-12	
Selenium	ND	11	6010B	2-14-12	2-14-12	
Silver	ND	0.55	6010B	2-14-12	2-14-12	

Lab ID:	02-085-02					
<b>Client ID:</b>	<b>DP13-0.0-1.5</b>					
Arsenic	ND	13	6010B	2-14-12	2-14-12	
Barium	97	3.1	6010B	2-14-12	2-14-12	
Cadmium	ND	0.63	6010B	2-14-12	2-14-12	
Chromium	37	0.63	6010B	2-14-12	2-14-12	
Lead	ND	6.3	6010B	2-14-12	2-14-12	
Mercury	ND	0.31	7471A	2-14-12	2-14-12	
Selenium	ND	13	6010B	2-14-12	2-14-12	
Silver	ND	0.63	6010B	2-14-12	2-14-12	

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 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-03					
<b>Client ID:</b>	<b>DP14-0.0-2.0</b>					
Arsenic	ND	11	6010B	2-14-12	2-14-12	
Barium	43	2.8	6010B	2-14-12	2-14-12	
Cadmium	ND	0.56	6010B	2-14-12	2-14-12	
Chromium	31	0.56	6010B	2-14-12	2-14-12	
Lead	ND	5.6	6010B	2-14-12	2-14-12	
Mercury	ND	0.28	7471A	2-14-12	2-14-12	
Selenium	ND	11	6010B	2-14-12	2-14-12	
Silver	ND	0.56	6010B	2-14-12	2-14-12	

Lab ID:	02-085-04					
<b>Client ID:</b>	<b>DP15-0.0-2.0</b>					
Arsenic	ND	12	6010B	2-14-12	2-14-12	
Barium	110	3.0	6010B	2-14-12	2-14-12	
Cadmium	ND	0.59	6010B	2-14-12	2-14-12	
Chromium	25	0.59	6010B	2-14-12	2-14-12	
Lead	ND	5.9	6010B	2-14-12	2-14-12	
Mercury	ND	0.30	7471A	2-14-12	2-14-12	
Selenium	ND	12	6010B	2-14-12	2-14-12	
Silver	ND	0.59	6010B	2-14-12	2-14-12	

Date of Report: February 16, 2012  
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 Laboratory Reference: 1202-085  
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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-05					
<b>Client ID:</b>	<b>DP16-0.0-2.0</b>					
Arsenic	ND	12	6010B	2-14-12	2-14-12	
Barium	45	2.9	6010B	2-14-12	2-14-12	
Cadmium	ND	0.58	6010B	2-14-12	2-14-12	
Chromium	28	0.58	6010B	2-14-12	2-14-12	
Lead	ND	5.8	6010B	2-14-12	2-14-12	
Mercury	ND	0.29	7471A	2-14-12	2-14-12	
Selenium	ND	12	6010B	2-14-12	2-14-12	
Silver	ND	0.58	6010B	2-14-12	2-14-12	

Lab ID:	02-085-06					
<b>Client ID:</b>	<b>DP17-0.0-2.0</b>					
Arsenic	ND	12	6010B	2-14-12	2-14-12	
Barium	55	2.9	6010B	2-14-12	2-14-12	
Cadmium	ND	0.58	6010B	2-14-12	2-14-12	
Chromium	31	0.58	6010B	2-14-12	2-14-12	
Lead	ND	5.8	6010B	2-14-12	2-14-12	
Mercury	ND	0.29	7471A	2-14-12	2-14-12	
Selenium	ND	12	6010B	2-14-12	2-14-12	
Silver	ND	0.58	6010B	2-14-12	2-14-12	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-10					
<b>Client ID:</b>	<b>DP21-0.0-2.0</b>					
Arsenic	ND	12	6010B	2-14-12	2-14-12	
Barium	81	2.9	6010B	2-14-12	2-14-12	
Cadmium	0.83	0.58	6010B	2-14-12	2-14-12	
Chromium	45	0.58	6010B	2-14-12	2-14-12	
Lead	48	5.8	6010B	2-14-12	2-14-12	
Mercury	ND	0.29	7471A	2-14-12	2-14-12	
Selenium	ND	12	6010B	2-14-12	2-14-12	
Silver	ND	0.58	6010B	2-14-12	2-14-12	

Lab ID:	02-085-12					
<b>Client ID:</b>	<b>DP12-5.0-7.5</b>					
Arsenic	ND	12	6010B	2-14-12	2-14-12	
Barium	29	3.0	6010B	2-14-12	2-14-12	
Cadmium	ND	0.59	6010B	2-14-12	2-14-12	
Chromium	35	0.59	6010B	2-14-12	2-14-12	
Lead	ND	5.9	6010B	2-14-12	2-14-12	
Mercury	ND	0.30	7471A	2-14-12	2-14-12	
Selenium	ND	12	6010B	2-14-12	2-14-12	
Silver	ND	0.59	6010B	2-14-12	2-14-12	



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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-19					
<b>Client ID:</b>	<b>DP15-5.0-7.0</b>					
Arsenic	ND	13	6010B	2-14-12	2-14-12	
Barium	62	3.2	6010B	2-14-12	2-14-12	
Cadmium	ND	0.65	6010B	2-14-12	2-14-12	
Chromium	26	0.65	6010B	2-14-12	2-14-12	
Lead	ND	6.5	6010B	2-14-12	2-14-12	
Mercury	ND	0.32	7471A	2-14-12	2-14-12	
Selenium	ND	13	6010B	2-14-12	2-14-12	
Silver	ND	0.65	6010B	2-14-12	2-14-12	

Lab ID:	02-085-27					
<b>Client ID:</b>	<b>DP21-2.0-3.0</b>					
Arsenic	ND	13	6010B	2-14-12	2-14-12	
Barium	86	3.2	6010B	2-14-12	2-14-12	
Cadmium	ND	0.63	6010B	2-14-12	2-14-12	
Chromium	44	0.63	6010B	2-14-12	2-14-12	
Lead	ND	6.3	6010B	2-14-12	2-14-12	
Mercury	ND	0.32	7471A	2-14-12	2-14-12	
Selenium	ND	13	6010B	2-14-12	2-14-12	
Silver	ND	0.63	6010B	2-14-12	2-14-12	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-28					
Client ID:	HA1-0.0-2.0					
Arsenic	ND	12	6010B	2-14-12	2-14-12	
Barium	120	3.1	6010B	2-14-12	2-14-12	
Cadmium	ND	0.62	6010B	2-14-12	2-14-12	
Chromium	33	0.62	6010B	2-14-12	2-14-12	
Lead	ND	6.2	6010B	2-14-12	2-14-12	
Mercury	ND	0.31	7471A	2-14-12	2-14-12	
Selenium	ND	12	6010B	2-14-12	2-14-12	
Silver	ND	0.62	6010B	2-14-12	2-14-12	

Lab ID:	02-085-29					
Client ID:	HA2-0.0-2.0					
Arsenic	ND	13	6010B	2-14-12	2-14-12	
Barium	93	3.4	6010B	2-14-12	2-14-12	
Cadmium	ND	0.67	6010B	2-14-12	2-14-12	
Chromium	37	0.67	6010B	2-14-12	2-14-12	
Lead	9.6	6.7	6010B	2-14-12	2-14-12	
Mercury	ND	0.34	7471A	2-14-12	2-14-12	
Selenium	ND	13	6010B	2-14-12	2-14-12	
Silver	ND	0.67	6010B	2-14-12	2-14-12	

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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-30					
Client ID:	HA3-0.0-2.0					
Arsenic	ND	12	6010B	2-14-12	2-14-12	
Barium	91	3.0	6010B	2-14-12	2-14-12	
Cadmium	2.7	0.60	6010B	2-14-12	2-14-12	
Chromium	58	0.60	6010B	2-14-12	2-14-12	
Lead	150	6.0	6010B	2-14-12	2-14-12	
Mercury	ND	0.30	7471A	2-14-12	2-14-12	
Selenium	ND	12	6010B	2-14-12	2-14-12	
Silver	ND	0.60	6010B	2-14-12	2-14-12	

Date of Report: February 16, 2012  
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**SOLUBLE HEXAVALENT CHROMIUM  
 WATER EXTRACTION  
 EPA 7196A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-085-31					
Client ID:	HA4-0.0-2.0					
Hexavalent Chromium	ND	1.3	7196A mod	2-15-12	2-15-12	

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**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0210S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	114	50-150				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	02-085-01					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil	<b>114</b>	<b>90.4</b>		23	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			102 104	50-150		
Laboratory ID:	02-085-05					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			105 98	50-150		

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**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0213S1					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	93	68-124				
Laboratory ID:	MB0213S2					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	93	68-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	02-085-01							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	30	
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				90	87	68-124		
Laboratory ID:	02-085-02							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	30	
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				101	99	68-124		

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0213S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chloromethane	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Vinyl Chloride	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromomethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chloroethane	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Trichlorofluoromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Acetone	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Iodomethane	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Carbon Disulfide	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Methylene Chloride	ND	0.0050	EPA 8260	2-13-12	2-13-12	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1-Dichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Vinyl Acetate	ND	0.0050	EPA 8260	2-13-12	2-13-12	
2,2-Dichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Butanone	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Bromochloromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chloroform	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Carbon Tetrachloride	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1-Dichloropropene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Benzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Trichloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Dibromomethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromodichloromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	2-13-12	2-13-12	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Toluene	ND	0.0050	EPA 8260	2-13-12	2-13-12	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	2-13-12	2-13-12	

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**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0213S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Tetrachloroethene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,3-Dichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Hexanone	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Dibromochloromethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dibromoethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Chlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Ethylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
m,p-Xylene	ND	0.0020	EPA 8260	2-13-12	2-13-12	
o-Xylene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Styrene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromoform	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Isopropylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Bromobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	2-13-12	2-13-12	
n-Propylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
2-Chlorotoluene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
4-Chlorotoluene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
tert-Butylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
sec-Butylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
p-Isopropyltoluene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
n-Butylbenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	2-13-12	2-13-12	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
Hexachlorobutadiene	ND	0.0050	EPA 8260	2-13-12	2-13-12	
Naphthalene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>114</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>110</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>93</i>	<i>55-121</i>				



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**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0214S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Chloromethane	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Vinyl Chloride	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Bromomethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Chloroethane	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Trichlorofluoromethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,1-Dichloroethene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Acetone	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Iodomethane	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Carbon Disulfide	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Methylene Chloride	ND	0.0050	EPA 8260	2-14-12	2-14-12	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,1-Dichloroethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Vinyl Acetate	ND	0.0050	EPA 8260	2-14-12	2-14-12	
2,2-Dichloropropane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
2-Butanone	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Bromochloromethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Chloroform	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Carbon Tetrachloride	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,1-Dichloropropene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Benzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2-Dichloroethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Trichloroethene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2-Dichloropropane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Dibromomethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Bromodichloromethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	2-14-12	2-14-12	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Toluene	ND	0.0050	EPA 8260	2-14-12	2-14-12	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	2-14-12	2-14-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0214S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Tetrachloroethene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,3-Dichloropropane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
2-Hexanone	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Dibromochloromethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2-Dibromoethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Chlorobenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Ethylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
m,p-Xylene	ND	0.0020	EPA 8260	2-14-12	2-14-12	
o-Xylene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Styrene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Bromoform	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Isopropylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Bromobenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	2-14-12	2-14-12	
n-Propylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
2-Chlorotoluene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
4-Chlorotoluene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
tert-Butylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
sec-Butylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
p-Isopropyltoluene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
n-Butylbenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	2-14-12	2-14-12	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
Hexachlorobutadiene	ND	0.0050	EPA 8260	2-14-12	2-14-12	
Naphthalene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>112</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>109</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>91</i>	<i>55-121</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
**SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
	SB	SBD	SB	SBD	SB	SBD				
Laboratory ID:	SB0213S1									
1,1-Dichloroethene	<b>0.0603</b>	<b>0.0593</b>	0.0500	0.0500	121	119	70-130	2	19	
Benzene	<b>0.0542</b>	<b>0.0554</b>	0.0500	0.0500	108	111	70-125	2	15	
Trichloroethene	<b>0.0494</b>	<b>0.0482</b>	0.0500	0.0500	99	96	70-122	2	14	
Toluene	<b>0.0512</b>	<b>0.0502</b>	0.0500	0.0500	102	100	73-120	2	16	
Chlorobenzene	<b>0.0500</b>	<b>0.0489</b>	0.0500	0.0500	100	98	74-109	2	12	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					100	101	63-127			
<i>Toluene-d8</i>					96	96	65-129			
<i>4-Bromofluorobenzene</i>					82	81	55-121			
Laboratory ID:	SB0214S1									
1,1-Dichloroethene	<b>0.0571</b>	<b>0.0582</b>	0.0500	0.0500	114	116	70-130	2	19	
Benzene	<b>0.0537</b>	<b>0.0534</b>	0.0500	0.0500	107	107	70-125	1	15	
Trichloroethene	<b>0.0498</b>	<b>0.0499</b>	0.0500	0.0500	100	100	70-122	0	14	
Toluene	<b>0.0503</b>	<b>0.0508</b>	0.0500	0.0500	101	102	73-120	1	16	
Chlorobenzene	<b>0.0497</b>	<b>0.0491</b>	0.0500	0.0500	99	98	74-109	1	12	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					100	103	63-127			
<i>Toluene-d8</i>					97	100	65-129			
<i>4-Bromofluorobenzene</i>					80	85	55-121			

Date of Report: February 16, 2012  
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 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0215S1					
n-Nitrosodimethylamine	ND	0.033	EPA 8270	2-15-12	2-15-12	
Pyridine	ND	0.33	EPA 8270	2-15-12	2-15-12	
Phenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
Aniline	ND	0.033	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270	2-15-12	2-15-12	
2-Chlorophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,3-Dichlorobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,4-Dichlorobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
Benzyl alcohol	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,2-Dichlorobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270	2-15-12	2-15-12	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270	2-15-12	2-15-12	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270	2-15-12	2-15-12	
Hexachloroethane	ND	0.033	EPA 8270	2-15-12	2-15-12	
Nitrobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
Isophorone	ND	0.033	EPA 8270	2-15-12	2-15-12	
2-Nitrophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,4-Dimethylphenol	ND	0.33	EPA 8270	2-15-12	2-15-12	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,4-Dichlorophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
Naphthalene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
4-Chloroaniline	ND	0.033	EPA 8270	2-15-12	2-15-12	
Hexachlorobutadiene	ND	0.033	EPA 8270	2-15-12	2-15-12	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,3-Dichloroaniline	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
2-Chloronaphthalene	ND	0.033	EPA 8270	2-15-12	2-15-12	
2-Nitroaniline	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,4-Dinitrobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
Dimethylphthalate	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,3-Dinitrobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,6-Dinitrotoluene	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,2-Dinitrobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
3-Nitroaniline	ND	0.033	EPA 8270	2-15-12	2-15-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
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 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0215S1					
2,4-Dinitrophenol	ND	0.17	EPA 8270	2-15-12	2-15-12	
Acenaphthene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
4-Nitrophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,4-Dinitrotoluene	ND	0.033	EPA 8270	2-15-12	2-15-12	
Dibenzofuran	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270	2-15-12	2-15-12	
Diethylphthalate	ND	0.17	EPA 8270	2-15-12	2-15-12	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270	2-15-12	2-15-12	
4-Nitroaniline	ND	0.033	EPA 8270	2-15-12	2-15-12	
Fluorene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270	2-15-12	2-15-12	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270	2-15-12	2-15-12	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270	2-15-12	2-15-12	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270	2-15-12	2-15-12	
Hexachlorobenzene	ND	0.033	EPA 8270	2-15-12	2-15-12	
Pentachlorophenol	ND	0.17	EPA 8270	2-15-12	2-15-12	
Phenanthrene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Anthracene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Carbazole	ND	0.033	EPA 8270	2-15-12	2-15-12	
Di-n-butylphthalate	ND	0.33	EPA 8270	2-15-12	2-15-12	
Fluoranthene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Benzidine	ND	0.33	EPA 8270	2-15-12	2-15-12	
Pyrene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Butylbenzylphthalate	ND	0.33	EPA 8270	2-15-12	2-15-12	
bis-2-Ethylhexyladipate	ND	0.17	EPA 8270	2-15-12	2-15-12	
3,3'-Dichlorobenzidine	ND	0.33	EPA 8270	2-15-12	2-15-12	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Chrysene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
bis(2-Ethylhexyl)phthalate	ND	0.17	EPA 8270	2-15-12	2-15-12	
Di-n-octylphthalate	ND	0.033	EPA 8270	2-15-12	2-15-12	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	2-15-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>58</i>	<i>30 - 97</i>				
<i>Phenol-d6</i>	<i>66</i>	<i>40 - 104</i>				
<i>Nitrobenzene-d5</i>	<i>61</i>	<i>35 - 102</i>				
<i>2-Fluorobiphenyl</i>	<i>70</i>	<i>44 - 97</i>				
<i>2,4,6-Tribromophenol</i>	<i>88</i>	<i>41 - 110</i>				
<i>Terphenyl-d14</i>	<i>82</i>	<i>53 - 107</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					Recovery	Limits	Limit			
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0215S1									
	SB	SBD	SB	SBD	SB	SBD				
Phenol	<b>0.985</b>	<b>0.923</b>	1.33	1.33	74	69	31 - 111	6	34	
2-Chlorophenol	<b>1.08</b>	<b>1.05</b>	1.33	1.33	81	79	29 - 112	3	37	
1,4-Dichlorobenzene	<b>0.534</b>	<b>0.529</b>	0.667	0.667	80	79	24 - 100	1	37	
n-Nitroso-di-n-propylamine	<b>0.507</b>	<b>0.501</b>	0.667	0.667	76	75	35 - 104	1	32	
1,2,4-Trichlorobenzene	<b>0.553</b>	<b>0.544</b>	0.667	0.667	83	82	29 - 110	2	35	
4-Chloro-3-methylphenol	<b>1.18</b>	<b>1.21</b>	1.33	1.33	89	91	53 - 104	3	25	
Acenaphthene	<b>0.595</b>	<b>0.580</b>	0.667	0.667	89	87	50 - 95	3	23	
4-Nitrophenol	<b>1.44</b>	<b>1.52</b>	1.33	1.33	108	114	42 - 126	5	30	
2,4-Dinitrotoluene	<b>0.692</b>	<b>0.707</b>	0.667	0.667	104	106	53 - 115	2	31	
Pentachlorophenol	<b>1.37</b>	<b>1.41</b>	1.33	1.33	103	106	50 - 116	3	30	
Pyrene	<b>0.620</b>	<b>0.649</b>	0.667	0.667	93	97	57 - 120	5	27	
<i>Surrogate:</i>										
2-Fluorophenol					65	62	30 - 97			
Phenol-d6					70	68	40 - 104			
Nitrobenzene-d5					65	65	35 - 102			
2-Fluorobiphenyl					73	72	44 - 97			
2,4,6-Tribromophenol					89	90	41 - 110			
Terphenyl-d14					79	83	53 - 107			

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 (with silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0214S2					
Naphthalene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Acenaphthene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Fluorene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Phenanthrene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Anthracene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Fluoranthene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Pyrene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Chrysene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	2-14-12	2-14-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>85</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>93</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>89</i>	<i>33 - 119</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0214S2									
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	<b>0.0696</b>	<b>0.0706</b>	0.0833	0.0833	84	85	43 - 108	1	27	
Acenaphthylene	<b>0.0733</b>	<b>0.0754</b>	0.0833	0.0833	88	91	52 - 120	3	21	
Acenaphthene	<b>0.0690</b>	<b>0.0704</b>	0.0833	0.0833	83	85	59 - 113	2	17	
Fluorene	<b>0.0730</b>	<b>0.0763</b>	0.0833	0.0833	88	92	64 - 117	4	14	
Phenanthrene	<b>0.0709</b>	<b>0.0717</b>	0.0833	0.0833	85	86	67 - 112	1	12	
Anthracene	<b>0.0780</b>	<b>0.0809</b>	0.0833	0.0833	94	97	59 - 110	4	16	
Fluoranthene	<b>0.0776</b>	<b>0.0795</b>	0.0833	0.0833	93	95	68 - 120	2	15	
Pyrene	<b>0.0794</b>	<b>0.0807</b>	0.0833	0.0833	95	97	66 - 121	2	17	
Benzo[a]anthracene	<b>0.0689</b>	<b>0.0716</b>	0.0833	0.0833	83	86	63 - 114	4	12	
Chrysene	<b>0.0768</b>	<b>0.0788</b>	0.0833	0.0833	92	95	67 - 118	3	12	
Benzo[b]fluoranthene	<b>0.0778</b>	<b>0.0793</b>	0.0833	0.0833	93	95	58 - 127	2	20	
Benzo(j,k)fluoranthene	<b>0.0809</b>	<b>0.0845</b>	0.0833	0.0833	97	101	42 - 134	4	26	
Benzo[a]pyrene	<b>0.0609</b>	<b>0.0606</b>	0.0833	0.0833	73	73	55 - 111	0	19	
Indeno(1,2,3-c,d)pyrene	<b>0.0617</b>	<b>0.0616</b>	0.0833	0.0833	74	74	60 - 125	0	20	
Dibenz[a,h]anthracene	<b>0.0721</b>	<b>0.0744</b>	0.0833	0.0833	87	89	62 - 125	3	19	
Benzo[g,h,i]perylene	<b>0.0711</b>	<b>0.0732</b>	0.0833	0.0833	85	88	61 - 124	3	19	
<i>Surrogate:</i>										
2-Fluorobiphenyl					74	82	43 - 109			
Pyrene-d10					85	85	38 - 128			
Terphenyl-d14					79	79	33 - 119			



Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0214S1					
Aroclor 1016	<b>ND</b>	0.050	EPA 8082	2-14-12	2-15-12	
Aroclor 1221	<b>ND</b>	0.050	EPA 8082	2-14-12	2-15-12	
Aroclor 1232	<b>ND</b>	0.050	EPA 8082	2-14-12	2-15-12	
Aroclor 1242	<b>ND</b>	0.050	EPA 8082	2-14-12	2-15-12	
Aroclor 1248	<b>ND</b>	0.050	EPA 8082	2-14-12	2-15-12	
Aroclor 1254	<b>ND</b>	0.050	EPA 8082	2-14-12	2-15-12	
Aroclor 1260	<b>ND</b>	0.050	EPA 8082	2-14-12	2-15-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	98		42-123			

Analyte	Result		Spike Level		Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>										
Laboratory ID:	02-093-01									
	MS	MSD	MS	MSD		MS	MSD			
Aroclor 1260	<b>0.561</b>	<b>0.503</b>	0.500	0.500	ND	<b>112</b>	<b>101</b>	44-125	11	15
<i>Surrogate:</i>										
DCB						102	91	42-123		

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**TOTAL METALS  
EPA 6010B  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-14-12  
Date Analyzed: 2-14-12  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB0214SM1

Analyte	Method	Result	PQL
Arsenic	6010B	<b>ND</b>	10
Barium	6010B	<b>ND</b>	2.5
Cadmium	6010B	<b>ND</b>	0.50
Chromium	6010B	<b>ND</b>	0.50
Lead	6010B	<b>ND</b>	5.0
Selenium	6010B	<b>ND</b>	10
Silver	6010B	<b>ND</b>	0.50

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**TOTAL MERCURY  
EPA 7471A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-10-12  
Date Analyzed: 2-10-12  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB0210S2

Analyte	Method	Result	PQL
Mercury	7471A	<b>ND</b>	0.25

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 2-14-12

Date Analyzed: 2-14-12

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 02-110-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	37.5	37.0	1	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	13.8	16.1	16	0.50	
Lead	ND	ND	NA	5.0	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**TOTAL MERCURY  
EPA 7471A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 2-10-12

Date Analyzed: 2-10-12

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 02-085-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Mercury	<b>ND</b>	<b>ND</b>	NA	0.25	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B  
 MS/MSD QUALITY CONTROL**

Date Extracted: 2-14-12

Date Analyzed: 2-14-12

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 02-110-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>100</b>	100	<b>101</b>	101	1	
Barium	100	<b>146</b>	109	<b>141</b>	104	4	
Cadmium	50.0	<b>49.4</b>	99	<b>50.2</b>	100	2	
Chromium	100	<b>114</b>	100	<b>114</b>	100	0	
Lead	250	<b>245</b>	98	<b>248</b>	99	1	
Selenium	100	<b>99.1</b>	99	<b>106</b>	106	7	
Silver	25.0	<b>23.3</b>	93	<b>24.0</b>	96	3	

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**TOTAL MERCURY  
EPA 7471A  
MS/MSD QUALITY CONTROL**

Date Extracted: 2-10-12

Date Analyzed: 2-10-12

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 02-085-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Mercury	0.500	<b>0.551</b>	110	<b>0.542</b>	108	2	

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**SOLUBLE HEXAVALENT CHROMIUM  
WATER EXTRACTION  
EPA 7196A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-15-12  
Date Analyzed: 2-15-12  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB0215S1

Analyte	Method	Result	PQL
Hexavalent Chromium	7196A mod	<b>ND</b>	1.0



Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**SOLUBLE HEXAVALENT CHROMIUM  
WATER EXTRACTION  
EPA 7196A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 2-15-12  
Date Analyzed: 2-15-12  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: 02-093-01

Analyte	Sample Result	Duplicate Result	RPD	Flags	PQL
Hexavalent Chromium	<b>ND</b>	<b>ND</b>	NA		1.0

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**SOLUBLE HEXAVALENT CHROMIUM  
WATER EXTRACTION  
EPA 7196A  
MS/MSD QUALITY CONTROL**

Date Extracted: 2-15-12  
Date Analyzed: 2-15-12  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: 02-093-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Hexavalent Chromium	5.00	<b>5.11</b>	102	<b>4.93</b>	99	4	

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085  
Project: 0180-292-00

**% MOISTURE**

Date Analyzed: 2-10,13&14-12

Client ID	Lab ID	% Moisture
DP12-0.0-2.0	02-085-01	10
DP13-0.0-1.5	02-085-02	20
DP14-0.0-2.0	02-085-03	11
DP15-0.0-2.0	02-085-04	16
DP16-0.0-2.0	02-085-05	13
DP17-0.0-2.0	02-085-06	14
DP18-0.0-2.0	02-085-07	13
DP19-0.0-2.0	02-085-08	11
DP20-0.0-2.0	02-085-09	15
DP21-0.0-2.0	02-085-10	14
DP12-5.0-7.5	02-085-12	15
DP15-5.0-7.0	02-085-19	23
DP21-2.0-3.0	02-085-27	21
HA1-0.0-2.0	02-085-28	20
HA2-0.0-2.0	02-085-29	26
HA3-0.0-2.0	02-085-30	17
HA4-0.0-2.0	02-085-31	21
SED9	02-085-32	60



#### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference





# TestAmerica

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Rush 4 days  
 Short Hold

Chain of  
 Custody Record  
 02-085

Client: Geotriggers Client Contact: Aaron Johnson Date: 2/9/12  
 Address: 1101 S. Fawcett Ave Ste 200 Telephone Number (Area Code)/Fax Number: (253) 383-4940 Lab Number: 12705

City: Tacoma State: WA Zip Code: 98402 Sampler: Aaron Johnson Lab Contact: David Bunker Page: 2 of 3  
 Project Name and Location (State): WSOOT Midway Metals Seguin WA Billing Contact: David Bunker

Contract/Purchase Order/Quote No.: 0180-292-00 Matrix: Unpres. H2SO4 HN03 HCl NaOH ZnAc/NaOH  
 Sample I.D. and Location/Description (Containers for each sample may be combined on one line)

Sample I.D. and Location/Description	Date	Time	Air	Aqueous	Sed.	Soil	Unpres.	H2SO4	HN03	HCl	NaOH	ZnAc/NaOH	Containers & Preservatives	Analysis (Attach list if more space is needed)	Special Instructions/Conditions of Receipt
13 DP12-7.5-8.5	2/8/12	1115				X	5							NWTPH-D <sub>1</sub> w/Sp/2a del cleanup	02-085
14 DP13-5.0-7.0		1135				X	5							NWTPH-6x	
15 DP13-8.0-1.0		1140				X	5							VOCs lowlevel	
16 DP14-2.0-3.0		1215				X	5							SVOCs lowlevel	
17 DP14-5.0-7.0		1230				X	5							PAHs lowlevel	
18 DP15-2.0-3.0		1320				X	5							RCRA 8metals	
19 DP15-5.0-7.0		1330				X	5							PCBs	
20 DP16-2.0-3.0		1400				X	5							Hold	
21 DP16-5.0-7.0		1410				X	5								
22 DP17-2.0-3.0		1435				X	5								
23 DP17-5.0-7.0		1440				X	5								
24 DP18-2.0-3.0		1525				X	5								

Cooler:  Yes  No Cooler Temp: \_\_\_\_\_ Possible Hazard Identification:  Non-Hazard  Flammable  Skin Irritant  Poison B  Unknown  Sample Disposal:  Return To Client  Disposal By Lab

Turn Around Time Required (business days):  24 Hours  48 Hours  5 Days  10 Days  15 Days  Other \_\_\_\_\_ QC Requirements (Specify): \_\_\_\_\_

(A fee may be assessed if samples are retained longer than 1 month)

1. Relinquished By Sign/Print: [Signature] Date: 2/9/12 Time: 1630 Received By Sign/Print: [Signature] Date: 2/9/12 Time: 1630

2. Relinquished By Sign/Print: [Signature] Date: \_\_\_\_\_ Time: \_\_\_\_\_ Received By Sign/Print: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

3. Relinquished By Sign/Print: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ Received By Sign/Print: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

Comments: \_\_\_\_\_







14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

December 29, 2011

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1112-098B

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on December 14, 2011.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures



Date of Report: December 29, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098B  
Project: 0180-292-00

### **Case Narrative**

Samples were collected on December 13 and 14, 2011 and received by the laboratory on December 14, 2011. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### NWTPH Gx Analysis

Sample analysis holding time for DP-8-2-3 was exceeded by approximately one day.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

#### NWTPH Dx Analysis

Samples DP-3-3-4, DP-4-5-6, DP-5-6-7, and DP-6-3.5-4.5 were extracted and analyzed out of holding time.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

#### Volatiles EPA 8260B Analysis

The holding times for samples DP-4-5-6, DP-5-6-7, and DP-6-3.5-4.5 were exceeded by one day.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

#### PAHs EPA 8270D/SIM Analysis

The sample DP-9-6-7 was extracted and analyzed one day out of holding time.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Date of Report: December 29, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098B  
Project: 0180-292-00

### ANALYTICAL REPORT FOR SAMPLES

Client ID	Laboratory ID	Matrix	Date Sampled	Date Received	Notes
DP-3-3-4	12-098-05	Soil	12-13-11	12-14-11	
DP-4-5-6	12-098-08	Soil	12-13-11	12-14-11	
DP-5-6-7	12-098-10	Soil	12-13-11	12-14-11	
DP-6-3.5-4.5	12-098-13	Soil	12-13-11	12-14-11	
DP-7-5-6	12-098-16	Soil	12-14-11	12-14-11	
DP-8-2-3	12-098-18	Soil	12-14-11	12-14-11	
DP-9-6-7	12-098-21	Soil	12-14-11	12-14-11	

Date of Report: December 29, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098B  
Project: 0180-292-00

**NWTPH-Gx**

Matrix: Soil  
Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-8-2-3</b>					
Laboratory ID:	12-098-18					
Gasoline	<b>ND</b>	5.2	NWTPH-Gx	12-29-11	12-29-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	68-124				

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-3-3-4</b>					
Laboratory ID:	12-098-05					
Diesel Range Organics	<b>ND</b>	29	NWTPH-Dx	12-28-11	12-28-11	
Lube Oil Range Organics	<b>ND</b>	58	NWTPH-Dx	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	110	50-150				
<b>Client ID:</b>	<b>DP-4-5-6</b>					
Laboratory ID:	12-098-08					
Diesel Range Organics	<b>ND</b>	32	NWTPH-Dx	12-28-11	12-28-11	
Lube Oil Range Organics	<b>ND</b>	64	NWTPH-Dx	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	101	50-150				
<b>Client ID:</b>	<b>DP-5-6-7</b>					
Laboratory ID:	12-098-10					
Diesel Range Organics	<b>ND</b>	30	NWTPH-Dx	12-28-11	12-28-11	
Lube Oil Range Organics	<b>ND</b>	60	NWTPH-Dx	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	102	50-150				
<b>Client ID:</b>	<b>DP-6-3.5-4.5</b>					
Laboratory ID:	12-098-13					
Diesel Range Organics	<b>ND</b>	29	NWTPH-Dx	12-28-11	12-28-11	
Lube Oil Range Organics	<b>ND</b>	59	NWTPH-Dx	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	103	50-150				

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
 page 1 of 2

Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-4-5-6</b>					
Laboratory ID:	12-098-08					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chloromethane	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Vinyl Chloride	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromomethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chloroethane	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Trichlorofluoromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Acetone	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Iodomethane	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Carbon Disulfide	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Methylene Chloride	ND	0.0056	EPA 8260	12-28-11	12-28-11	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Vinyl Acetate	ND	0.0056	EPA 8260	12-28-11	12-28-11	
2,2-Dichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Butanone	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Bromochloromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chloroform	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Carbon Tetrachloride	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1-Dichloropropene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Benzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Trichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Dibromomethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromodichloromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Chloroethyl Vinyl Ether	ND	0.0056	EPA 8260	12-28-11	12-28-11	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Methyl Isobutyl Ketone	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Toluene	ND	0.0056	EPA 8260	12-28-11	12-28-11	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-28-11	12-28-11	

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 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-4-5-6</b>					
Laboratory ID:	12-098-08					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Tetrachloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,3-Dichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Hexanone	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Dibromochloromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dibromoethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Ethylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
m,p-Xylene	ND	0.0023	EPA 8260	12-28-11	12-28-11	
o-Xylene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Styrene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromoform	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Isopropylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
n-Propylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Chlorotoluene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
4-Chlorotoluene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
tert-Butylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
sec-Butylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
p-Isopropyltoluene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
n-Butylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dibromo-3-chloropropane	ND	0.0056	EPA 8260	12-28-11	12-28-11	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Hexachlorobutadiene	ND	0.0056	EPA 8260	12-28-11	12-28-11	
Naphthalene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>107</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>111</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>86</i>	<i>55-121</i>				

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 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-5-6-7</b>					
Laboratory ID:	12-098-10					
Dichlorodifluoromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chloromethane	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Vinyl Chloride	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromomethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chloroethane	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Trichlorofluoromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Acetone	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Iodomethane	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Carbon Disulfide	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Methylene Chloride	ND	0.0054	EPA 8260	12-28-11	12-28-11	
(trans) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Methyl t-Butyl Ether	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Vinyl Acetate	ND	0.0054	EPA 8260	12-28-11	12-28-11	
2,2-Dichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
(cis) 1,2-Dichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Butanone	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Bromochloromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chloroform	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1,1-Trichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Carbon Tetrachloride	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1-Dichloropropene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Benzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Trichloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Dibromomethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromodichloromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Chloroethyl Vinyl Ether	ND	0.0054	EPA 8260	12-28-11	12-28-11	
(cis) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Methyl Isobutyl Ketone	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Toluene	ND	0.0054	EPA 8260	12-28-11	12-28-11	
(trans) 1,3-Dichloropropene	ND	0.0011	EPA 8260	12-28-11	12-28-11	

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 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-5-6-7</b>					
Laboratory ID:	12-098-10					
1,1,2-Trichloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Tetrachloroethene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,3-Dichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Hexanone	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Dibromochloromethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dibromoethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Chlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1,1,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Ethylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
m,p-Xylene	ND	0.0022	EPA 8260	12-28-11	12-28-11	
o-Xylene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Styrene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromoform	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Isopropylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Bromobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,1,2,2-Tetrachloroethane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichloropropane	ND	0.0011	EPA 8260	12-28-11	12-28-11	
n-Propylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
2-Chlorotoluene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
4-Chlorotoluene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,3,5-Trimethylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
tert-Butylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2,4-Trimethylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
sec-Butylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,3-Dichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
p-Isopropyltoluene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,4-Dichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
n-Butylbenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2-Dibromo-3-chloropropane	ND	0.0054	EPA 8260	12-28-11	12-28-11	
1,2,4-Trichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
Hexachlorobutadiene	ND	0.0054	EPA 8260	12-28-11	12-28-11	
Naphthalene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichlorobenzene	ND	0.0011	EPA 8260	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>109</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>110</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>86</i>	<i>55-121</i>				



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 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-6-3.5-4.5</b>					
Laboratory ID:	12-098-13					
Dichlorodifluoromethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Chloromethane	ND	0.0048	EPA 8260	12-28-11	12-28-11	
Vinyl Chloride	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Bromomethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Chloroethane	ND	0.0048	EPA 8260	12-28-11	12-28-11	
Trichlorofluoromethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Acetone	0.016	0.0048	EPA 8260	12-28-11	12-28-11	
Iodomethane	ND	0.0048	EPA 8260	12-28-11	12-28-11	
Carbon Disulfide	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Methylene Chloride	ND	0.0048	EPA 8260	12-28-11	12-28-11	
(trans) 1,2-Dichloroethene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Methyl t-Butyl Ether	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Vinyl Acetate	ND	0.0048	EPA 8260	12-28-11	12-28-11	
2,2-Dichloropropane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
(cis) 1,2-Dichloroethene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
2-Butanone	ND	0.0048	EPA 8260	12-28-11	12-28-11	
Bromochloromethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Chloroform	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,1,1-Trichloroethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Carbon Tetrachloride	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,1-Dichloropropene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Benzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2-Dichloroethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Trichloroethene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2-Dichloropropane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Dibromomethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Bromodichloromethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
2-Chloroethyl Vinyl Ether	ND	0.0048	EPA 8260	12-28-11	12-28-11	
(cis) 1,3-Dichloropropene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Methyl Isobutyl Ketone	ND	0.0048	EPA 8260	12-28-11	12-28-11	
Toluene	ND	0.0048	EPA 8260	12-28-11	12-28-11	
(trans) 1,3-Dichloropropene	ND	0.00097	EPA 8260	12-28-11	12-28-11	

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 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-6-3.5-4.5</b>					
Laboratory ID:	12-098-13					
1,1,2-Trichloroethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Tetrachloroethene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,3-Dichloropropane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
2-Hexanone	ND	0.0048	EPA 8260	12-28-11	12-28-11	
Dibromochloromethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2-Dibromoethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Chlorobenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,1,1,2-Tetrachloroethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Ethylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
m,p-Xylene	ND	0.0019	EPA 8260	12-28-11	12-28-11	
o-Xylene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Styrene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Bromoform	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Isopropylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Bromobenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,1,2,2-Tetrachloroethane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichloropropane	ND	0.00097	EPA 8260	12-28-11	12-28-11	
n-Propylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
2-Chlorotoluene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
4-Chlorotoluene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,3,5-Trimethylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
tert-Butylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2,4-Trimethylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
sec-Butylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,3-Dichlorobenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
p-Isopropyltoluene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,4-Dichlorobenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2-Dichlorobenzene	0.0019	0.00097	EPA 8260	12-28-11	12-28-11	
n-Butylbenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2-Dibromo-3-chloropropane	ND	0.0048	EPA 8260	12-28-11	12-28-11	
1,2,4-Trichlorobenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
Hexachlorobutadiene	ND	0.0048	EPA 8260	12-28-11	12-28-11	
Naphthalene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichlorobenzene	ND	0.00097	EPA 8260	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>110</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>111</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>87</i>	<i>55-121</i>				

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP-9-6-7</b>					
Laboratory ID:	12-098-21					
Naphthalene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
2-Methylnaphthalene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
1-Methylnaphthalene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Acenaphthylene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Acenaphthene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Fluorene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Phenanthrene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Anthracene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Fluoranthene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Pyrene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[a]anthracene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Chrysene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[b]fluoranthene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[a]pyrene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0083	EPA 8270/SIM	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>54</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>54</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>52</i>	<i>33 - 119</i>				

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-4-5-6</b>					
Laboratory ID:	12-098-08					
Aroclor 1016	ND	0.064	EPA 8082	12-28-11	12-28-11	
Aroclor 1221	ND	0.064	EPA 8082	12-28-11	12-28-11	
Aroclor 1232	ND	0.064	EPA 8082	12-28-11	12-28-11	
Aroclor 1242	ND	0.064	EPA 8082	12-28-11	12-28-11	
Aroclor 1248	ND	0.064	EPA 8082	12-28-11	12-28-11	
Aroclor 1254	ND	0.064	EPA 8082	12-28-11	12-28-11	
Aroclor 1260	ND	0.064	EPA 8082	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	73	42-123				
<b>Client ID:</b>	<b>DP-5-6-7</b>					
Laboratory ID:	12-098-10					
Aroclor 1016	ND	0.060	EPA 8082	12-28-11	12-28-11	
Aroclor 1221	ND	0.060	EPA 8082	12-28-11	12-28-11	
Aroclor 1232	ND	0.060	EPA 8082	12-28-11	12-28-11	
Aroclor 1242	ND	0.060	EPA 8082	12-28-11	12-28-11	
Aroclor 1248	ND	0.060	EPA 8082	12-28-11	12-28-11	
Aroclor 1254	ND	0.060	EPA 8082	12-28-11	12-28-11	
Aroclor 1260	ND	0.060	EPA 8082	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	81	42-123				
<b>Client ID:</b>	<b>DP-6-3.5-4.5</b>					
Laboratory ID:	12-098-13					
Aroclor 1016	ND	0.059	EPA 8082	12-28-11	12-28-11	
Aroclor 1221	ND	0.059	EPA 8082	12-28-11	12-28-11	
Aroclor 1232	ND	0.059	EPA 8082	12-28-11	12-28-11	
Aroclor 1242	ND	0.059	EPA 8082	12-28-11	12-28-11	
Aroclor 1248	ND	0.059	EPA 8082	12-28-11	12-28-11	
Aroclor 1254	ND	0.059	EPA 8082	12-28-11	12-28-11	
Aroclor 1260	ND	0.059	EPA 8082	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	79	42-123				

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP-7-5-6</b>					
Laboratory ID:	12-098-16					
Aroclor 1016	<b>ND</b>	0.061	EPA 8082	12-28-11	12-28-11	
Aroclor 1221	<b>ND</b>	0.061	EPA 8082	12-28-11	12-28-11	
Aroclor 1232	<b>ND</b>	0.061	EPA 8082	12-28-11	12-28-11	
Aroclor 1242	<b>ND</b>	0.061	EPA 8082	12-28-11	12-28-11	
Aroclor 1248	<b>ND</b>	0.061	EPA 8082	12-28-11	12-28-11	
Aroclor 1254	<b>ND</b>	0.061	EPA 8082	12-28-11	12-28-11	
Aroclor 1260	<b>ND</b>	0.061	EPA 8082	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>81</i>	<i>42-123</i>				
<b>Client ID:</b>	<b>DP-8-2-3</b>					
Laboratory ID:	12-098-18					
Aroclor 1016	<b>ND</b>	0.057	EPA 8082	12-28-11	12-28-11	
Aroclor 1221	<b>ND</b>	0.057	EPA 8082	12-28-11	12-28-11	
Aroclor 1232	<b>ND</b>	0.057	EPA 8082	12-28-11	12-28-11	
Aroclor 1242	<b>ND</b>	0.057	EPA 8082	12-28-11	12-28-11	
Aroclor 1248	<b>ND</b>	0.057	EPA 8082	12-28-11	12-28-11	
Aroclor 1254	<b>ND</b>	0.057	EPA 8082	12-28-11	12-28-11	
Aroclor 1260	<b>ND</b>	0.057	EPA 8082	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>80</i>	<i>42-123</i>				

Date of Report: December 29, 2011  
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 Project: 0180-292-00

**TOTAL METALS**  
**EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-08					
<b>Client ID:</b>	<b>DP-4-5-6</b>					
Arsenic	<b>ND</b>	13	6010B	12-28-11	12-28-11	
Barium	<b>54</b>	3.2	6010B	12-28-11	12-28-11	
Cadmium	<b>ND</b>	0.64	6010B	12-28-11	12-28-11	
Chromium	<b>46</b>	0.64	6010B	12-28-11	12-28-11	
Lead	<b>ND</b>	6.4	6010B	12-28-11	12-28-11	
Mercury	<b>ND</b>	0.32	7471A	12-28-11	12-28-11	
Selenium	<b>ND</b>	13	6010B	12-28-11	12-28-11	
Silver	<b>ND</b>	0.64	6010B	12-28-11	12-28-11	

Lab ID:	12-098-10					
<b>Client ID:</b>	<b>DP-5-6-7</b>					
Arsenic	<b>ND</b>	12	6010B	12-28-11	12-28-11	
Barium	<b>72</b>	3.0	6010B	12-28-11	12-28-11	
Cadmium	<b>ND</b>	0.60	6010B	12-28-11	12-28-11	
Chromium	<b>50</b>	0.60	6010B	12-28-11	12-28-11	
Lead	<b>ND</b>	6.0	6010B	12-28-11	12-28-11	
Mercury	<b>ND</b>	0.30	7471A	12-28-11	12-28-11	
Selenium	<b>ND</b>	12	6010B	12-28-11	12-28-11	
Silver	<b>ND</b>	0.60	6010B	12-28-11	12-28-11	

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-098-13					
Client ID:	DP-6-3.5-4.5					
Arsenic	ND	12	6010B	12-28-11	12-28-11	
Barium	46	2.9	6010B	12-28-11	12-28-11	
Cadmium	ND	0.59	6010B	12-28-11	12-28-11	
Chromium	34	0.59	6010B	12-28-11	12-28-11	
Lead	ND	5.9	6010B	12-28-11	12-28-11	
Mercury	ND	0.29	7471A	12-28-11	12-28-11	
Selenium	ND	12	6010B	12-28-11	12-28-11	
Silver	ND	0.59	6010B	12-28-11	12-28-11	

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1229S1					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	12-29-11	12-29-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	94	68-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-098-18							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				91	95	68-124		



Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1228S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	12-28-11	12-28-11	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	117	50-150				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-098-13					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			103 105	50-150		

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1228S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Chloromethane	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Vinyl Chloride	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Bromomethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Chloroethane	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Trichlorofluoromethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Acetone	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Iodomethane	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Carbon Disulfide	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Methylene Chloride	ND	0.0050	EPA 8260	12-28-11	12-28-11	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,1-Dichloroethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Vinyl Acetate	ND	0.0050	EPA 8260	12-28-11	12-28-11	
2,2-Dichloropropane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
2-Butanone	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Bromochloromethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Chloroform	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Carbon Tetrachloride	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,1-Dichloropropene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Benzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2-Dichloroethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Trichloroethene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2-Dichloropropane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Dibromomethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Bromodichloromethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	12-28-11	12-28-11	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Toluene	ND	0.0050	EPA 8260	12-28-11	12-28-11	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-28-11	12-28-11	

Date of Report: December 29, 2011  
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 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1228S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Tetrachloroethene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,3-Dichloropropane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
2-Hexanone	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Dibromochloromethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2-Dibromoethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Chlorobenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Ethylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
m,p-Xylene	ND	0.0020	EPA 8260	12-28-11	12-28-11	
o-Xylene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Styrene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Bromoform	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Isopropylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Bromobenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	12-28-11	12-28-11	
n-Propylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
2-Chlorotoluene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
4-Chlorotoluene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
tert-Butylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
sec-Butylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
p-Isopropyltoluene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
n-Butylbenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	12-28-11	12-28-11	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
Hexachlorobutadiene	ND	0.0050	EPA 8260	12-28-11	12-28-11	
Naphthalene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>111</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>113</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>91</i>	<i>55-121</i>				

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1228S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0636</b>	<b>0.0633</b>	0.0500	0.0500	127	127	70-130	0	19	
Benzene	<b>0.0547</b>	<b>0.0564</b>	0.0500	0.0500	109	113	70-125	3	15	
Trichloroethene	<b>0.0484</b>	<b>0.0490</b>	0.0500	0.0500	97	98	70-122	1	14	
Toluene	<b>0.0518</b>	<b>0.0512</b>	0.0500	0.0500	104	102	73-120	1	16	
Chlorobenzene	<b>0.0428</b>	<b>0.0431</b>	0.0500	0.0500	86	86	74-109	1	12	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					96	99	63-127			
<i>Toluene-d8</i>					100	100	65-129			
<i>4-Bromofluorobenzene</i>					79	79	55-121			

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 (with silica gel clean-up)

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1228S1					
Naphthalene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>83</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>81</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>79</i>	<i>33 - 119</i>				

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent		Recovery	RPD	RPD	Flags
	MS	MSD	MS	MSD	Result	Recovery	Limits	RPD	Limit		
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-170-20										
	MS	MSD	MS	MSD		MS	MSD				
Naphthalene	<b>0.0522</b>	<b>0.0601</b>	0.0833	0.0833	ND	63	72	39 - 110	14	21	
Acenaphthylene	<b>0.0562</b>	<b>0.0637</b>	0.0833	0.0833	ND	67	76	47 - 124	13	21	
Acenaphthene	<b>0.0572</b>	<b>0.0640</b>	0.0833	0.0833	ND	69	77	50 - 120	11	20	
Fluorene	<b>0.0565</b>	<b>0.0644</b>	0.0833	0.0833	ND	68	77	52 - 126	13	21	
Phenanthrene	<b>0.0547</b>	<b>0.0628</b>	0.0833	0.0833	ND	66	75	41 - 130	14	22	
Anthracene	<b>0.0534</b>	<b>0.0613</b>	0.0833	0.0833	ND	64	74	48 - 124	14	23	
Fluoranthene	<b>0.0567</b>	<b>0.0651</b>	0.0833	0.0833	ND	68	78	40 - 137	14	23	
Pyrene	<b>0.0572</b>	<b>0.0651</b>	0.0833	0.0833	ND	69	78	36 - 139	13	23	
Benzo[a]anthracene	<b>0.0654</b>	<b>0.0754</b>	0.0833	0.0833	ND	79	91	43 - 127	14	21	
Chrysene	<b>0.0564</b>	<b>0.0628</b>	0.0833	0.0833	ND	68	75	41 - 133	11	19	
Benzo[b]fluoranthene	<b>0.0551</b>	<b>0.0630</b>	0.0833	0.0833	ND	66	76	40 - 132	13	25	
Benzo(j,k)fluoranthene	<b>0.0546</b>	<b>0.0606</b>	0.0833	0.0833	ND	66	73	35 - 132	10	25	
Benzo[a]pyrene	<b>0.0559</b>	<b>0.0632</b>	0.0833	0.0833	ND	67	76	37 - 131	12	26	
Indeno(1,2,3-c,d)pyrene	<b>0.0578</b>	<b>0.0638</b>	0.0833	0.0833	ND	69	77	39 - 134	10	23	
Dibenz[a,h]anthracene	<b>0.0590</b>	<b>0.0666</b>	0.0833	0.0833	ND	71	80	40 - 137	12	21	
Benzo[g,h,i]perylene	<b>0.0588</b>	<b>0.0667</b>	0.0833	0.0833	ND	71	80	35 - 135	13	22	
<i>Surrogate:</i>											
<i>2-Fluorobiphenyl</i>						<i>64</i>	<i>71</i>	<i>43 - 109</i>			
<i>Pyrene-d10</i>						<i>70</i>	<i>79</i>	<i>38 - 128</i>			
<i>Terphenyl-d14</i>						<i>68</i>	<i>78</i>	<i>33 - 119</i>			

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1228S1					
Aroclor 1016	ND	0.050	EPA 8082	12-28-11	12-28-11	
Aroclor 1221	ND	0.050	EPA 8082	12-28-11	12-28-11	
Aroclor 1232	ND	0.050	EPA 8082	12-28-11	12-28-11	
Aroclor 1242	ND	0.050	EPA 8082	12-28-11	12-28-11	
Aroclor 1248	ND	0.050	EPA 8082	12-28-11	12-28-11	
Aroclor 1254	ND	0.050	EPA 8082	12-28-11	12-28-11	
Aroclor 1260	ND	0.050	EPA 8082	12-28-11	12-28-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	78		42-123			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-170-20										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.403	0.356	0.500	0.500	ND	81	71	44-125	12	15	
<i>Surrogate:</i>											
DCB						85	78	42-123			

Date of Report: December 29, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098B  
Project: 0180-292-00

**TOTAL METALS  
EPA 6010B/7471A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-28-11  
Date Analyzed: 12-28-11  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB1228S1&MB1228SM2

Analyte	Method	Result	PQL
Arsenic	6010B	ND	10
Barium	6010B	ND	2.5
Cadmium	6010B	ND	0.50
Chromium	6010B	ND	0.50
Lead	6010B	ND	5.0
Mercury	7471A	ND	0.25
Selenium	6010B	ND	10
Silver	6010B	ND	0.50



Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-28-11  
 Date Analyzed: 12-28-11  
  
 Matrix: Soil  
 Units: mg/kg (ppm)  
  
 Lab ID: 12-098-13

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	39.4	44.1	11	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	29.3	28.8	2	0.50	
Lead	ND	ND	NA	5.0	
Mercury	ND	ND	NA	0.25	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	

Date of Report: December 29, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-098B  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-28-11

Date Analyzed: 12-28-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 12-098-13

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>90.3</b>	90	<b>90.0</b>	90	0	
Barium	100	<b>133</b>	94	<b>132</b>	92	1	
Cadmium	50.0	<b>46.5</b>	93	<b>46.4</b>	93	0	
Chromium	100	<b>123</b>	94	<b>117</b>	88	5	
Lead	250	<b>224</b>	90	<b>223</b>	89	1	
Mercury	0.500	<b>0.506</b>	101	<b>0.485</b>	97	4	
Selenium	100	<b>90.2</b>	90	<b>90.8</b>	91	1	
Silver	25.0	<b>19.7</b>	79	<b>19.7</b>	79	0	

Date of Report: December 29, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-098B  
Project: 0180-292-00

**% MOISTURE**

Date Analyzed: 12-28-11

Client ID	Lab ID	% Moisture
DP-3-3-4	12-098-05	14
DP-4-5-6	12-098-08	22
DP-5-6-7	12-098-10	17
DP-6-3.5-4.5	12-098-13	15
DP-7-5-6	12-098-16	18
DP-8-2-3	12-098-18	12
DP-9-6-7	12-098-21	19



### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference



**OnSite Environmental Inc.**  
 Analytical Laboratory Testing Services  
 14648 NE 95th Street • Redmond, WA 98052  
 Phone: (425) 863-3881 • www.onsite-env.com

# Chain of Custody

Turnaround Request  
(in working days)

Laboratory Number:

**12-098**

(Check One)

- Same Day     1 Day  
 2 Days     3 Days  
 Standard (7 Days) (TPH analysis 5 Days)

(other)

Company: GeoEngineers/WS DOT  
 Project Number: 0180-292-00  
 Project Name: Milway Metals  
 Project Manager: Aaron Wagoner  
 Sampled by: [Signature]

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	No. of Cont.	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx	Volatiles 8260B	Halogenated Volatiles 8260B	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHs 8270D/SIM (low-level)	PCBs 8082	Organochlorine Pesticides 8081A	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664	% Moisture
1	DP-1-0-2	12/13/11	0920	Soil	5			X	X	X		X	X	X				X	X			X
2	DP-1-7-8		0940					X	X	X		X	X	X				X	X			X
3	DP-1-13-15		0930					X	X	X		X	X	X				X	X			X
4	DP-3-0-2		1055					X	X	X		X	X	X				X	X			X
5	DP-3-3-4		1120					X	X	X		X	X	X				X	X			X
6	DP-3-9-10		1146					X	X	X		X	X	X				X	X			X
7	DP-4-6-2		1315					X	X	X		X	X	X				X	X			X
8	DP-4-5-6		1325					X	X	X		X	X	X				X	X			X
9	DP-5-0-2		1435					X	X	X		X	X	X				X	X			X
10	DP-5-6-7		1445					X	X	X		X	X	X				X	X			X

Signature: [Signature] Company: Geo Date: 12/14/11 Time: 1348 Comments/Special Instructions: Added 12/15/11 m. DB (STA)

Relinquished: [Signature]  
 Received: [Signature]  
 Relinquished: [Signature]  
 Received: [Signature]  
 Relinquished: [Signature]  
 Received: [Signature]  
 Reviewed/Date: \_\_\_\_\_









14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

February 24, 2012

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1202-085B

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on February 9, 2012.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures



Date of Report: February 24, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085B  
Project: 0180-292-00

#### **Case Narrative**

Sample was collected on February 9, 2012 and received by the laboratory on February 9, 2012. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### PAHs by EPA 8270D/SIM Analysis

Sample HA4-0.0-2.0 was extracted outside of the holding time.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Date of Report: February 24, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-085B  
Project: 0180-292-00

#### ANALYTICAL REPORT FOR SAMPLES

Client ID	Laboratory ID	Matrix	Date Sampled	Date Received	Notes
HA4-0.0-2.0	02-085-31	Soil	2-9-12	2-9-12	

Date of Report: February 24, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085B  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>HA4-0.0-2.0</b>					
Laboratory ID:	02-085-31					
Naphthalene	<b>0.021</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
2-Methylnaphthalene	<b>0.019</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
1-Methylnaphthalene	<b>0.010</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Acenaphthylene	<b>ND</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Acenaphthene	<b>ND</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Fluorene	<b>ND</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Phenanthrene	<b>0.031</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Anthracene	<b>ND</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Fluoranthene	<b>0.036</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Pyrene	<b>0.038</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[a]anthracene	<b>0.013</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Chrysene	<b>0.017</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[b]fluoranthene	<b>0.018</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[a]pyrene	<b>0.013</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Indeno(1,2,3-c,d)pyrene	<b>0.013</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Dibenz[a,h]anthracene	<b>ND</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[g,h,i]perylene	<b>0.019</b>	0.0084	EPA 8270/SIM	2-23-12	2-24-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>47</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>75</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>49</i>	<i>33 - 119</i>				

Date of Report: February 24, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085B  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0223S1					
Naphthalene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Acenaphthene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Fluorene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Phenanthrene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Anthracene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Fluoranthene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Pyrene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Chrysene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	2-23-12	2-24-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>70</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>115</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>102</i>	<i>33 - 119</i>				

Date of Report: February 24, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-085B  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL  
 (with silica gel clean-up)**

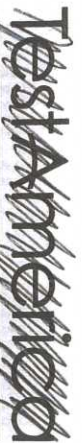
Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0223S1									
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	<b>0.0736</b>	<b>0.0781</b>	0.0833	0.0833	88	94	43 - 108	6	27	
Acenaphthylene	<b>0.0671</b>	<b>0.0701</b>	0.0833	0.0833	81	84	52 - 120	4	21	
Acenaphthene	<b>0.0615</b>	<b>0.0653</b>	0.0833	0.0833	74	78	59 - 113	6	17	
Fluorene	<b>0.0659</b>	<b>0.0720</b>	0.0833	0.0833	79	86	64 - 117	9	14	
Phenanthrene	<b>0.0613</b>	<b>0.0658</b>	0.0833	0.0833	74	79	67 - 112	7	12	
Anthracene	<b>0.0678</b>	<b>0.0728</b>	0.0833	0.0833	81	87	59 - 110	7	16	
Fluoranthene	<b>0.0790</b>	<b>0.0849</b>	0.0833	0.0833	95	102	68 - 120	7	15	
Pyrene	<b>0.0771</b>	<b>0.0819</b>	0.0833	0.0833	93	98	66 - 121	6	17	
Benzo[a]anthracene	<b>0.0734</b>	<b>0.0781</b>	0.0833	0.0833	88	94	63 - 114	6	12	
Chrysene	<b>0.0666</b>	<b>0.0708</b>	0.0833	0.0833	80	85	67 - 118	6	12	
Benzo[b]fluoranthene	<b>0.0645</b>	<b>0.0695</b>	0.0833	0.0833	77	83	58 - 127	7	20	
Benzo(j,k)fluoranthene	<b>0.0617</b>	<b>0.0670</b>	0.0833	0.0833	74	80	42 - 134	8	26	
Benzo[a]pyrene	<b>0.0675</b>	<b>0.0720</b>	0.0833	0.0833	81	86	55 - 111	6	19	
Indeno(1,2,3-c,d)pyrene	<b>0.0670</b>	<b>0.0718</b>	0.0833	0.0833	80	86	60 - 125	7	20	
Dibenz[a,h]anthracene	<b>0.0727</b>	<b>0.0771</b>	0.0833	0.0833	87	93	62 - 125	6	19	
Benzo[g,h,i]perylene	<b>0.0712</b>	<b>0.0757</b>	0.0833	0.0833	85	91	61 - 124	6	19	
<i>Surrogate:</i>										
2-Fluorobiphenyl					71	75	43 - 109			
Pyrene-d10					113	121	38 - 128			
Terphenyl-d14					96	103	33 - 119			



### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference



THE LEADER IN ENVIRONMENTAL TESTING  
On SITE Environmental

TestAmerica Seattle  
5755 8th Street E.  
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www.testamericainc.com

Rush 4 Days  
 Short Hold

Chain of  
Custody Record

Client: Geo Engineers Client Contact: Aaron Wegman Date: 2/9/12 Chain of Custody Number: 12706

Address: 1101 S. Everett Ave Ste 200 Telephone Number (Area Code): (253) 383 4940 Lab Number: \_\_\_\_\_ Page 1 of 3

City: Tacoma State: WA Zip Code: 98402 Project Name and Location (State): WSDOT McKinley Metals, Security, WA Billing Contact: Aaron Wegman Lab Contact: Daniel Brunsen

Contract/Purchase Order/Quote No.: 0180-292-00 Matrix: \_\_\_\_\_ Containers & Preservatives: \_\_\_\_\_

Sample I.D. and Location/Description (Containers for each sample may be combined on one line)	Date	Time	Matrix			Containers & Preservatives						Analysis (Attach list if more space is needed)													
			Air	Aqueous	Sed.	Soil	Unpres.	H2SO4	HNO3	HCl	NaOH	ZnAc/NaOH	NWTPH-Dt w/ silica gel cleanup	NWTPH-Gx	VOCs Lowlevel	SVOCs Lowlevel	PAHs Lowlevel	Dissolved RCRA metals	Total RCRA metals	PCB	Lead	RCRA metals	HEX CHROME	EMERSON	
1 DP12-0.0-2.0	2/8/12	1100				X																			
2 DP13-0.0-1.0g		1125				X																			
3 DP14-0.0-2.0		1155				X																			
4 DP15-0.0-2.0		1335				X																			
5 DP16-0.0-2.0		1355				X																			
6 DP17-0.0-2.0		1420				X																			
7 DP18-0.0-2.0		1520				X																			
8 DP19-0.0-2.0		1530				X																			
9 DP20-0.0-2.0		1540				X																			
10 DP21-0.0-2.0		1615				X																			
11 DP12-2.5-3.5		1105				X																			
12 DP12-5.0-7.5		1110				X																			

Cooler:  Yes  No Cooler Temp: \_\_\_\_\_ Possible Hazard Identification:  Non-Hazard  Flammable  Skin Irritant  Poison B  Unknown

Turn Around Time Required (business days):  24 Hours  48 Hours  5 Days  10 Days  15 Days  Other \_\_\_\_\_ QC Requirements (Specify): \_\_\_\_\_

1. Relinquished By Sign/Print: Aaron Wegman Date: 2/9/12 Time: 1630 1. Received By Sign/Print: M VORND Date: 2/9/12 Time: 1636

2. Relinquished By Sign/Print: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ 2. Received By Sign/Print: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

3. Relinquished By Sign/Print: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_ 3. Received By Sign/Print: \_\_\_\_\_ Date: \_\_\_\_\_ Time: \_\_\_\_\_

Comments: \_\_\_\_\_ Distribution: WHITE - Stays with the Samples; CANARY - Returned to Client with Report; PINK - Field Copy











14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

December 23, 2011

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1112-099

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on December 14, 2011.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal flourish extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-099  
Project: 0180-292-00

### **Case Narrative**

Samples were collected on December 13, 2011 and received by the laboratory on December 14, 2011. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### NWTPH Gx and Volatiles EPA 8260B (soil) Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-099  
Project: 0180-292-00

**ANALYTICAL REPORT FOR SAMPLES**

<b>Client ID</b>	<b>Laboratory ID</b>	<b>Matrix</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Notes</b>
SW2	12-099-01	Water	12-13-11	12-14-11	
SED3	12-099-03	Soil	12-13-11	12-14-11	
SED4	12-099-04	Soil	12-13-11	12-14-11	
SED5	12-099-05	Soil	12-13-11	12-14-11	
SED6	12-099-06	Soil	12-13-11	12-14-11	
SED7	12-099-07	Soil	12-13-11	12-14-11	
SED8	12-099-08	Soil	12-13-11	12-14-11	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-099  
Project: 0180-292-00

**NWTPH-Gx**

Matrix: Water  
Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SW2</b>					
Laboratory ID:	12-099-01					
Gasoline	<b>ND</b>	100	NWTPH-Gx	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	92	73-121				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

### NWTPH-Gx

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED3</b>					
Laboratory ID:	12-099-03					
Gasoline	<b>ND</b>	9.0	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	98	68-124				
<b>Client ID:</b>	<b>SED4</b>					
Laboratory ID:	12-099-04					
Gasoline	<b>ND</b>	9.6	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	96	68-124				
<b>Client ID:</b>	<b>SED5</b>					
Laboratory ID:	12-099-05					
Gasoline	<b>ND</b>	7.6	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	96	68-124				
<b>Client ID:</b>	<b>SED6</b>					
Laboratory ID:	12-099-06					
Gasoline	<b>ND</b>	9.6	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	104	68-124				
<b>Client ID:</b>	<b>SED7</b>					
Laboratory ID:	12-099-07					
Gasoline	<b>ND</b>	5.3	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	98	68-124				
<b>Client ID:</b>	<b>SED8</b>					
Laboratory ID:	12-099-08					
Gasoline	<b>ND</b>	8.0	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	99	68-124				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**NWTPH-Dx**  
**(with acid/silica gel clean-up)**

Matrix: Water  
 Units: mg/L (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SW2</b>					
Laboratory ID:	12-099-01					
Diesel Range Organics	<b>ND</b>	0.29	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	0.46	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>71</i>	<i>50-150</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**NWTPH-Dx**  
**(with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED3</b>					
Laboratory ID:	12-099-03					
Diesel Range Organics	<b>ND</b>	38	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	77	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	105	50-150				
<b>Client ID:</b>	<b>SED4</b>					
Laboratory ID:	12-099-04					
Diesel Range Organics	<b>ND</b>	40	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil	<b>270</b>	79	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	109	50-150				
<b>Client ID:</b>	<b>SED5</b>					
Laboratory ID:	12-099-05					
Diesel Range Organics	<b>ND</b>	120	NWTPH-Dx	12-19-11	12-19-11	U1
Lube Oil	<b>950</b>	66	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	107	50-150				
<b>Client ID:</b>	<b>SED6</b>					
Laboratory ID:	12-099-06					
Diesel Range Organics	<b>ND</b>	37	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	75	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	103	50-150				
<b>Client ID:</b>	<b>SED7</b>					
Laboratory ID:	12-099-07					
Diesel Range Organics	<b>ND</b>	28	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	57	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	109	50-150				
<b>Client ID:</b>	<b>SED8</b>					
Laboratory ID:	12-099-08					
Diesel Range Organics	<b>ND</b>	33	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	65	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	103	50-150				



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
 Page 1 of 2

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW2</b>					
Laboratory ID:	12-099-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Acetone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	2.0	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Toluene	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
 Page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW2</b>					
Laboratory ID:	12-099-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.40	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	1.0	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>84</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>84</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>86</i>	<i>65-120</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
 Page 1 of 2

Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SED3</b>					
Laboratory ID:	12-099-03					
Dichlorodifluoromethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0064	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0064	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0064	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0064	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0013	EPA 8260	12-15-11	12-15-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED3</b>					
Laboratory ID:	12-099-03					
1,1,2-Trichloroethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0026	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0013	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0064	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0064	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0013	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>93</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>97</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>90</i>	<i>55-121</i>				

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED4</b>					
Laboratory ID:	12-099-04					
Dichlorodifluoromethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0089	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0089	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0089	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0089	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0018	EPA 8260	12-15-11	12-15-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED4</b>					
Laboratory ID:	12-099-04					
1,1,2-Trichloroethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0036	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0018	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0089	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0089	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0018	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>89</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>94</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>74</i>	<i>55-121</i>				

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 Project: 0180-292-00

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SED5</b>					
Laboratory ID:	12-099-05					
Dichlorodifluoromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0068	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0068	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0068	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0068	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED5</b>					
Laboratory ID:	12-099-05					
1,1,2-Trichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0027	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0068	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0068	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>87</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>91</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>74</i>	<i>55-121</i>				



Date of Report: December 23, 2011  
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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED6</b>					
Laboratory ID:	12-099-06					
Dichlorodifluoromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0069	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0069	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0069	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0069	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED6</b>					
Laboratory ID:	12-099-06					
1,1,2-Trichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0027	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>84</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>87</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>72</i>	<i>55-121</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

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Matrix: Soil  
 Units: mg/kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SED7</b>					
Laboratory ID:	12-099-07					
Dichlorodifluoromethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0049	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0049	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0049	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0049	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.00097	EPA 8260	12-15-11	12-15-11	

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 Project: 0180-292-00

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED7</b>					
Laboratory ID:	12-099-07					
1,1,2-Trichloroethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0019	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.00097	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0049	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0049	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.00097	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>91</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>98</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>94</i>	<i>55-121</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED8</b>					
Laboratory ID:	12-099-08					
Dichlorodifluoromethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0059	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0059	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0059	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0059	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0012	EPA 8260	12-15-11	12-15-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED8</b>					
Laboratory ID:	12-099-08					
1,1,2-Trichloroethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0024	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0012	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0059	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0059	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0012	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>95</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>98</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>82</i>	<i>55-121</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW2</b>					
Laboratory ID:	12-099-01					
n-Nitrosodimethylamine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Pyridine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Phenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
Aniline	ND	4.8	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Chlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,3-Dichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,4-Dichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Benzyl alcohol	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2-Dichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.95	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.95	EPA 8270	12-16-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.95	EPA 8270	12-16-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Hexachloroethane	ND	0.95	EPA 8270	12-16-11	12-20-11	
Nitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Isophorone	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Nitrophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4-Dimethylphenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4-Dichlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Naphthalene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
4-Chloroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
Hexachlorobutadiene	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Methylnaphthalene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Hexachlorocyclopentadiene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,3-Dichloroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Chloronaphthalene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Nitroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,4-Dinitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Dimethylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,3-Dinitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,6-Dinitrotoluene	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2-Dinitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Acenaphthylene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
3-Nitroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	

Date of Report: December 23, 2011  
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 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW2</b>					
Laboratory ID:	12-099-01					
2,4-Dinitrophenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
Acenaphthene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
4-Nitrophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4-Dinitrotoluene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Dibenzofuran	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
Diethylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Nitroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
Fluorene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
4,6-Dinitro-2-methylphenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.95	EPA 8270	12-16-11	12-20-11	
Hexachlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Pentachlorophenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
Phenanthrene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Carbazole	ND	0.95	EPA 8270	12-16-11	12-20-11	
Di-n-butylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
Fluoranthene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Benzidine	ND	4.8	EPA 8270	12-16-11	12-20-11	
Pyrene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Butylbenzylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
bis-2-Ethylhexyladipate	ND	4.8	EPA 8270	12-16-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Benzo[a]anthracene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
bis(2-Ethylhexyl)phthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
Di-n-octylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Indeno[1,2,3-cd]pyrene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	32	18 - 86				
Phenol-d6	23	10 - 88				
Nitrobenzene-d5	51	37 - 112				
2-Fluorobiphenyl	66	42 - 108				
2,4,6-Tribromophenol	79	39 - 118				
Terphenyl-d14	61	49 - 122				



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED5</b>					
Laboratory ID:	12-099-05					
n-Nitrosodimethylamine	ND	0.22	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	2.2	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.22	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.22	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.22	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.22	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.22	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.22	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.22	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.22	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	2.2	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
Naphthalene	<b>0.093</b>	0.044	EPA 8270/SIM	12-19-11	12-20-11	
4-Chloroaniline	ND	0.22	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.22	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.094</b>	0.044	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	ND	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Hexachlorocyclopentadiene	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.22	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	<b>0.061</b>	0.044	EPA 8270/SIM	12-19-11	12-20-11	
3-Nitroaniline	ND	0.22	EPA 8270	12-19-11	12-20-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED5</b>					
Laboratory ID:	12-099-05					
2,4-Dinitrophenol	ND	1.1	EPA 8270	12-19-11	12-20-11	
Acenaphthene	ND	0.044	EPA 8270/SIM	12-19-11	12-20-11	
4-Nitrophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.22	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.22	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	1.1	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.22	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.22	EPA 8270	12-19-11	12-20-11	
Fluorene	ND	0.044	EPA 8270/SIM	12-19-11	12-20-11	
4,6-Dinitro-2-methylphenol	ND	1.1	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.22	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.22	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.22	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.22	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	1.1	EPA 8270	12-19-11	12-20-11	
Phenanthrene	0.19	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	0.078	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Carbazole	ND	0.22	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	2.2	EPA 8270	12-19-11	12-20-11	
Fluoranthene	0.30	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Benzidine	ND	2.2	EPA 8270	12-19-11	12-20-11	
Pyrene	0.29	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Butylbenzylphthalate	ND	2.2	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	1.1	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	2.2	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	0.16	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	0.20	0.044	EPA 8270/SIM	12-19-11	12-20-11	
bis(2-Ethylhexyl)phthalate	ND	1.1	EPA 8270	12-19-11	12-20-11	
Di-n-octylphthalate	ND	0.22	EPA 8270	12-19-11	12-20-11	
Benzo[b]fluoranthene	0.33	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	0.11	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	0.27	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Indeno[1,2,3-cd]pyrene	0.20	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	0.049	0.044	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	0.23	0.044	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	67	30 - 97				
Phenol-d6	72	40 - 104				
Nitrobenzene-d5	64	35 - 102				
2-Fluorobiphenyl	95	44 - 97				
2,4,6-Tribromophenol	136	41 - 110				
Terphenyl-d14	84	53 - 107				

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Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED6</b>					
Laboratory ID:	12-099-06					
n-Nitrosodimethylamine	ND	0.050	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	0.50	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.050	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.050	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	<b>0.064</b>	0.050	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.050	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.050	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.050	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.050	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.050	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	0.50	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
Naphthalene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
4-Chloroaniline	ND	0.050	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.050	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Hexachlorocyclopentadiene	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.050	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
3-Nitroaniline	ND	0.050	EPA 8270	12-19-11	12-20-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED6</b>					
Laboratory ID:	12-099-06					
2,4-Dinitrophenol	ND	0.25	EPA 8270	12-19-11	12-20-11	
Acenaphthene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
4-Nitrophenol	0.13	0.050	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.050	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.050	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	0.25	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.050	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.050	EPA 8270	12-19-11	12-20-11	
Fluorene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
4,6-Dinitro-2-methylphenol	ND	0.25	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.050	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.050	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.050	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.050	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	0.25	EPA 8270	12-19-11	12-20-11	
Phenanthrene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Carbazole	ND	0.050	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	0.50	EPA 8270	12-19-11	12-20-11	
Fluoranthene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Benzidine	ND	0.50	EPA 8270	12-19-11	12-20-11	
Pyrene	0.093	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Butylbenzylphthalate	ND	0.50	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	0.25	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.50	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
bis(2-Ethylhexyl)phthalate	ND	0.25	EPA 8270	12-19-11	12-20-11	
Di-n-octylphthalate	ND	0.050	EPA 8270	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Indeno[1,2,3-cd]pyrene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.050	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	61	30 - 97				
Phenol-d6	59	40 - 104				
Nitrobenzene-d5	56	35 - 102				
2-Fluorobiphenyl	70	44 - 97				
2,4,6-Tribromophenol	100	41 - 110				
Terphenyl-d14	66	53 - 107				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED3</b>					
Laboratory ID:	12-099-03					
Naphthalene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>54</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>60</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>53</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SED4</b>					
Laboratory ID:	12-099-04					
Naphthalene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.011	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>81</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>91</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>92</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
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 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SED7</b>					
Laboratory ID:	12-099-07					
Naphthalene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	ND	0.0075	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>64</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>76</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>74</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
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**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED8</b>					
Laboratory ID:	12-099-08					
Naphthalene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	ND	0.0087	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>60</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>67</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>65</i>	<i>33 - 119</i>				



Date of Report: December 23, 2011  
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### PCBs by EPA 8082

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW2</b>					
Laboratory ID:	12-099-01					
Aroclor 1016	<b>ND</b>	0.048	EPA 8082	12-16-11	12-16-11	
Aroclor 1221	<b>ND</b>	0.048	EPA 8082	12-16-11	12-16-11	
Aroclor 1232	<b>ND</b>	0.048	EPA 8082	12-16-11	12-16-11	
Aroclor 1242	<b>ND</b>	0.048	EPA 8082	12-16-11	12-16-11	
Aroclor 1248	<b>ND</b>	0.048	EPA 8082	12-16-11	12-16-11	
Aroclor 1254	<b>ND</b>	0.048	EPA 8082	12-16-11	12-16-11	
Aroclor 1260	<b>ND</b>	0.048	EPA 8082	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>80</i>	<i>36-127</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED3</b>					
Laboratory ID:	12-099-03					
Aroclor 1016	ND	0.076	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	ND	0.076	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	ND	0.076	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	ND	0.076	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	ND	0.076	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	ND	0.076	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	ND	0.076	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	73	42-123				
<b>Client ID:</b>	<b>SED4</b>					
Laboratory ID:	12-099-04					
Aroclor 1016	ND	0.079	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	ND	0.079	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	ND	0.079	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	ND	0.079	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	ND	0.079	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	ND	0.079	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	ND	0.079	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	79	42-123				
<b>Client ID:</b>	<b>SED5</b>					
Laboratory ID:	12-099-05					
Aroclor 1016	ND	0.066	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	ND	0.066	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	ND	0.066	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	ND	0.066	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	ND	0.066	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	0.16	0.066	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	ND	0.066	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	85	42-123				

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 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED6</b>					
Laboratory ID:	12-099-06					
Aroclor 1016	ND	0.075	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	ND	0.075	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	ND	0.075	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	ND	0.075	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	ND	0.075	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	ND	0.075	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	ND	0.075	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	75	42-123				
<b>Client ID:</b>	<b>SED7</b>					
Laboratory ID:	12-099-07					
Aroclor 1016	ND	0.057	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	ND	0.057	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	ND	0.057	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	ND	0.057	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	ND	0.057	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	ND	0.057	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	ND	0.057	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	79	42-123				
<b>Client ID:</b>	<b>SED8</b>					
Laboratory ID:	12-099-08					
Aroclor 1016	ND	0.065	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	ND	0.065	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	ND	0.065	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	ND	0.065	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	ND	0.065	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	ND	0.065	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	ND	0.065	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	82	42-123				

Date of Report: December 23, 2011  
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**TOTAL METALS  
 EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	12-099-01					
<b>Client ID:</b>	<b>SW2</b>					
Arsenic	<b>ND</b>	3.3	200.8	12-16-11	12-16-11	
Barium	<b>31</b>	28	200.8	12-16-11	12-16-11	
Cadmium	<b>ND</b>	4.4	200.8	12-16-11	12-16-11	
Chromium	<b>ND</b>	11	200.8	12-16-11	12-16-11	
Lead	<b>4.0</b>	1.1	200.8	12-16-11	12-16-11	
Mercury	<b>ND</b>	0.50	7470A	12-16-11	12-16-11	
Selenium	<b>ND</b>	5.6	200.8	12-16-11	12-16-11	
Silver	<b>ND</b>	11	200.8	12-16-11	12-16-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
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**DISSOLVED METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-099-01					
Client ID:	SW2					
Arsenic	ND	3.0	200.8	12-14-11	12-15-11	
Barium	ND	25	200.8	12-14-11	12-15-11	
Cadmium	ND	4.0	200.8	12-14-11	12-15-11	
Chromium	ND	10	200.8	12-14-11	12-15-11	
Lead	1.4	1.0	200.8	12-14-11	12-15-11	
Mercury	ND	0.50	7470A	12-14-11	12-20-11	
Selenium	ND	5.0	200.8	12-14-11	12-15-11	
Silver	ND	10	200.8	12-14-11	12-15-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-099-03					
<b>Client ID:</b>	<b>SED3</b>					
Arsenic	ND	15	6010B	12-19-11	12-19-11	
Barium	64	3.8	6010B	12-19-11	12-19-11	
Cadmium	ND	0.76	6010B	12-19-11	12-19-11	
Chromium	41	0.76	6010B	12-19-11	12-19-11	
Lead	ND	7.6	6010B	12-19-11	12-19-11	
Mercury	ND	0.38	7471A	12-19-11	12-19-11	
Selenium	ND	15	6010B	12-19-11	12-19-11	
Silver	ND	0.76	6010B	12-19-11	12-19-11	

Lab ID:	12-099-04					
<b>Client ID:</b>	<b>SED4</b>					
Arsenic	ND	16	6010B	12-19-11	12-19-11	
Barium	88	4.0	6010B	12-19-11	12-19-11	
Cadmium	ND	0.79	6010B	12-19-11	12-19-11	
Chromium	48	0.79	6010B	12-19-11	12-19-11	
Lead	39	7.9	6010B	12-19-11	12-19-11	
Mercury	ND	0.4	7471A	12-19-11	12-19-11	
Selenium	ND	16	6010B	12-19-11	12-19-11	
Silver	ND	0.79	6010B	12-19-11	12-19-11	

Date of Report: December 23, 2011  
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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-099-05					
<b>Client ID:</b>	<b>SED5</b>					
Arsenic	<b>ND</b>	13	6010B	12-19-11	12-19-11	
Barium	<b>75</b>	3.3	6010B	12-19-11	12-19-11	
Cadmium	<b>0.89</b>	0.66	6010B	12-19-11	12-19-11	
Chromium	<b>57</b>	0.66	6010B	12-19-11	12-19-11	
Lead	<b>120</b>	6.6	6010B	12-19-11	12-19-11	
Mercury	<b>ND</b>	0.33	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	13	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.66	6010B	12-19-11	12-19-11	

Lab ID:	12-099-06					
<b>Client ID:</b>	<b>SED6</b>					
Arsenic	<b>ND</b>	15	6010B	12-19-11	12-19-11	
Barium	<b>100</b>	3.7	6010B	12-19-11	12-19-11	
Cadmium	<b>0.90</b>	0.75	6010B	12-19-11	12-19-11	
Chromium	<b>37</b>	0.75	6010B	12-19-11	12-19-11	
Lead	<b>13</b>	7.5	6010B	12-19-11	12-19-11	
Mercury	<b>ND</b>	0.37	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	15	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.75	6010B	12-19-11	12-19-11	

Date of Report: December 23, 2011  
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 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-099-07					
<b>Client ID:</b>	<b>SED7</b>					
Arsenic	ND	11	6010B	12-19-11	12-19-11	
Barium	41	2.8	6010B	12-19-11	12-19-11	
Cadmium	ND	0.57	6010B	12-19-11	12-19-11	
Chromium	30	0.57	6010B	12-19-11	12-19-11	
Lead	ND	5.7	6010B	12-19-11	12-19-11	
Mercury	ND	0.28	7471A	12-19-11	12-19-11	
Selenium	ND	11	6010B	12-19-11	12-19-11	
Silver	ND	0.57	6010B	12-19-11	12-19-11	

Lab ID:	12-099-08					
<b>Client ID:</b>	<b>SED8</b>					
Arsenic	ND	13	6010B	12-19-11	12-19-11	
Barium	78	3.3	6010B	12-19-11	12-19-11	
Cadmium	ND	0.65	6010B	12-19-11	12-19-11	
Chromium	29	0.65	6010B	12-19-11	12-19-11	
Lead	ND	6.5	6010B	12-19-11	12-19-11	
Mercury	ND	0.33	7471A	12-19-11	12-19-11	
Selenium	ND	13	6010B	12-19-11	12-19-11	
Silver	ND	0.65	6010B	12-19-11	12-19-11	



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1215W2					
Gasoline	<b>ND</b>	100	NWTPH-Gx	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	73-121				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-100-01							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				91	90	73-121		

Date of Report: December 23, 2011  
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**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1216S2					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	87	68-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-100-03							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				102	106	68-124		

Date of Report: December 23, 2011  
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**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1219S1					
Diesel Range Organics	<b>ND</b>	25	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	50	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>120</i>	<i>50-150</i>				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-110-01					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			109 114	50-150		

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1221W1					
Diesel Range Organics	<b>ND</b>	0.25	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	0.40	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>101</i>	<i>50-150</i>				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-105-02					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			100 105	50-150		

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1215W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Acetone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	2.0	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Toluene	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:		MB1215W1				
1,1,2-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.40	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	1.0	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>85</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>90</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>88</i>	<i>65-120</i>				

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**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1215W1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>11.0</b>	<b>10.4</b>	10.0	10.0	110	104	70-130	6	11	
Benzene	<b>9.09</b>	<b>8.95</b>	10.0	10.0	91	90	75-123	2	8	
Trichloroethene	<b>10.1</b>	<b>10.3</b>	10.0	10.0	101	103	80-113	2	9	
Toluene	<b>9.74</b>	<b>9.71</b>	10.0	10.0	97	97	80-113	0	8	
Chlorobenzene	<b>10.8</b>	<b>10.5</b>	10.0	10.0	108	105	80-111	3	8	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					<i>77</i>	<i>77</i>	<i>68-120</i>			
<i>Toluene-d8</i>					<i>82</i>	<i>85</i>	<i>73-120</i>			
<i>4-Bromofluorobenzene</i>					<i>79</i>	<i>85</i>	<i>65-120</i>			

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1215S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0050	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0050	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0050	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-15-11	12-15-11	



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**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1215S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0020	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>93</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>98</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>96</i>	<i>55-121</i>				

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 Project: 0180-292-00

**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1215S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0573</b>	<b>0.0584</b>	0.0500	0.0500	115	117	70-130	2	19	
Benzene	<b>0.0499</b>	<b>0.0504</b>	0.0500	0.0500	100	101	70-125	1	15	
Trichloroethene	<b>0.0474</b>	<b>0.0476</b>	0.0500	0.0500	95	95	70-122	0	14	
Toluene	<b>0.0493</b>	<b>0.0495</b>	0.0500	0.0500	99	99	73-120	0	16	
Chlorobenzene	<b>0.0451</b>	<b>0.0451</b>	0.0500	0.0500	90	90	74-109	0	12	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					<i>85</i>	<i>84</i>	<i>63-127</i>			
<i>Toluene-d8</i>					<i>94</i>	<i>90</i>	<i>65-129</i>			
<i>4-Bromofluorobenzene</i>					<i>91</i>	<i>84</i>	<i>55-121</i>			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216W2					
n-Nitrosodimethylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Pyridine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Phenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
Aniline	ND	5.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethyl)ether	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Chlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,3-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,4-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzyl alcohol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	1.0	EPA 8270	12-16-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	1.0	EPA 8270	12-16-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachloroethane	ND	1.0	EPA 8270	12-16-11	12-20-11	
Nitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Isophorone	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Nitrophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dimethylphenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2,4-Trichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Naphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4-Chloroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachlorobutadiene	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Chloro-3-methylphenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Hexachlorocyclopentadiene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4,6-Trichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3-Dichloroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4,5-Trichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Chloronaphthalene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,4-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Dimethylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,3-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,6-Dinitrotoluene	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Acenaphthylene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
3-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216W2					
2,4-Dinitrophenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
Acenaphthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4-Nitrophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dinitrotoluene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Dibenzofuran	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
Diethylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Chlorophenyl-phenylether	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
Fluorene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4,6-Dinitro-2-methylphenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
n-Nitrosodiphenylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Diphenylhydrazine	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Bromophenyl-phenylether	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Pentachlorophenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
Phenanthrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Carbazole	ND	1.0	EPA 8270	12-16-11	12-20-11	
Di-n-butylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Fluoranthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Benzidine	ND	5.0	EPA 8270	12-16-11	12-20-11	
Pyrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Butylbenzylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis-2-Ethylhexyladipate	ND	5.0	EPA 8270	12-16-11	12-20-11	
3,3'-Dichlorobenzidine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzo[a]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
bis(2-Ethylhexyl)phthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Di-n-octylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzo[b]fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Indeno[1,2,3-cd]pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	57	18 - 86				
Phenol-d6	41	10 - 88				
Nitrobenzene-d5	76	37 - 112				
2-Fluorobiphenyl	92	42 - 108				
2,4,6-Tribromophenol	108	39 - 118				
Terphenyl-d14	83	49 - 122				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
	SB	SBD	SB	SBD	SB	SBD				
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1216W2									
Phenol	15.5	13.3	40.0	40.0	39	33	26 - 60	15	29	
2-Chlorophenol	29.6	24.4	40.0	40.0	74	61	46 - 104	19	34	
1,4-Dichlorobenzene	13.0	10.6	20.0	20.0	65	53	46 - 92	20	29	
n-Nitroso-di-n-propylamine	12.1	10.2	20.0	20.0	61	51	30 - 102	17	25	
1,2,4-Trichlorobenzene	18.3	15.4	20.0	20.0	92	77	45 - 92	17	25	
4-Chloro-3-methylphenol	34.4	32.1	40.0	40.0	86	80	53 - 104	7	18	
Acenaphthene	16.0	14.6	20.0	20.0	80	73	57 - 95	9	15	
4-Nitrophenol	22.4	20.8	40.0	40.0	56	52	21 - 75	7	33	
2,4-Dinitrotoluene	21.4	20.3	20.0	20.0	107	102	60 - 108	5	20	
Pentachlorophenol	26.3	25.2	40.0	40.0	66	63	48 - 119	4	31	
Pyrene	13.8	13.3	20.0	20.0	69	67	62 - 111	4	19	
<i>Surrogate:</i>										
2-Fluorophenol					58	46	18 - 86			
Phenol-d6					42	34	10 - 88			
Nitrobenzene-d5					73	61	37 - 112			
2-Fluorobiphenyl					90	79	42 - 108			
2,4,6-Tribromophenol					103	100	39 - 118			
Terphenyl-d14					80	78	49 - 122			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S2					
n-Nitrosodimethylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
Pyridine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Phenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
Aniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Chlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,3-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,4-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Benzyl alcohol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Dichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.033	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.033	EPA 8270	12-19-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.033	EPA 8270	12-19-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachloroethane	ND	0.033	EPA 8270	12-19-11	12-20-11	
Nitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Isophorone	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Nitrophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dimethylphenol	ND	0.33	EPA 8270	12-19-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Naphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4-Chloroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachlorobutadiene	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Hexachlorocyclopentadiene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3-Dichloroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Chloronaphthalene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,4-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Dimethylphthalate	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,3-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,6-Dinitrotoluene	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Dinitrobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
3-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S2					
2,4-Dinitrophenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4-Nitrophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,4-Dinitrotoluene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Dibenzofuran	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.033	EPA 8270	12-19-11	12-20-11	
Diethylphthalate	ND	0.17	EPA 8270	12-19-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Nitroaniline	ND	0.033	EPA 8270	12-19-11	12-20-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
4,6-Dinitro-2-methylphenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.033	EPA 8270	12-19-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.033	EPA 8270	12-19-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.033	EPA 8270	12-19-11	12-20-11	
Hexachlorobenzene	ND	0.033	EPA 8270	12-19-11	12-20-11	
Pentachlorophenol	ND	0.17	EPA 8270	12-19-11	12-20-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Carbazole	ND	0.033	EPA 8270	12-19-11	12-20-11	
Di-n-butylphthalate	ND	0.33	EPA 8270	12-19-11	12-20-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzidine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Butylbenzylphthalate	ND	0.33	EPA 8270	12-19-11	12-20-11	
bis-2-Ethylhexyladipate	ND	0.17	EPA 8270	12-19-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.33	EPA 8270	12-19-11	12-20-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
bis(2-Ethylhexyl)phthalate	ND	0.17	EPA 8270	12-19-11	12-20-11	
Di-n-octylphthalate	ND	0.033	EPA 8270	12-19-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Indeno[1,2,3-cd]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	85	30 - 97				
Phenol-d6	80	40 - 104				
Nitrobenzene-d5	74	35 - 102				
2-Fluorobiphenyl	92	44 - 97				
2,4,6-Tribromophenol	103	41 - 110				
Terphenyl-d14	79	53 - 107				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
	SB	SBD	SB	SBD	SB	SBD				
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1219S2									
Phenol	<b>0.925</b>	<b>0.950</b>	1.33	1.33	70	71	31 - 111	3	34	
2-Chlorophenol	<b>1.04</b>	<b>1.06</b>	1.33	1.33	78	80	29 - 112	2	37	
1,4-Dichlorobenzene	<b>0.522</b>	<b>0.528</b>	0.667	0.667	78	79	24 - 100	1	37	
n-Nitroso-di-n-propylamine	<b>0.396</b>	<b>0.405</b>	0.667	0.667	59	61	35 - 104	2	32	
1,2,4-Trichlorobenzene	<b>0.685</b>	<b>0.713</b>	0.667	0.667	103	107	29 - 110	4	35	
4-Chloro-3-methylphenol	<b>1.13</b>	<b>1.19</b>	1.33	1.33	85	89	53 - 104	5	25	
Acenaphthene	<b>0.544</b>	<b>0.563</b>	0.667	0.667	82	84	50 - 95	3	23	
4-Nitrophenol	<b>1.22</b>	<b>1.30</b>	1.33	1.33	92	98	42 - 126	6	30	
2,4-Dinitrotoluene	<b>0.722</b>	<b>0.749</b>	0.667	0.667	108	112	53 - 115	4	31	
Pentachlorophenol	<b>0.939</b>	<b>1.03</b>	1.33	1.33	71	77	50 - 116	9	30	
Pyrene	<b>0.446</b>	<b>0.468</b>	0.667	0.667	67	70	57 - 120	5	27	
<i>Surrogate:</i>										
2-Fluorophenol					83	82	30 - 97			
Phenol-d6					76	78	40 - 104			
Nitrobenzene-d5					72	74	35 - 102			
2-Fluorobiphenyl					91	91	44 - 97			
2,4,6-Tribromophenol					105	111	41 - 110			
Terphenyl-d14					78	82	53 - 107			



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
Naphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>75</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>85</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>77</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD	RPD	Flags
					Result	Recovery	Limits			
<b>MATRIX SPIKES</b>										
Laboratory ID:	12-110-03									
	MS	MSD	MS	MSD		MS	MSD			
Naphthalene	<b>0.0546</b>	<b>0.0593</b>	0.0833	0.0833	ND	66	71	39 - 110	8	21
Acenaphthylene	<b>0.0530</b>	<b>0.0593</b>	0.0833	0.0833	ND	64	71	47 - 124	11	21
Acenaphthene	<b>0.0489</b>	<b>0.0560</b>	0.0833	0.0833	ND	59	67	50 - 120	14	20
Fluorene	<b>0.0487</b>	<b>0.0568</b>	0.0833	0.0833	ND	58	68	52 - 126	15	21
Phenanthrene	<b>0.0478</b>	<b>0.0551</b>	0.0833	0.0833	ND	57	66	41 - 130	14	22
Anthracene	<b>0.0473</b>	<b>0.0554</b>	0.0833	0.0833	ND	57	67	48 - 124	16	23
Fluoranthene	<b>0.0495</b>	<b>0.0581</b>	0.0833	0.0833	ND	59	70	40 - 137	16	23
Pyrene	<b>0.0487</b>	<b>0.0579</b>	0.0833	0.0833	ND	58	70	36 - 139	17	23
Benzo[a]anthracene	<b>0.0496</b>	<b>0.0593</b>	0.0833	0.0833	ND	60	71	43 - 127	18	21
Chrysene	<b>0.0476</b>	<b>0.0570</b>	0.0833	0.0833	ND	57	68	41 - 133	18	19
Benzo[b]fluoranthene	<b>0.0452</b>	<b>0.0555</b>	0.0833	0.0833	ND	54	67	40 - 132	20	25
Benzo(j,k)fluoranthene	<b>0.0436</b>	<b>0.0526</b>	0.0833	0.0833	ND	52	63	35 - 132	19	25
Benzo[a]pyrene	<b>0.0471</b>	<b>0.0563</b>	0.0833	0.0833	ND	57	68	37 - 131	18	26
Indeno(1,2,3-c,d)pyrene	<b>0.0444</b>	<b>0.0541</b>	0.0833	0.0833	ND	53	65	39 - 134	20	23
Dibenz[a,h]anthracene	<b>0.0457</b>	<b>0.0553</b>	0.0833	0.0833	ND	55	66	40 - 137	19	21
Benzo[g,h,i]perylene	<b>0.0456</b>	<b>0.0550</b>	0.0833	0.0833	ND	55	66	35 - 135	19	22
<i>Surrogate:</i>										
<i>2-Fluorobiphenyl</i>						<i>54</i>	<i>60</i>	<i>43 - 109</i>		
<i>Pyrene-d10</i>						<i>56</i>	<i>67</i>	<i>38 - 128</i>		
<i>Terphenyl-d14</i>						<i>53</i>	<i>63</i>	<i>33 - 119</i>		

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1216W1					
Aroclor 1016	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1221	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1232	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1242	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1248	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1254	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1260	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
<i>DCB</i>	78		36-127			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>SPIKE BLANKS</b>											
Laboratory ID:	SB1216W1										
	SB	SBD	SB	SBD		SB	SBD				
Aroclor 1260	<b>0.387</b>	<b>0.430</b>	0.500	0.500	N/A	<b>77</b>	<b>86</b>	57-122	11	11	
<i>Surrogate:</i>											
<i>DCB</i>						80	85	36-127			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1215S1					
Aroclor 1016	ND	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	ND	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	ND	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	ND	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	ND	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	ND	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	ND	0.050	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	91		42-123			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-099-07										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	0.534	0.460	0.500	0.500	ND	107	92	44-125	15	15	
<i>Surrogate:</i>											
DCB						96	83	42-123			

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-099  
Project: 0180-292-00

**TOTAL METALS  
EPA 200.8/7470A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-16-11  
Date Analyzed: 12-16-11  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB1216WM1&MB1216W1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	11
Lead	200.8	<b>ND</b>	1.1
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-16-11  
 Date Analyzed: 12-16-11  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: 12-091-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.3	
Barium	ND	ND	NA	28	
Cadmium	ND	ND	NA	4.4	
Chromium	ND	ND	NA	11	
Lead	ND	ND	NA	1.1	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.6	
Silver	ND	ND	NA	11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-16-11  
 Date Analyzed: 12-16-11  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: 12-091-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	111	<b>108</b>	97	<b>109</b>	98	1	
Barium	111	<b>113</b>	102	<b>115</b>	104	1	
Cadmium	111	<b>105</b>	95	<b>106</b>	95	1	
Chromium	111	<b>113</b>	102	<b>113</b>	102	0	
Lead	111	<b>111</b>	100	<b>111</b>	100	0	
Mercury	12.5	<b>12.0</b>	96	<b>12.2</b>	97	2	
Selenium	111	<b>101</b>	91	<b>105</b>	94	3	
Silver	111	<b>100</b>	90	<b>103</b>	92	3	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**DISSOLVED METALS  
 EPA 200.8/7470A  
 METHOD BLANK QUALITY CONTROL**

Date Filtered: 12-14-11  
 Date Analyzed: 12-15&20-11  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: MB1214F1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.0
Barium	200.8	<b>ND</b>	25
Cadmium	200.8	<b>ND</b>	4.0
Chromium	200.8	<b>ND</b>	10
Lead	200.8	<b>ND</b>	1.0
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.0
Silver	200.8	<b>ND</b>	10



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
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**DISSOLVED METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Filtered: 12-14-11  
 Date Analyzed: 12-15&20-11

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 12-091-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.0	
Barium	ND	ND	NA	25	
Cadmium	ND	ND	NA	4.0	
Chromium	ND	ND	NA	10	
Lead	ND	ND	NA	1.0	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.0	
Silver	ND	ND	NA	10	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
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**DISSOLVED METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Filtered: 12-14-11  
 Date Analyzed: 12-15&20-11

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 12-091-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	200	<b>209</b>	105	<b>205</b>	102	2	
Barium	200	<b>205</b>	102	<b>207</b>	104	1	
Cadmium	200	<b>207</b>	104	<b>204</b>	102	2	
Chromium	200	<b>188</b>	94	<b>188</b>	94	0	
Lead	200	<b>201</b>	101	<b>195</b>	98	3	
Mercury	12.5	<b>12.2</b>	98	<b>12.1</b>	97	1	
Selenium	200	<b>209</b>	104	<b>210</b>	105	1	
Silver	200	<b>180</b>	90	<b>181</b>	91	1	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-099  
Project: 0180-292-00

**TOTAL METALS  
EPA 6010B/7471A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-19-11  
Date Analyzed: 12-19-11  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB1219SM1&MB1219S1

Analyte	Method	Result	PQL
Arsenic	6010B	ND	10
Barium	6010B	ND	2.5
Cadmium	6010B	ND	0.50
Chromium	6010B	ND	0.50
Lead	6010B	ND	5.0
Mercury	7471A	ND	0.25
Selenium	6010B	ND	10
Silver	6010B	ND	0.50

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-19-11

Date Analyzed: 12-19-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 12-098-17

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	42.3	47.4	11	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	26.3	27.9	6	0.50	
Lead	21.6	22.6	5	5.0	
Mercury	ND	ND	NA	0.25	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-099  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-19-11

Date Analyzed: 12-19-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 12-098-17

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>88.7</b>	89	<b>89.0</b>	89	0	
Barium	100	<b>142</b>	100	<b>140</b>	97	2	
Cadmium	50.0	<b>47.0</b>	94	<b>47.2</b>	94	0	
Chromium	100	<b>119</b>	92	<b>121</b>	94	2	
Lead	250	<b>253</b>	93	<b>241</b>	88	5	
Mercury	0.500	<b>0.512</b>	102	<b>0.490</b>	98	4	
Selenium	100	<b>91.6</b>	92	<b>92.3</b>	92	1	
Silver	25.0	<b>21.0</b>	84	<b>21.1</b>	84	0	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
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**% MOISTURE**

Date Analyzed: 12-15-11

Client ID	Lab ID	% Moisture
SED3	12-099-03	35
SED4	12-099-04	37
SED5	12-099-05	25
SED6	12-099-06	33
SED7	12-099-07	12
SED8	12-099-08	24



### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference







14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

December 23, 2011

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1112-100

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on December 14, 2011.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-100  
Project: 0180-292-00

### Case Narrative

Samples were collected on December 13, 2011 and received by the laboratory on December 14, 2011. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

#### NWTPH Gx and Volatiles EPA 8260B Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-100  
Project: 0180-292-00

**ANALYTICAL REPORT FOR SAMPLES**

<b>Client ID</b>	<b>Laboratory ID</b>	<b>Matrix</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Notes</b>
DP1-W	12-100-01	Water	12-13-11	12-14-11	
SW1	12-100-02	Water	12-13-11	12-14-11	
SED1	12-100-03	Soil	12-13-11	12-14-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
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 Project: 0180-292-00

**NWTPH-Gx**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP1-W</b>					
Laboratory ID:	12-100-01					
Gasoline	<b>ND</b>	100	NWTPH-Gx	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	91	73-121				
<b>Client ID:</b>	<b>SW1</b>					
Laboratory ID:	12-100-02					
Gasoline	<b>ND</b>	100	NWTPH-Gx	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	73-121				

Date of Report: December 23, 2011  
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 Project: 0180-292-00

**NWTPH-Gx**

Matrix: Soil  
 Units: mg/kg (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SED1</b>					
Laboratory ID:	12-100-03					
Gasoline	<b>ND</b>	7.5	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	<i>102</i>	<i>68-124</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP1-W</b>					
Laboratory ID:	12-100-01					
Diesel Range Organics	<b>ND</b>	0.27	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	0.43	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>106</i>	<i>50-150</i>				
<b>Client ID:</b>	<b>SW1</b>					
Laboratory ID:	12-100-02					
Diesel Range Organics	<b>ND</b>	0.29	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	0.47	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>84</i>	<i>50-150</i>				

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**NWTPH-Dx**  
**(with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED1</b>					
Laboratory ID:	12-100-03					
Diesel Range Organics	<b>ND</b>	35	NWTPH-Dx	12-19-11	12-19-11	
Lube Oil Range Organics	<b>ND</b>	70	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>106</i>	<i>50-150</i>				

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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP1-W</b>					
Laboratory ID:	12-100-01					
Dichlorodifluoromethane	0.52	0.20	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Acetone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	2.0	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Toluene	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	



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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP1-W</b>					
Laboratory ID:	12-100-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.40	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	1.0	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>80</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>83</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>83</i>	<i>65-120</i>				

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Matrix: Water  
 Units: ug/L

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SW1</b>					
Laboratory ID:	12-100-02					
Dichlorodifluoromethane	0.25	0.20	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	0.62	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Acetone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	2.0	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Toluene	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW1</b>					
Laboratory ID:	12-100-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.40	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	1.0	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>84</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>83</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>85</i>	<i>65-120</i>				

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED1</b>					
Laboratory ID:	12-100-03					
Dichlorodifluoromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0069	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0069	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0069	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0069	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0014	EPA 8260	12-15-11	12-15-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED1</b>					
Laboratory ID:	12-100-03					
1,1,2-Trichloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0028	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0014	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0069	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0069	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0014	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>91</i>	<i>63-127</i>				
<i>Toluene-d8</i>	<i>95</i>	<i>65-129</i>				
<i>4-Bromofluorobenzene</i>	<i>88</i>	<i>55-121</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP1-W</b>					
<b>Laboratory ID:</b>	<b>12-100-01</b>					
n-Nitrosodimethylamine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Pyridine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Phenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
Aniline	ND	4.8	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Chlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,3-Dichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,4-Dichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Benzyl alcohol	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2-Dichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.95	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.95	EPA 8270	12-16-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.95	EPA 8270	12-16-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Hexachloroethane	ND	0.95	EPA 8270	12-16-11	12-20-11	
Nitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Isophorone	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Nitrophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4-Dimethylphenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4-Dichlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Naphthalene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
4-Chloroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
Hexachlorobutadiene	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Methylnaphthalene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Hexachlorocyclopentadiene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,3-Dichloroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Chloronaphthalene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2-Nitroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,4-Dinitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Dimethylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,3-Dinitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,6-Dinitrotoluene	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2-Dinitrobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Acenaphthylene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
3-Nitroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	

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**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP1-W</b>					
Laboratory ID:	12-100-01					
2,4-Dinitrophenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
Acenaphthene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
4-Nitrophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,4-Dinitrotoluene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Dibenzofuran	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.95	EPA 8270	12-16-11	12-20-11	
Diethylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Nitroaniline	ND	0.95	EPA 8270	12-16-11	12-20-11	
Fluorene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
4,6-Dinitro-2-methylphenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.95	EPA 8270	12-16-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.95	EPA 8270	12-16-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.95	EPA 8270	12-16-11	12-20-11	
Hexachlorobenzene	ND	0.95	EPA 8270	12-16-11	12-20-11	
Pentachlorophenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
Phenanthrene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Carbazole	ND	0.95	EPA 8270	12-16-11	12-20-11	
Di-n-butylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
Fluoranthene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Benzidine	ND	4.8	EPA 8270	12-16-11	12-20-11	
Pyrene	ND	0.095	EPA 8270/SIM	12-16-11	12-16-11	
Butylbenzylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
bis-2-Ethylhexyladipate	ND	4.8	EPA 8270	12-16-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.95	EPA 8270	12-16-11	12-20-11	
Benzo[a]anthracene	<b>0.0097</b>	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
bis(2-Ethylhexyl)phthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
Di-n-octylphthalate	ND	0.95	EPA 8270	12-16-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Indeno[1,2,3-cd]pyrene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	ND	0.0095	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	43	18 - 86				
Phenol-d6	30	10 - 88				
Nitrobenzene-d5	64	37 - 112				
2-Fluorobiphenyl	82	42 - 108				
2,4,6-Tribromophenol	109	39 - 118				
Terphenyl-d14	67	49 - 122				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW1</b>					
Laboratory ID:	12-100-02					
Naphthalene	<b>0.12</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
2-Methylnaphthalene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Acenaphthylene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Acenaphthene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Fluorene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Phenanthrene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Fluoranthene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Pyrene	<b>ND</b>	0.098	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]anthracene	<b>0.010</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	<b>ND</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[b]fluoranthene	<b>ND</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	<b>ND</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	<b>ND</b>	0.0098	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>82</i>	<i>38 - 105</i>				
<i>Pyrene-d10</i>	<i>87</i>	<i>37 - 121</i>				
<i>Terphenyl-d14</i>	<i>100</i>	<i>32 - 112</i>				



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SED1</b>					
Laboratory ID:	12-100-03					
Naphthalene	<b>0.048</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
2-Methylnaphthalene	<b>0.022</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
1-Methylnaphthalene	<b>0.014</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthylene	<b>ND</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Acenaphthene	<b>0.021</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Fluorene	<b>0.014</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Phenanthrene	<b>0.059</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Anthracene	<b>ND</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Fluoranthene	<b>0.030</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Pyrene	<b>0.040</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]anthracene	<b>0.011</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Chrysene	<b>0.016</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[b]fluoranthene	<b>0.013</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[a]pyrene	<b>0.012</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Dibenz[a,h]anthracene	<b>ND</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
Benzo[g,h,i]perylene	<b>0.0096</b>	0.0094	EPA 8270/SIM	12-19-11	12-20-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>59</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>65</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>63</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
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 Laboratory Reference: 1112-100  
 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP1-W</b>					
Laboratory ID:	12-100-01					
Aroclor 1016	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1221	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1232	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1242	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1248	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1254	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1260	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>71</i>	<i>36-127</i>				
<b>Client ID:</b>	<b>SW1</b>					
Laboratory ID:	12-100-02					
Aroclor 1016	<b>ND</b>	0.049	EPA 8082	12-16-11	12-17-11	
Aroclor 1221	<b>ND</b>	0.049	EPA 8082	12-16-11	12-17-11	
Aroclor 1232	<b>ND</b>	0.049	EPA 8082	12-16-11	12-17-11	
Aroclor 1242	<b>0.074</b>	0.049	EPA 8082	12-16-11	12-17-11	
Aroclor 1248	<b>ND</b>	0.049	EPA 8082	12-16-11	12-17-11	
Aroclor 1254	<b>ND</b>	0.049	EPA 8082	12-16-11	12-17-11	
Aroclor 1260	<b>ND</b>	0.049	EPA 8082	12-16-11	12-17-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>81</i>	<i>36-127</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SED1</b>					
Laboratory ID:	12-100-03					
Aroclor 1016	<b>ND</b>	0.070	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	<b>ND</b>	0.070	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	<b>ND</b>	0.070	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	<b>ND</b>	0.070	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	<b>ND</b>	0.070	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	<b>0.13</b>	0.070	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	<b>ND</b>	0.070	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>85</i>	<i>42-123</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	12-100-01					
<b>Client ID:</b>	<b>DP1-W</b>					
Arsenic	<b>18</b>	3.3	200.8	12-16-11	12-16-11	
Barium	<b>30</b>	28	200.8	12-16-11	12-16-11	
Cadmium	<b>ND</b>	4.4	200.8	12-16-11	12-16-11	
Chromium	<b>26</b>	11	200.8	12-16-11	12-16-11	
Lead	<b>ND</b>	1.1	200.8	12-16-11	12-16-11	
Mercury	<b>ND</b>	0.50	7470A	12-16-11	12-16-11	
Selenium	<b>ND</b>	5.6	200.8	12-16-11	12-16-11	
Silver	<b>ND</b>	11	200.8	12-16-11	12-16-11	

Lab ID:	12-100-02					
<b>Client ID:</b>	<b>SW1</b>					
Arsenic	<b>ND</b>	3.3	200.8	12-16-11	12-16-11	
Barium	<b>41</b>	28	200.8	12-16-11	12-16-11	
Cadmium	<b>ND</b>	4.4	200.8	12-16-11	12-16-11	
Chromium	<b>ND</b>	11	200.8	12-16-11	12-16-11	
Lead	<b>7.4</b>	1.1	200.8	12-16-11	12-16-11	
Mercury	<b>ND</b>	0.50	7470A	12-16-11	12-16-11	
Selenium	<b>ND</b>	5.6	200.8	12-16-11	12-16-11	
Silver	<b>ND</b>	11	200.8	12-16-11	12-16-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**DISSOLVED METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	12-100-01					
<b>Client ID:</b>	<b>DP1-W</b>					
Arsenic	<b>21</b>	3.0	200.8		12-15-11	
Barium	<b>ND</b>	25	200.8		12-15-11	
Cadmium	<b>ND</b>	4.0	200.8		12-15-11	
Chromium	<b>22</b>	10	200.8		12-15-11	
Lead	<b>ND</b>	1.0	200.8		12-15-11	
Mercury	<b>ND</b>	0.50	7470A		12-20-11	
Selenium	<b>ND</b>	5.0	200.8		12-15-11	
Silver	<b>ND</b>	10	200.8		12-15-11	

Lab ID:	12-100-02					
<b>Client ID:</b>	<b>SW1</b>					
Arsenic	<b>ND</b>	3.0	200.8	12-14-11	12-15-11	
Barium	<b>31</b>	25	200.8	12-14-11	12-15-11	
Cadmium	<b>ND</b>	4.0	200.8	12-14-11	12-15-11	
Chromium	<b>ND</b>	10	200.8	12-14-11	12-15-11	
Lead	<b>1.4</b>	1.0	200.8	12-14-11	12-15-11	
Mercury	<b>ND</b>	0.50	7470A	12-14-11	12-20-11	
Selenium	<b>ND</b>	5.0	200.8	12-14-11	12-15-11	
Silver	<b>ND</b>	10	200.8	12-14-11	12-15-11	

Date of Report: December 23, 2011  
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**TOTAL METALS  
 EPA 6010B/7471A**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	12-100-03					
<b>Client ID:</b>	<b>SED1</b>					
Arsenic	<b>ND</b>	14	6010B	12-19-11	12-19-11	
Barium	<b>85</b>	3.5	6010B	12-19-11	12-19-11	
Cadmium	<b>ND</b>	0.70	6010B	12-19-11	12-19-11	
Chromium	<b>36</b>	0.70	6010B	12-19-11	12-19-11	
Lead	<b>17</b>	7.0	6010B	12-19-11	12-19-11	
Mercury	<b>ND</b>	0.35	7471A	12-19-11	12-19-11	
Selenium	<b>ND</b>	14	6010B	12-19-11	12-19-11	
Silver	<b>ND</b>	0.70	6010B	12-19-11	12-19-11	

Date of Report: December 23, 2011  
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 Laboratory Reference: 1112-100  
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**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1215W2					
Gasoline	<b>ND</b>	100	NWTPH-Gx	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	73-121				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-100-01							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				91	90	73-121		

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**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1216S2					
Gasoline	<b>ND</b>	5.0	NWTPH-Gx	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	87	68-124				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-100-03							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				102	106	68-124		



Date of Report: December 23, 2011  
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 Project: 0180-292-00

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1221W1					
Diesel Range Organics	<b>ND</b>	0.25	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	0.40	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>101</i>	<i>50-150</i>				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-105-02					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			100 105	50-150		

Date of Report: December 23, 2011  
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 Project: 0180-292-00

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1219S1					
Mineral Oil	<b>ND</b>	25	NWTPH-Dx	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	120	50-150				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-110-05					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>			NA	NA
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>			NA	NA
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			104	102	50-150	

Date of Report: December 23, 2011  
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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Water

Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1215W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Acetone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	2.0	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Toluene	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:		MB1215W1				
1,1,2-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.40	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	1.0	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>85</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>90</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>88</i>	<i>65-120</i>				

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**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1215W1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>11.0</b>	<b>10.4</b>	10.0	10.0	110	104	70-130	6	11	
Benzene	<b>9.09</b>	<b>8.95</b>	10.0	10.0	91	90	75-123	2	8	
Trichloroethene	<b>10.1</b>	<b>10.3</b>	10.0	10.0	101	103	80-113	2	9	
Toluene	<b>9.74</b>	<b>9.71</b>	10.0	10.0	97	97	80-113	0	8	
Chlorobenzene	<b>10.8</b>	<b>10.5</b>	10.0	10.0	108	105	80-111	3	8	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					<i>77</i>	<i>77</i>	<i>68-120</i>			
<i>Toluene-d8</i>					<i>82</i>	<i>85</i>	<i>73-120</i>			
<i>4-Bromofluorobenzene</i>					<i>79</i>	<i>85</i>	<i>65-120</i>			

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Soil  
 Units: mg/kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1215S1					
Dichlorodifluoromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Acetone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	0.0050	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	0.0050	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	0.0050	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Toluene	ND	0.0050	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.0010	EPA 8260	12-15-11	12-15-11	

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**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1215S1					
1,1,2-Trichloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.0020	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.0010	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	0.0050	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.0050	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.0010	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	93	63-127				
<i>Toluene-d8</i>	98	65-129				
<i>4-Bromofluorobenzene</i>	96	55-121				

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**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Soil  
 Units: mg/kg

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					Recovery	Limits	RPD	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1215S1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	<b>0.0573</b>	<b>0.0584</b>	0.0500	0.0500	115	117	70-130	2	19	
Benzene	<b>0.0499</b>	<b>0.0504</b>	0.0500	0.0500	100	101	70-125	1	15	
Trichloroethene	<b>0.0474</b>	<b>0.0476</b>	0.0500	0.0500	95	95	70-122	0	14	
Toluene	<b>0.0493</b>	<b>0.0495</b>	0.0500	0.0500	99	99	73-120	0	16	
Chlorobenzene	<b>0.0451</b>	<b>0.0451</b>	0.0500	0.0500	90	90	74-109	0	12	
<i>Surrogate:</i>										
<i>Dibromofluoromethane</i>					<i>85</i>	<i>84</i>	<i>63-127</i>			
<i>Toluene-d8</i>					<i>94</i>	<i>90</i>	<i>65-129</i>			
<i>4-Bromofluorobenzene</i>					<i>91</i>	<i>84</i>	<i>55-121</i>			



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL**

page 1 of 2

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216W2					
n-Nitrosodimethylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Pyridine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Phenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
Aniline	ND	5.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethyl)ether	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Chlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,3-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,4-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzyl alcohol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	1.0	EPA 8270	12-16-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	1.0	EPA 8270	12-16-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachloroethane	ND	1.0	EPA 8270	12-16-11	12-20-11	
Nitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Isophorone	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Nitrophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dimethylphenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2,4-Trichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Naphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4-Chloroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachlorobutadiene	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Chloro-3-methylphenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Hexachlorocyclopentadiene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4,6-Trichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3-Dichloroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4,5-Trichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Chloronaphthalene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,4-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Dimethylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,3-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,6-Dinitrotoluene	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Acenaphthylene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
3-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	

Date of Report: December 23, 2011  
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 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216W2					
2,4-Dinitrophenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
Acenaphthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4-Nitrophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dinitrotoluene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Dibenzofuran	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
Diethylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Chlorophenyl-phenylether	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
Fluorene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4,6-Dinitro-2-methylphenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
n-Nitrosodiphenylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Diphenylhydrazine	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Bromophenyl-phenylether	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Pentachlorophenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
Phenanthrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Carbazole	ND	1.0	EPA 8270	12-16-11	12-20-11	
Di-n-butylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Fluoranthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Benzidine	ND	5.0	EPA 8270	12-16-11	12-20-11	
Pyrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Butylbenzylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis-2-Ethylhexyladipate	ND	5.0	EPA 8270	12-16-11	12-20-11	
3,3'-Dichlorobenzidine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzo[a]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
bis(2-Ethylhexyl)phthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Di-n-octylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzo[b]fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Indeno[1,2,3-cd]pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	57	18 - 86				
Phenol-d6	41	10 - 88				
Nitrobenzene-d5	76	37 - 112				
2-Fluorobiphenyl	92	42 - 108				
2,4,6-Tribromophenol	108	39 - 118				
Terphenyl-d14	83	49 - 122				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

<b>Analyte</b>	<b>Result</b>		<b>Spike Level</b>		<b>Percent Recovery</b>		<b>Recovery Limits</b>	<b>RPD</b>	<b>RPD Limit</b>	<b>Flags</b>
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1216W2									
	SB	SBD	SB	SBD	SB	SBD				
Phenol	15.5	13.3	40.0	40.0	39	33	26 - 60	15	29	
2-Chlorophenol	29.6	24.4	40.0	40.0	74	61	46 - 104	19	34	
1,4-Dichlorobenzene	13.0	10.6	20.0	20.0	65	53	46 - 92	20	29	
n-Nitroso-di-n-propylamine	12.1	10.2	20.0	20.0	61	51	30 - 102	17	25	
1,2,4-Trichlorobenzene	18.3	15.4	20.0	20.0	92	77	45 - 92	17	25	
4-Chloro-3-methylphenol	34.4	32.1	40.0	40.0	86	80	53 - 104	7	18	
Acenaphthene	16.0	14.6	20.0	20.0	80	73	57 - 95	9	15	
4-Nitrophenol	22.4	20.8	40.0	40.0	56	52	21 - 75	7	33	
2,4-Dinitrotoluene	21.4	20.3	20.0	20.0	107	102	60 - 108	5	20	
Pentachlorophenol	26.3	25.2	40.0	40.0	66	63	48 - 119	4	31	
Pyrene	13.8	13.3	20.0	20.0	69	67	62 - 111	4	19	
<i>Surrogate:</i>										
2-Fluorophenol					58	46	18 - 86			
Phenol-d6					42	34	10 - 88			
Nitrobenzene-d5					73	61	37 - 112			
2-Fluorobiphenyl					90	79	42 - 108			
2,4,6-Tribromophenol					103	100	39 - 118			
Terphenyl-d14					80	78	49 - 122			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216W1					
Naphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
2-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Acenaphthylene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Acenaphthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Fluorene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Phenanthrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Fluoranthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Pyrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[b]fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>69</i>	<i>38 - 105</i>				
<i>Pyrene-d10</i>	<i>74</i>	<i>37 - 121</i>				
<i>Terphenyl-d14</i>	<i>87</i>	<i>32 - 112</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1216W1									
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	0.412	0.448	0.500	0.500	82	90	38 - 110	8	35	
Acenaphthylene	0.510	0.541	0.500	0.500	102	108	47 - 120	6	30	
Acenaphthene	0.470	0.507	0.500	0.500	94	101	46 - 113	8	26	
Fluorene	0.465	0.492	0.500	0.500	93	98	60 - 104	6	25	
Phenanthrene	0.435	0.473	0.500	0.500	87	95	61 - 99	8	19	
Anthracene	0.458	0.499	0.500	0.500	92	100	55 - 122	9	19	
Fluoranthene	0.448	0.488	0.500	0.500	90	98	58 - 129	9	18	
Pyrene	0.438	0.532	0.500	0.500	88	106	57 - 126	19	22	
Benzo[a]anthracene	0.441	0.478	0.500	0.500	88	96	51 - 124	8	18	
Chrysene	0.434	0.478	0.500	0.500	87	96	53 - 123	10	20	
Benzo[b]fluoranthene	0.449	0.503	0.500	0.500	90	101	53 - 126	11	18	
Benzo(j,k)fluoranthene	0.443	0.474	0.500	0.500	89	95	51 - 126	7	23	
Benzo[a]pyrene	0.472	0.514	0.500	0.500	94	103	52 - 127	9	21	
Indeno(1,2,3-c,d)pyrene	0.444	0.489	0.500	0.500	89	98	49 - 123	10	26	
Dibenz[a,h]anthracene	0.435	0.479	0.500	0.500	87	96	39 - 125	10	31	
Benzo[g,h,i]perylene	0.421	0.464	0.500	0.500	84	93	40 - 125	10	30	
<i>Surrogate:</i>										
2-Fluorobiphenyl					81	85	38 - 105			
Pyrene-d10					83	90	37 - 121			
Terphenyl-d14					104	108	32 - 112			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1219S1					
Naphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
2-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
1-Methylnaphthalene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Acenaphthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluorene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Phenanthrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Chrysene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[b]fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo(j,k)fluoranthene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[a]pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Indeno(1,2,3-c,d)pyrene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Dibenz[a,h]anthracene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
Benzo[g,h,i]perylene	ND	0.0067	EPA 8270/SIM	12-19-11	12-19-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>75</i>	<i>43 - 109</i>				
<i>Pyrene-d10</i>	<i>85</i>	<i>38 - 128</i>				
<i>Terphenyl-d14</i>	<i>77</i>	<i>33 - 119</i>				

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 MS/MSD QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Soil  
 Units: mg/Kg

Analyte	Result		Spike Level		Source	Percent	Recovery	RPD		Flags
					Result	Recovery	Limits	RPD	Limit	
<b>MATRIX SPIKES</b>										
Laboratory ID:	12-110-03									
	MS	MSD	MS	MSD		MS	MSD			
Naphthalene	<b>0.0546</b>	<b>0.0593</b>	0.0833	0.0833	ND	66	71	39 - 110	8	21
Acenaphthylene	<b>0.0530</b>	<b>0.0593</b>	0.0833	0.0833	ND	64	71	47 - 124	11	21
Acenaphthene	<b>0.0489</b>	<b>0.0560</b>	0.0833	0.0833	ND	59	67	50 - 120	14	20
Fluorene	<b>0.0487</b>	<b>0.0568</b>	0.0833	0.0833	ND	58	68	52 - 126	15	21
Phenanthrene	<b>0.0478</b>	<b>0.0551</b>	0.0833	0.0833	ND	57	66	41 - 130	14	22
Anthracene	<b>0.0473</b>	<b>0.0554</b>	0.0833	0.0833	ND	57	67	48 - 124	16	23
Fluoranthene	<b>0.0495</b>	<b>0.0581</b>	0.0833	0.0833	ND	59	70	40 - 137	16	23
Pyrene	<b>0.0487</b>	<b>0.0579</b>	0.0833	0.0833	ND	58	70	36 - 139	17	23
Benzo[a]anthracene	<b>0.0496</b>	<b>0.0593</b>	0.0833	0.0833	ND	60	71	43 - 127	18	21
Chrysene	<b>0.0476</b>	<b>0.0570</b>	0.0833	0.0833	ND	57	68	41 - 133	18	19
Benzo[b]fluoranthene	<b>0.0452</b>	<b>0.0555</b>	0.0833	0.0833	ND	54	67	40 - 132	20	25
Benzo(j,k)fluoranthene	<b>0.0436</b>	<b>0.0526</b>	0.0833	0.0833	ND	52	63	35 - 132	19	25
Benzo[a]pyrene	<b>0.0471</b>	<b>0.0563</b>	0.0833	0.0833	ND	57	68	37 - 131	18	26
Indeno(1,2,3-c,d)pyrene	<b>0.0444</b>	<b>0.0541</b>	0.0833	0.0833	ND	53	65	39 - 134	20	23
Dibenz[a,h]anthracene	<b>0.0457</b>	<b>0.0553</b>	0.0833	0.0833	ND	55	66	40 - 137	19	21
Benzo[g,h,i]perylene	<b>0.0456</b>	<b>0.0550</b>	0.0833	0.0833	ND	55	66	35 - 135	19	22
<i>Surrogate:</i>										
<i>2-Fluorobiphenyl</i>						<i>54</i>	<i>60</i>	<i>43 - 109</i>		
<i>Pyrene-d10</i>						<i>56</i>	<i>67</i>	<i>38 - 128</i>		
<i>Terphenyl-d14</i>						<i>53</i>	<i>63</i>	<i>33 - 119</i>		

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1216W1					
Aroclor 1016	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1221	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1232	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1242	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1248	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1254	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1260	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	78	36-127				

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>SPIKE BLANKS</b>											
Laboratory ID:	SB1216W1										
	SB	SBD	SB	SBD		SB	SBD				
Aroclor 1260	<b>0.387</b>	<b>0.430</b>	0.500	0.500	N/A	<b>77</b>	<b>86</b>	57-122	11	11	
<i>Surrogate:</i>											
DCB						80	85	36-127			



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Soil  
 Units: mg/Kg (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1215S1					
Aroclor 1016	<b>ND</b>	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1221	<b>ND</b>	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1232	<b>ND</b>	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1242	<b>ND</b>	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1248	<b>ND</b>	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1254	<b>ND</b>	0.050	EPA 8082	12-15-11	12-16-11	
Aroclor 1260	<b>ND</b>	0.050	EPA 8082	12-15-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	91		42-123			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>MATRIX SPIKES</b>											
Laboratory ID:	12-099-07										
	MS	MSD	MS	MSD		MS	MSD				
Aroclor 1260	<b>0.534</b>	<b>0.460</b>	0.500	0.500	ND	<b>107</b>	<b>92</b>	44-125	15	15	
<i>Surrogate:</i>											
DCB						96	83	42-123			

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-100  
Project: 0180-292-00

**TOTAL METALS  
EPA 200.8/7470A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-16-11  
Date Analyzed: 12-16-11  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB1216WM1&MB1216W1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	11
Lead	200.8	<b>ND</b>	1.1
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-16-11  
 Date Analyzed: 12-16-11  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: 12-091-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.3	
Barium	ND	ND	NA	28	
Cadmium	ND	ND	NA	4.4	
Chromium	ND	ND	NA	11	
Lead	ND	ND	NA	1.1	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.6	
Silver	ND	ND	NA	11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-16-11  
 Date Analyzed: 12-16-11  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: 12-091-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	111	<b>108</b>	97	<b>109</b>	98	1	
Barium	111	<b>113</b>	102	<b>115</b>	104	1	
Cadmium	111	<b>105</b>	95	<b>106</b>	95	1	
Chromium	111	<b>113</b>	102	<b>113</b>	102	0	
Lead	111	<b>111</b>	100	<b>111</b>	100	0	
Mercury	12.5	<b>12.0</b>	96	<b>12.2</b>	97	2	
Selenium	111	<b>101</b>	91	<b>105</b>	94	3	
Silver	111	<b>100</b>	90	<b>103</b>	92	3	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-100  
Project: 0180-292-00

**DISSOLVED METALS  
EPA 200.8/7470A  
METHOD BLANK QUALITY CONTROL**

Date Filtered: 12-14-11  
Date Analyzed: 12-15&20-11  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB1214F1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.0
Barium	200.8	<b>ND</b>	25
Cadmium	200.8	<b>ND</b>	4.0
Chromium	200.8	<b>ND</b>	10
Lead	200.8	<b>ND</b>	1.0
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.0
Silver	200.8	<b>ND</b>	10

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**DISSOLVED METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Filtered: 12-14-11  
 Date Analyzed: 12-15&20-11

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 12-091-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.0	
Barium	ND	ND	NA	25	
Cadmium	ND	ND	NA	4.0	
Chromium	ND	ND	NA	10	
Lead	ND	ND	NA	1.0	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.0	
Silver	ND	ND	NA	10	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**DISSOLVED METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Filtered: 12-14-11  
 Date Analyzed: 12-15&20-11

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 12-091-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	200	<b>209</b>	105	<b>205</b>	102	2	
Barium	200	<b>205</b>	102	<b>207</b>	104	1	
Cadmium	200	<b>207</b>	104	<b>204</b>	102	2	
Chromium	200	<b>188</b>	94	<b>188</b>	94	0	
Lead	200	<b>201</b>	101	<b>195</b>	98	3	
Mercury	12.5	<b>12.2</b>	98	<b>12.1</b>	97	1	
Selenium	200	<b>209</b>	104	<b>210</b>	105	1	
Silver	200	<b>180</b>	90	<b>181</b>	91	1	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-100  
Project: 0180-292-00

**TOTAL METALS  
EPA 6010B/7471A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-19-11  
Date Analyzed: 12-19-11  
  
Matrix: Soil  
Units: mg/kg (ppm)  
  
Lab ID: MB1219SM1&MB1219S1

Analyte	Method	Result	PQL
Arsenic	6010B	ND	10
Barium	6010B	ND	2.5
Cadmium	6010B	ND	0.50
Chromium	6010B	ND	0.50
Lead	6010B	ND	5.0
Mercury	7471A	ND	0.25
Selenium	6010B	ND	10
Silver	6010B	ND	0.50



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-19-11  
 Date Analyzed: 12-19-11  
  
 Matrix: Soil  
 Units: mg/kg (ppm)  
  
 Lab ID: 12-098-17

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	10	
Barium	42.3	47.4	11	2.5	
Cadmium	ND	ND	NA	0.50	
Chromium	26.3	27.9	6	0.50	
Lead	21.6	22.6	5	5.0	
Mercury	ND	ND	NA	0.25	
Selenium	ND	ND	NA	10	
Silver	ND	ND	NA	0.50	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-100  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 6010B/7471A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-19-11

Date Analyzed: 12-19-11

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 12-098-17

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	100	<b>88.7</b>	89	<b>89.0</b>	89	0	
Barium	100	<b>142</b>	100	<b>140</b>	97	2	
Cadmium	50.0	<b>47.0</b>	94	<b>47.2</b>	94	0	
Chromium	100	<b>119</b>	92	<b>121</b>	94	2	
Lead	250	<b>253</b>	93	<b>241</b>	88	5	
Mercury	0.500	<b>0.512</b>	102	<b>0.490</b>	98	4	
Selenium	100	<b>91.6</b>	92	<b>92.3</b>	92	1	
Silver	25.0	<b>21.0</b>	84	<b>21.1</b>	84	0	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-100  
Project: 0180-292-00

**% MOISTURE**

Date Analyzed: 12-15-11

Client ID	Lab ID	% Moisture
SED1	12-100-03	29



#### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference





14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

December 23, 2011

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1112-101

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on December 14, 2011.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

A handwritten signature in black ink, appearing to read "DB", with a long horizontal stroke extending to the right.

David Baumeister  
Project Manager

Enclosures

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-101  
Project: 0180-292-00

### **Case Narrative**

Samples were collected on December 14, 2011 and received by the laboratory on December 14, 2011. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-101  
Project: 0180-292-00

**ANALYTICAL REPORT FOR SAMPLES**

<b>Client ID</b>	<b>Laboratory ID</b>	<b>Matrix</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Notes</b>
Well1-W	12-101-01	Water	12-14-11	12-14-11	



Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-101  
Project: 0180-292-00

**NWTPH-Gx**

Matrix: Water  
Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>Well1-W</b>					
Laboratory ID:	12-101-01					
Gasoline	<b>ND</b>	100	NWTPH-Gx	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	91	73-121				

Date of Report: December 23, 2011  
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**NWTPH-Dx**  
**(with acid/silica gel clean-up)**

Matrix: Water  
 Units: mg/L (ppm)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>Well1-W</b>					
Laboratory ID:	12-101-01					
Diesel Range Organics	<b>ND</b>	0.26	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	0.41	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>101</i>	<i>50-150</i>				

Date of Report: December 23, 2011  
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**VOLATILES by EPA 8260B**  
 Page 1 of 2

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Well1-W</b>					
Laboratory ID:	12-101-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Acetone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	2.0	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Toluene	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	

Date of Report: December 23, 2011  
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**VOLATILES by EPA 8260B**  
 Page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Well1-W</b>					
Laboratory ID:	12-101-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.40	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	1.0	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>83</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>82</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>84</i>	<i>65-120</i>				

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 Laboratory Reference: 1112-101  
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**SEMIVOLATILES by EPA 8270D/SIM**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Well1-W</b>					
<b>Laboratory ID:</b>	<b>12-101-01</b>					
n-Nitrosodimethylamine	ND	0.96	EPA 8270	12-16-11	12-20-11	
Pyridine	ND	0.96	EPA 8270	12-16-11	12-20-11	
Phenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
Aniline	ND	4.8	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethyl)ether	ND	0.96	EPA 8270	12-16-11	12-20-11	
2-Chlorophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,3-Dichlorobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,4-Dichlorobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
Benzyl alcohol	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,2-Dichlorobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	0.96	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	0.96	EPA 8270	12-16-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.96	EPA 8270	12-16-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	0.96	EPA 8270	12-16-11	12-20-11	
Hexachloroethane	ND	0.96	EPA 8270	12-16-11	12-20-11	
Nitrobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
Isophorone	ND	0.96	EPA 8270	12-16-11	12-20-11	
2-Nitrophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,4-Dimethylphenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,4-Dichlorophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,2,4-Trichlorobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
Naphthalene	0.19	0.096	EPA 8270/SIM	12-16-11	12-16-11	
4-Chloroaniline	ND	0.96	EPA 8270	12-16-11	12-20-11	
Hexachlorobutadiene	ND	0.96	EPA 8270	12-16-11	12-20-11	
4-Chloro-3-methylphenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
2-Methylnaphthalene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
Hexachlorocyclopentadiene	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,4,6-Trichlorophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,3-Dichloroaniline	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,4,5-Trichlorophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
2-Chloronaphthalene	ND	0.96	EPA 8270	12-16-11	12-20-11	
2-Nitroaniline	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,4-Dinitrobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
Dimethylphthalate	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,3-Dinitrobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,6-Dinitrotoluene	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,2-Dinitrobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
Acenaphthylene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
3-Nitroaniline	ND	0.96	EPA 8270	12-16-11	12-20-11	

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**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Well1-W</b>					
Laboratory ID:	12-101-01					
2,4-Dinitrophenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
Acenaphthene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
4-Nitrophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,4-Dinitrotoluene	ND	0.96	EPA 8270	12-16-11	12-20-11	
Dibenzofuran	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	0.96	EPA 8270	12-16-11	12-20-11	
Diethylphthalate	ND	0.96	EPA 8270	12-16-11	12-20-11	
4-Chlorophenyl-phenylether	ND	0.96	EPA 8270	12-16-11	12-20-11	
4-Nitroaniline	ND	0.96	EPA 8270	12-16-11	12-20-11	
Fluorene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
4,6-Dinitro-2-methylphenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
n-Nitrosodiphenylamine	ND	0.96	EPA 8270	12-16-11	12-20-11	
1,2-Diphenylhydrazine	ND	0.96	EPA 8270	12-16-11	12-20-11	
4-Bromophenyl-phenylether	ND	0.96	EPA 8270	12-16-11	12-20-11	
Hexachlorobenzene	ND	0.96	EPA 8270	12-16-11	12-20-11	
Pentachlorophenol	ND	4.8	EPA 8270	12-16-11	12-20-11	
Phenanthrene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
Carbazole	ND	0.96	EPA 8270	12-16-11	12-20-11	
Di-n-butylphthalate	ND	0.96	EPA 8270	12-16-11	12-20-11	
Fluoranthene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
Benzidine	ND	4.8	EPA 8270	12-16-11	12-20-11	
Pyrene	ND	0.096	EPA 8270/SIM	12-16-11	12-16-11	
Butylbenzylphthalate	ND	0.96	EPA 8270	12-16-11	12-20-11	
bis-2-Ethylhexyladipate	ND	4.8	EPA 8270	12-16-11	12-20-11	
3,3'-Dichlorobenzidine	ND	0.96	EPA 8270	12-16-11	12-20-11	
Benzo[a]anthracene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
bis(2-Ethylhexyl)phthalate	ND	0.96	EPA 8270	12-16-11	12-20-11	
Di-n-octylphthalate	ND	0.96	EPA 8270	12-16-11	12-20-11	
Benzo[b]fluoranthene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
Indeno[1,2,3-cd]pyrene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	ND	0.0096	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	50	18 - 86				
Phenol-d6	34	10 - 88				
Nitrobenzene-d5	67	37 - 112				
2-Fluorobiphenyl	86	42 - 108				
2,4,6-Tribromophenol	111	39 - 118				
Terphenyl-d14	79	49 - 122				

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 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>Well1-W</b>					
Laboratory ID:	12-101-01					
Aroclor 1016	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1221	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1232	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1242	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1248	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1254	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
Aroclor 1260	<b>ND</b>	0.048	EPA 8082	12-16-11	12-17-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>DCB</i>	<i>87</i>	<i>36-127</i>				

Date of Report: December 23, 2011  
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 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	12-101-01					
<b>Client ID:</b>	<b>Well1-W</b>					
Arsenic	ND	3.3	200.8	12-16-11	12-16-11	
Barium	ND	28	200.8	12-16-11	12-16-11	
Cadmium	ND	4.4	200.8	12-16-11	12-16-11	
Chromium	ND	11	200.8	12-16-11	12-16-11	
Lead	ND	1.1	200.8	12-16-11	12-16-11	
Mercury	ND	0.50	7470A	12-16-11	12-16-11	
Selenium	ND	5.6	200.8	12-16-11	12-16-11	
Silver	ND	11	200.8	12-16-11	12-16-11	



Date of Report: December 23, 2011  
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**DISSOLVED METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>EPA Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
Lab ID:	12-101-01					
<b>Client ID:</b>	<b>Well1-W</b>					
Arsenic	<b>ND</b>	3.0	200.8		12-15-11	
Barium	<b>ND</b>	25	200.8		12-15-11	
Cadmium	<b>ND</b>	4.0	200.8		12-15-11	
Chromium	<b>ND</b>	10	200.8		12-15-11	
Lead	<b>ND</b>	1.0	200.8		12-15-11	
Mercury	<b>ND</b>	0.50	7470A		12-20-11	
Selenium	<b>ND</b>	5.0	200.8		12-15-11	
Silver	<b>ND</b>	10	200.8		12-15-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1215W2					
Gasoline	<b>ND</b>	100	NWTPH-Gx	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	73-121				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	12-100-01							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				91	90	73-121		

Date of Report: December 23, 2011  
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 Project: 0180-292-00

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1221W1					
Diesel Range Organics	<b>ND</b>	0.25	NWTPH-Dx	12-21-11	12-21-11	
Lube Oil Range Organics	<b>ND</b>	0.40	NWTPH-Dx	12-21-11	12-21-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	<i>101</i>	<i>50-150</i>				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	12-101-01					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			101 104	50-150		

Date of Report: December 23, 2011  
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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1215W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloromethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Vinyl Chloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Trichlorofluoromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Acetone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Iodomethane	ND	1.0	EPA 8260	12-15-11	12-15-11	
Carbon Disulfide	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methylene Chloride	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Vinyl Acetate	ND	2.0	EPA 8260	12-15-11	12-15-11	
2,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Butanone	ND	5.0	EPA 8260	12-15-11	12-15-11	
Bromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chloroform	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Carbon Tetrachloride	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Benzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Trichloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Dibromomethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromodichloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	12-15-11	12-15-11	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Toluene	ND	1.0	EPA 8260	12-15-11	12-15-11	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	12-15-11	12-15-11	

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:		MB1215W1				
1,1,2-Trichloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Tetrachloroethene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Hexanone	ND	2.0	EPA 8260	12-15-11	12-15-11	
Dibromochloromethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromoethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Chlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
Ethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
m,p-Xylene	ND	0.40	EPA 8260	12-15-11	12-15-11	
o-Xylene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Styrene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromoform	ND	1.0	EPA 8260	12-15-11	12-15-11	
Isopropylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Bromobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Propylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
2-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
4-Chlorotoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
tert-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
sec-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
p-Isopropyltoluene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
n-Butylbenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Hexachlorobutadiene	ND	0.20	EPA 8260	12-15-11	12-15-11	
Naphthalene	ND	1.0	EPA 8260	12-15-11	12-15-11	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	12-15-11	12-15-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>85</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>90</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>88</i>	<i>65-120</i>				

Date of Report: December 23, 2011  
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 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**VOLATILES by EPA 8260B  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD		Flags
					SB	SBD	Limits	RPD	Limit	
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1215W1									
	SB	SBD	SB	SBD	SB	SBD				
1,1-Dichloroethene	11.0	10.4	10.0	10.0	110	104	70-130	6	11	
Benzene	9.09	8.95	10.0	10.0	91	90	75-123	2	8	
Trichloroethene	10.1	10.3	10.0	10.0	101	103	80-113	2	9	
Toluene	9.74	9.71	10.0	10.0	97	97	80-113	0	8	
Chlorobenzene	10.8	10.5	10.0	10.0	108	105	80-111	3	8	
<i>Surrogate:</i>										
					77	77	68-120			
					82	85	73-120			
					79	85	65-120			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 1 of 2

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216W2					
n-Nitrosodimethylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Pyridine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Phenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
Aniline	ND	5.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethyl)ether	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Chlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,3-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,4-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzyl alcohol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Dichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Methylphenol (o-Cresol)	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroisopropyl)ether	ND	1.0	EPA 8270	12-16-11	12-20-11	
(3+4)-Methylphenol (m,p-Cresol)	ND	1.0	EPA 8270	12-16-11	12-20-11	
n-Nitroso-di-n-propylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachloroethane	ND	1.0	EPA 8270	12-16-11	12-20-11	
Nitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Isophorone	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Nitrophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dimethylphenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis(2-Chloroethoxy)methane	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2,4-Trichlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Naphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4-Chloroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachlorobutadiene	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Chloro-3-methylphenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
1-Methylnaphthalene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Hexachlorocyclopentadiene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4,6-Trichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3-Dichloroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4,5-Trichlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Chloronaphthalene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,4-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Dimethylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,3-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,6-Dinitrotoluene	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Dinitrobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Acenaphthylene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
3-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB1216W2					
2,4-Dinitrophenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
Acenaphthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4-Nitrophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,4-Dinitrotoluene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Dibenzofuran	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3,5,6-Tetrachlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
2,3,4,6-Tetrachlorophenol	ND	1.0	EPA 8270	12-16-11	12-20-11	
Diethylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Chlorophenyl-phenylether	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Nitroaniline	ND	1.0	EPA 8270	12-16-11	12-20-11	
Fluorene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
4,6-Dinitro-2-methylphenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
n-Nitrosodiphenylamine	ND	1.0	EPA 8270	12-16-11	12-20-11	
1,2-Diphenylhydrazine	ND	1.0	EPA 8270	12-16-11	12-20-11	
4-Bromophenyl-phenylether	ND	1.0	EPA 8270	12-16-11	12-20-11	
Hexachlorobenzene	ND	1.0	EPA 8270	12-16-11	12-20-11	
Pentachlorophenol	ND	5.0	EPA 8270	12-16-11	12-20-11	
Phenanthrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Anthracene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Carbazole	ND	1.0	EPA 8270	12-16-11	12-20-11	
Di-n-butylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Fluoranthene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Benzidine	ND	5.0	EPA 8270	12-16-11	12-20-11	
Pyrene	ND	0.10	EPA 8270/SIM	12-16-11	12-16-11	
Butylbenzylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
bis-2-Ethylhexyladipate	ND	5.0	EPA 8270	12-16-11	12-20-11	
3,3'-Dichlorobenzidine	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzo[a]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Chrysene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
bis(2-Ethylhexyl)phthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Di-n-octylphthalate	ND	1.0	EPA 8270	12-16-11	12-20-11	
Benzo[b]fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[a]pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Indeno[1,2,3-cd]pyrene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270/SIM	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
2-Fluorophenol	57	18 - 86				
Phenol-d6	41	10 - 88				
Nitrobenzene-d5	76	37 - 112				
2-Fluorobiphenyl	92	42 - 108				
2,4,6-Tribromophenol	108	39 - 118				
Terphenyl-d14	83	49 - 122				



Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
	SB	SBD	SB	SBD	SB	SBD				
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB1216W2									
Phenol	15.5	13.3	40.0	40.0	39	33	26 - 60	15	29	
2-Chlorophenol	29.6	24.4	40.0	40.0	74	61	46 - 104	19	34	
1,4-Dichlorobenzene	13.0	10.6	20.0	20.0	65	53	46 - 92	20	29	
n-Nitroso-di-n-propylamine	12.1	10.2	20.0	20.0	61	51	30 - 102	17	25	
1,2,4-Trichlorobenzene	18.3	15.4	20.0	20.0	92	77	45 - 92	17	25	
4-Chloro-3-methylphenol	34.4	32.1	40.0	40.0	86	80	53 - 104	7	18	
Acenaphthene	16.0	14.6	20.0	20.0	80	73	57 - 95	9	15	
4-Nitrophenol	22.4	20.8	40.0	40.0	56	52	21 - 75	7	33	
2,4-Dinitrotoluene	21.4	20.3	20.0	20.0	107	102	60 - 108	5	20	
Pentachlorophenol	26.3	25.2	40.0	40.0	66	63	48 - 119	4	31	
Pyrene	13.8	13.3	20.0	20.0	69	67	62 - 111	4	19	
<i>Surrogate:</i>										
2-Fluorophenol					58	46	18 - 86			
Phenol-d6					42	34	10 - 88			
Nitrobenzene-d5					73	61	37 - 112			
2-Fluorobiphenyl					90	79	42 - 108			
2,4,6-Tribromophenol					103	100	39 - 118			
Terphenyl-d14					80	78	49 - 122			

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB1216W1					
Aroclor 1016	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1221	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1232	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1242	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1248	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1254	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
Aroclor 1260	<b>ND</b>	0.050	EPA 8082	12-16-11	12-16-11	
<i>Surrogate:</i>	<i>Percent Recovery</i>		<i>Control Limits</i>			
<i>DCB</i>	78		36-127			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>SPIKE BLANKS</b>											
Laboratory ID:	SB1216W1										
	SB	SBD	SB	SBD		SB	SBD				
Aroclor 1260	<b>0.387</b>	<b>0.430</b>	0.500	0.500	N/A	<b>77</b>	<b>86</b>	57-122	11	11	
<i>Surrogate:</i>											
<i>DCB</i>						80	85	36-127			

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-101  
Project: 0180-292-00

**TOTAL METALS  
EPA 200.8/7470A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 12-16-11  
Date Analyzed: 12-16-11  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB1216WM1&MB1216W1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	11
Lead	200.8	<b>ND</b>	1.1
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 12-16-11  
 Date Analyzed: 12-16-11  
  
 Matrix: Water  
 Units: ug/L (ppb)  
  
 Lab ID: 12-091-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.3	
Barium	ND	ND	NA	28	
Cadmium	ND	ND	NA	4.4	
Chromium	ND	ND	NA	11	
Lead	ND	ND	NA	1.1	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.6	
Silver	ND	ND	NA	11	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Extracted: 12-16-11  
 Date Analyzed: 12-16-11  
 Matrix: Water  
 Units: ug/L (ppb)  
 Lab ID: 12-091-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	111	<b>108</b>	97	<b>109</b>	98	1	
Barium	111	<b>113</b>	102	<b>115</b>	104	1	
Cadmium	111	<b>105</b>	95	<b>106</b>	95	1	
Chromium	111	<b>113</b>	102	<b>113</b>	102	0	
Lead	111	<b>111</b>	100	<b>111</b>	100	0	
Mercury	12.5	<b>12.0</b>	96	<b>12.2</b>	97	2	
Selenium	111	<b>101</b>	91	<b>105</b>	94	3	
Silver	111	<b>100</b>	90	<b>103</b>	92	3	

Date of Report: December 23, 2011  
Samples Submitted: December 14, 2011  
Laboratory Reference: 1112-101  
Project: 0180-292-00

**DISSOLVED METALS  
EPA 200.8/7470A  
METHOD BLANK QUALITY CONTROL**

Date Filtered: 12-14-11  
Date Analyzed: 12-15&20-11  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB1214F1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.0
Barium	200.8	<b>ND</b>	25
Cadmium	200.8	<b>ND</b>	4.0
Chromium	200.8	<b>ND</b>	10
Lead	200.8	<b>ND</b>	1.0
Mercury	7470A	<b>ND</b>	0.50
Selenium	200.8	<b>ND</b>	5.0
Silver	200.8	<b>ND</b>	10

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**DISSOLVED METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Filtered: 12-14-11  
 Date Analyzed: 12-15&20-11

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 12-091-02

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.0	
Barium	ND	ND	NA	25	
Cadmium	ND	ND	NA	4.0	
Chromium	ND	ND	NA	10	
Lead	ND	ND	NA	1.0	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.0	
Silver	ND	ND	NA	10	

Date of Report: December 23, 2011  
 Samples Submitted: December 14, 2011  
 Laboratory Reference: 1112-101  
 Project: 0180-292-00

**DISSOLVED METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Filtered: 12-14-11  
 Date Analyzed: 12-15&20-11

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 12-091-02

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	200	<b>209</b>	105	<b>205</b>	102	2	
Barium	200	<b>205</b>	102	<b>207</b>	104	1	
Cadmium	200	<b>207</b>	104	<b>204</b>	102	2	
Chromium	200	<b>188</b>	94	<b>188</b>	94	0	
Lead	200	<b>201</b>	101	<b>195</b>	98	3	
Mercury	12.5	<b>12.2</b>	98	<b>12.1</b>	97	1	
Selenium	200	<b>209</b>	104	<b>210</b>	105	1	
Silver	200	<b>180</b>	90	<b>181</b>	91	1	





### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference



**Onsite Environmental Inc.**  
 Analytical Laboratory Testing Services  
 14648 NE 95th Street • Redmond, WA 98052  
 Phone: (425) 883-3881 • www.onsite-env.com

**Chain of Custody**

Turnaround Request  
(in working days)

Laboratory Number:

12-101

(Check One)

Same Day  1 Day

2 Days  3 Days

Standard (7 Days) (TPH analysis 5 Days)

(other) \_\_\_\_\_

Company: *GeoEngineers*  
 Project Number: *0180-292-00*  
 Project Name: *WSDOT - Midway Metals*  
 Project Manager: *Aaron Messner*  
 Sampled by: \_\_\_\_\_

Lab ID	Sample Identification	Date Sampled	Time Sampled	Matrix	No. of Cont.	NWTPH-HCID	NWTPH-Gx/BTEX	NWTPH-Gx	NWTPH-Dx	Volatiles 8260B	Halogenated Volatiles 8260B	Semivolatiles 8270D/SIM (with low-level PAHs)	PAHs 8270D/SIM (low-level)	PCBs 8082	Organochlorine Pesticides 8081A	Organophosphorus Pesticides 8270D/SIM	Chlorinated Acid Herbicides 8151A	Total RCRA Metals	Total MTCA Metals	TCLP Metals	HEM (oil and grease) 1664	% Moisture
1	<i>Well 1-W</i>	<i>12/14/11</i>	<i>11:00 AM</i>	<i>W</i>	<i>13</i>			<i>X</i>	<i>X</i>	<i>X</i>	<i>X</i>	<i>X</i>	<i>X</i>	<i>X</i>				<i>X</i>	<i>X</i>			

	Signature	Company	Date	Time	Comments/Special Instructions
Relinquished	<i>[Signature]</i>	<i>Geo</i>	<i>12/14/11</i>	<i>1340</i>	<i>(X) Added 12/15/11. DB (STA)</i>
Received	<i>[Signature]</i>	<i>PEER</i>	<i>12/14/11</i>	<i>142</i>	
Relinquished	<i>[Signature]</i>	<i>Geo</i>	<i>12/14/11</i>	<i>1620</i>	
Received	<i>[Signature]</i>	<i>Geo</i>	<i>12/14/11</i>	<i>1620</i>	
Relinquished	<i>[Signature]</i>	<i>Geo</i>			
Received					
Reviewed/Date					



14648 NE 95<sup>th</sup> Street, Redmond, WA 98052 • (425) 883-3881

February 16, 2012

Aaron Waggoner  
GeoEngineers, Inc.  
1101 Fawcett Avenue South, Suite 200  
Tacoma, WA 98402

Re: Analytical Data for Project 0180-292-00  
Laboratory Reference No. 1202-084

Dear Aaron:

Enclosed are the analytical results and associated quality control data for samples submitted on February 9, 2012.

The standard policy of OnSite Environmental, Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister  
Project Manager

Enclosures

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-084  
Project: 0180-292-00

### **Case Narrative**

Samples were collected on February 8 and 9, 2012 and received by the laboratory on February 9, 2012. They were maintained at the laboratory at a temperature of 2°C to 6°C.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-084  
Project: 0180-292-00

**ANALYTICAL REPORT FOR SAMPLES**

<b>Client ID</b>	<b>Laboratory ID</b>	<b>Matrix</b>	<b>Date Sampled</b>	<b>Date Received</b>	<b>Notes</b>
DP12-W	02-084-01	Water	2-8-12	2-9-12	
DP15-W	02-084-02	Water	2-8-12	2-9-12	
SW9	02-084-03	Water	2-9-12	2-9-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**TOTAL METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-084-01					
<b>Client ID:</b>	<b>DP12-W</b>					
Arsenic	<b>12</b>	3.3	200.8	2-10-12	2-10-12	
Barium	<b>300</b>	28	200.8	2-10-12	2-10-12	
Cadmium	<b>ND</b>	4.4	200.8	2-10-12	2-10-12	
Chromium	<b>140</b>	22	200.8	2-10-12	2-10-12	
Lead	<b>12</b>	1.1	200.8	2-10-12	2-10-12	
Mercury	<b>ND</b>	0.50	7470A	2-10-12	2-10-12	
Selenium	<b>ND</b>	5.6	200.8	2-10-12	2-10-12	
Silver	<b>ND</b>	11	200.8	2-10-12	2-10-12	

Lab ID:	02-084-02					
<b>Client ID:</b>	<b>DP15-W</b>					
Arsenic	<b>27</b>	3.3	200.8	2-10-12	2-10-12	
Barium	<b>400</b>	28	200.8	2-10-12	2-10-12	
Cadmium	<b>ND</b>	4.4	200.8	2-10-12	2-10-12	
Chromium	<b>180</b>	22	200.8	2-10-12	2-10-12	
Lead	<b>16</b>	1.1	200.8	2-10-12	2-10-12	
Mercury	<b>ND</b>	0.50	7470A	2-10-12	2-10-12	
Selenium	<b>6.8</b>	5.6	200.8	2-10-12	2-10-12	
Silver	<b>ND</b>	11	200.8	2-10-12	2-10-12	

Date of Report: February 16, 2012  
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 Project: 0180-292-00

**TOTAL METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date	Date	Flags
				Prepared	Analyzed	
Lab ID:	02-084-03					
Client ID:	SW9					
Arsenic	ND	3.3	200.8	2-10-12	2-10-12	
Barium	ND	28	200.8	2-10-12	2-10-12	
Cadmium	ND	4.4	200.8	2-10-12	2-10-12	
Chromium	ND	11	200.8	2-10-12	2-10-12	
Lead	8.3	1.1	200.8	2-10-12	2-10-12	
Mercury	ND	0.50	7470A	2-10-12	2-10-12	
Selenium	ND	5.6	200.8	2-10-12	2-10-12	
Silver	ND	11	200.8	2-10-12	2-10-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**DISSOLVED METALS**  
**EPA 200.8/7470A**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	EPA Method	Date Prepared	Date Analyzed	Flags
Lab ID:	02-084-01					
<b>Client ID:</b>	<b>DP12-W</b>					
Arsenic	ND	3.0	200.8		2-10-12	
Barium	ND	25	200.8		2-10-12	
Cadmium	ND	4.0	200.8		2-10-12	
Chromium	ND	10	200.8		2-10-12	
Lead	ND	1.0	200.8		2-10-12	
Mercury	ND	0.50	7470A		2-10-12	
Selenium	ND	5.0	200.8		2-10-12	
Silver	ND	10	200.8		2-10-12	
Lab ID:	02-084-02					
<b>Client ID:</b>	<b>DP15-W</b>					
Arsenic	ND	3.0	200.8	2-9-12	2-10-12	
Barium	ND	25	200.8	2-9-12	2-10-12	
Cadmium	ND	4.0	200.8	2-9-12	2-10-12	
Chromium	ND	10	200.8	2-9-12	2-10-12	
Lead	ND	1.0	200.8	2-9-12	2-10-12	
Mercury	ND	0.50	7470A	2-9-12	2-10-12	
Selenium	ND	5.0	200.8	2-9-12	2-10-12	
Silver	ND	10	200.8	2-9-12	2-10-12	
Lab ID:	02-084-03					
<b>Client ID:</b>	<b>SW9</b>					
Arsenic	ND	3.0	200.8	2-9-12	2-10-12	
Barium	ND	25	200.8	2-9-12	2-10-12	
Cadmium	ND	4.0	200.8	2-9-12	2-10-12	
Chromium	ND	10	200.8	2-9-12	2-10-12	
Lead	1.6	1.0	200.8	2-9-12	2-10-12	
Mercury	ND	0.50	7470A	2-9-12	2-10-12	
Selenium	ND	5.0	200.8	2-9-12	2-10-12	
Silver	ND	10	200.8	2-9-12	2-10-12	



Date of Report: February 16, 2012  
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**NWTPH-Gx**

Matrix: Water  
 Units: ug/L (ppb)

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP12-W</b>					
Laboratory ID:	02-084-01					
Gasoline	<b>ND</b>	100	NWTPH-Gx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	73-121				
<b>Client ID:</b>	<b>DP15-W</b>					
Laboratory ID:	02-084-02					
Gasoline	<b>ND</b>	100	NWTPH-Gx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	88	73-121				
<b>Client ID:</b>	<b>SW9</b>					
Laboratory ID:	02-084-03					
Gasoline	<b>ND</b>	100	NWTPH-Gx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	88	73-121				

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 Project: 0180-292-00

**NWTPH-Dx**  
 (with acid/silica gel clean-up)

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-W</b>					
Laboratory ID:	02-084-01					
Diesel Range Organics	<b>ND</b>	0.27	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	0.43	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	98	50-150				
<b>Client ID:</b>	<b>DP15-W</b>					
Laboratory ID:	02-084-02					
Diesel Range Organics	<b>ND</b>	0.30	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	0.48	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	99	50-150				
<b>Client ID:</b>	<b>SW9</b>					
Laboratory ID:	02-084-03					
Diesel Range Organics	<b>ND</b>	0.27	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	0.43	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	103	50-150				

Date of Report: February 16, 2012  
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 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Matrix: Water  
 Units: ug/L

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP12-W</b>					
Laboratory ID:	02-084-01					
Dichlorodifluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloromethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Vinyl Chloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Trichlorofluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Acetone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Iodomethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Carbon Disulfide	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methylene Chloride	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Vinyl Acetate	ND	2.0	EPA 8260	2-10-12	2-10-12	
2,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Butanone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Bromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroform	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Carbon Tetrachloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Benzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Trichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Dibromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromodichloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	2-10-12	2-10-12	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Toluene	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	

Date of Report: February 16, 2012  
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**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-W</b>					
Laboratory ID:	02-084-01					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Tetrachloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Hexanone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Dibromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromoethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Ethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
m,p-Xylene	ND	0.40	EPA 8260	2-10-12	2-10-12	
o-Xylene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Styrene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromoform	ND	1.0	EPA 8260	2-10-12	2-10-12	
Isopropylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Propylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
4-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
tert-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
sec-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
p-Isopropyltoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Hexachlorobutadiene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Naphthalene	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>88</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>88</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>85</i>	<i>65-120</i>				

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**VOLATILES by EPA 8260B**  
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Matrix: Water  
 Units: ug/L

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP15-W</b>					
Laboratory ID:	02-084-02					
Dichlorodifluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloromethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Vinyl Chloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Trichlorofluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Acetone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Iodomethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Carbon Disulfide	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methylene Chloride	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Vinyl Acetate	ND	2.0	EPA 8260	2-10-12	2-10-12	
2,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Butanone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Bromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroform	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Carbon Tetrachloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Benzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Trichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Dibromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromodichloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	2-10-12	2-10-12	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Toluene	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	

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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP15-W</b>					
Laboratory ID:	02-084-02					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Tetrachloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Hexanone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Dibromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromoethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Ethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
m,p-Xylene	ND	0.40	EPA 8260	2-10-12	2-10-12	
o-Xylene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Styrene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromoform	ND	1.0	EPA 8260	2-10-12	2-10-12	
Isopropylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Propylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
4-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
tert-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
sec-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
p-Isopropyltoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Hexachlorobutadiene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Naphthalene	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>89</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>89</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>86</i>	<i>65-120</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
 Page 1 of 2

Matrix: Water  
 Units: ug/L

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SW9</b>					
Laboratory ID:	02-084-03					
Dichlorodifluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloromethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Vinyl Chloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Trichlorofluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Acetone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Iodomethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Carbon Disulfide	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methylene Chloride	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Vinyl Acetate	ND	2.0	EPA 8260	2-10-12	2-10-12	
2,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Butanone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Bromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroform	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Carbon Tetrachloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Benzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Trichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Dibromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromodichloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	2-10-12	2-10-12	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Toluene	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>SW9</b>					
Laboratory ID:	02-084-03					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Tetrachloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Hexanone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Dibromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromoethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Ethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
m,p-Xylene	ND	0.40	EPA 8260	2-10-12	2-10-12	
o-Xylene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Styrene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromoform	ND	1.0	EPA 8260	2-10-12	2-10-12	
Isopropylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Propylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
4-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
tert-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
sec-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
p-Isopropyltoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Hexachlorobutadiene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Naphthalene	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>87</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>87</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>86</i>	<i>65-120</i>				



Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
 page 1 of 2

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-W</b>					
Laboratory ID:	02-084-01					
n-Nitrosodimethylamine	ND	0.98	EPA 8270	2-10-12	2-10-12	
Pyridine	ND	0.98	EPA 8270	2-10-12	2-10-12	
Phenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
Aniline	ND	4.9	EPA 8270	2-10-12	2-10-12	
bis(2-Chloroethyl)ether	ND	0.98	EPA 8270	2-10-12	2-10-12	
2-Chlorophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,3-Dichlorobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,4-Dichlorobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
Benzyl alcohol	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,2-Dichlorobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
2-Methylphenol (o-Cresol)	ND	0.98	EPA 8270	2-10-12	2-10-12	
bis(2-Chloroisopropyl)ether	ND	0.98	EPA 8270	2-10-12	2-10-12	
(3+4)-Methylphenol (m,p-Cresol)	ND	0.98	EPA 8270	2-10-12	2-10-12	
n-Nitroso-di-n-propylamine	ND	0.98	EPA 8270	2-10-12	2-10-12	
Hexachloroethane	ND	0.98	EPA 8270	2-10-12	2-10-12	
Nitrobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
Isophorone	ND	0.98	EPA 8270	2-10-12	2-10-12	
2-Nitrophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,4-Dimethylphenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
bis(2-Chloroethoxy)methane	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,4-Dichlorophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,2,4-Trichlorobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
Naphthalene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
4-Chloroaniline	ND	0.98	EPA 8270	2-10-12	2-10-12	
Hexachlorobutadiene	ND	0.98	EPA 8270	2-10-12	2-10-12	
4-Chloro-3-methylphenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
2-Methylnaphthalene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
1-Methylnaphthalene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
Hexachlorocyclopentadiene	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,4,6-Trichlorophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,3-Dichloroaniline	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,4,5-Trichlorophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
2-Chloronaphthalene	ND	0.98	EPA 8270	2-10-12	2-10-12	
2-Nitroaniline	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,4-Dinitrobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
Dimethylphthalate	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,3-Dinitrobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,6-Dinitrotoluene	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,2-Dinitrobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
Acenaphthylene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
3-Nitroaniline	ND	0.98	EPA 8270	2-10-12	2-10-12	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**SEMIVOLATILES by EPA 8270D/SIM**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-W</b>					
Laboratory ID:	02-084-01					
2,4-Dinitrophenol	ND	4.9	EPA 8270	2-10-12	2-10-12	
Acenaphthene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
4-Nitrophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,4-Dinitrotoluene	ND	0.98	EPA 8270	2-10-12	2-10-12	
Dibenzofuran	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,3,5,6-Tetrachlorophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
2,3,4,6-Tetrachlorophenol	ND	0.98	EPA 8270	2-10-12	2-10-12	
Diethylphthalate	ND	0.98	EPA 8270	2-10-12	2-10-12	
4-Chlorophenyl-phenylether	ND	0.98	EPA 8270	2-10-12	2-10-12	
4-Nitroaniline	ND	0.98	EPA 8270	2-10-12	2-10-12	
Fluorene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
4,6-Dinitro-2-methylphenol	ND	4.9	EPA 8270	2-10-12	2-10-12	
n-Nitrosodiphenylamine	ND	0.98	EPA 8270	2-10-12	2-10-12	
1,2-Diphenylhydrazine	ND	0.98	EPA 8270	2-10-12	2-10-12	
4-Bromophenyl-phenylether	ND	0.98	EPA 8270	2-10-12	2-10-12	
Hexachlorobenzene	ND	0.98	EPA 8270	2-10-12	2-10-12	
Pentachlorophenol	ND	4.9	EPA 8270	2-10-12	2-10-12	
Phenanthrene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
Anthracene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
Carbazole	ND	0.98	EPA 8270	2-10-12	2-10-12	
Di-n-butylphthalate	ND	0.98	EPA 8270	2-10-12	2-10-12	
Fluoranthene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
Benzidine	ND	4.9	EPA 8270	2-10-12	2-10-12	
Pyrene	ND	0.098	EPA 8270/SIM	2-10-12	2-10-12	
Butylbenzylphthalate	ND	0.98	EPA 8270	2-10-12	2-10-12	
bis(2-Ethylhexyl)adipate	ND	4.9	EPA 8270	2-10-12	2-10-12	
3,3'-Dichlorobenzidine	ND	0.98	EPA 8270	2-10-12	2-10-12	
Benzo[a]anthracene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
Chrysene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
bis(2-Ethylhexyl)phthalate	ND	0.98	EPA 8270	2-10-12	2-10-12	
Di-n-octylphthalate	ND	0.98	EPA 8270	2-10-12	2-10-12	
Benzo[b]fluoranthene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
Benzo(j,k)fluoranthene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
Benzo[a]pyrene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
Indeno[1,2,3-cd]pyrene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
Dibenz[a,h]anthracene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
Benzo[g,h,i]perylene	ND	0.0098	EPA 8270/SIM	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>37</i>	<i>18 - 86</i>				
<i>Phenol-d6</i>	<i>30</i>	<i>10 - 88</i>				
<i>Nitrobenzene-d5</i>	<i>66</i>	<i>37 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>73</i>	<i>42 - 108</i>				
<i>2,4,6-Tribromophenol</i>	<i>71</i>	<i>39 - 118</i>				
<i>Terphenyl-d14</i>	<i>75</i>	<i>49 - 122</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

### PCBs by EPA 8082

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>Client ID:</b>	<b>DP12-W</b>					
Laboratory ID:	02-084-01					
Aroclor 1016	ND	0.048	EPA 8082	2-10-12	2-10-12	
Aroclor 1221	ND	0.048	EPA 8082	2-10-12	2-10-12	
Aroclor 1232	ND	0.048	EPA 8082	2-10-12	2-10-12	
Aroclor 1242	ND	0.048	EPA 8082	2-10-12	2-10-12	
Aroclor 1248	ND	0.048	EPA 8082	2-10-12	2-10-12	
Aroclor 1254	ND	0.048	EPA 8082	2-10-12	2-10-12	
Aroclor 1260	ND	0.048	EPA 8082	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	86	36-127				
<b>Client ID:</b>	<b>DP15-W</b>					
Laboratory ID:	02-084-02					
Aroclor 1016	ND	0.10	EPA 8082	2-10-12	2-10-12	
Aroclor 1221	ND	0.10	EPA 8082	2-10-12	2-10-12	
Aroclor 1232	ND	0.10	EPA 8082	2-10-12	2-10-12	
Aroclor 1242	ND	0.10	EPA 8082	2-10-12	2-10-12	
Aroclor 1248	ND	0.10	EPA 8082	2-10-12	2-10-12	
Aroclor 1254	ND	0.10	EPA 8082	2-10-12	2-10-12	
Aroclor 1260	ND	0.10	EPA 8082	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	81	36-127				
<b>Client ID:</b>	<b>SW9</b>					
Laboratory ID:	02-084-03					
Aroclor 1016	ND	0.053	EPA 8082	2-10-12	2-10-12	
Aroclor 1221	ND	0.053	EPA 8082	2-10-12	2-10-12	
Aroclor 1232	ND	0.053	EPA 8082	2-10-12	2-10-12	
Aroclor 1242	ND	0.053	EPA 8082	2-10-12	2-10-12	
Aroclor 1248	ND	0.053	EPA 8082	2-10-12	2-10-12	
Aroclor 1254	ND	0.053	EPA 8082	2-10-12	2-10-12	
Aroclor 1260	ND	0.053	EPA 8082	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
DCB	92	36-127				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Water  
 Units: ug/L

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>DP15-W</b>					
Laboratory ID:	02-084-02					
Naphthalene	<b>0.10</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
2-Methylnaphthalene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
1-Methylnaphthalene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Acenaphthylene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Acenaphthene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Fluorene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Phenanthrene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Anthracene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Fluoranthene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Pyrene	<b>ND</b>	0.092	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[a]anthracene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
Chrysene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[b]fluoranthene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
Benzo(j,k)fluoranthene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[a]pyrene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
Dibenz[a,h]anthracene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[g,h,i]perylene	<b>ND</b>	0.0092	EPA 8270/SIM	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>62</i>	<i>38 - 105</i>				
<i>Pyrene-d10</i>	<i>64</i>	<i>37 - 121</i>				
<i>Terphenyl-d14</i>	<i>78</i>	<i>32 - 112</i>				

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 (with silica gel clean-up)**

Matrix: Water

Units: ug/L

<b>Analyte</b>	<b>Result</b>	<b>PQL</b>	<b>Method</b>	<b>Date Prepared</b>	<b>Date Analyzed</b>	<b>Flags</b>
<b>Client ID:</b>	<b>SW9</b>					
Laboratory ID:	02-084-03					
Naphthalene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
2-Methylnaphthalene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
1-Methylnaphthalene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Acenaphthylene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Acenaphthene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Fluorene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Phenanthrene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Anthracene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Fluoranthene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Pyrene	<b>ND</b>	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[a]anthracene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Chrysene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[b]fluoranthene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo(j,k)fluoranthene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[a]pyrene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Indeno(1,2,3-c,d)pyrene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Dibenz[a,h]anthracene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[g,h,i]perylene	<b>ND</b>	0.010	EPA 8270/SIM	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>74</i>	<i>38 - 105</i>				
<i>Pyrene-d10</i>	<i>84</i>	<i>37 - 121</i>				
<i>Terphenyl-d14</i>	<i>83</i>	<i>32 - 112</i>				

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-084  
Project: 0180-292-00

**TOTAL METALS  
EPA 200.8  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-10-12  
Date Analyzed: 2-10-12  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB0210WM1

Analyte	Method	Result	PQL
Arsenic	200.8	<b>ND</b>	3.3
Barium	200.8	<b>ND</b>	28
Cadmium	200.8	<b>ND</b>	4.4
Chromium	200.8	<b>ND</b>	22
Lead	200.8	<b>ND</b>	1.1
Selenium	200.8	<b>ND</b>	5.6
Silver	200.8	<b>ND</b>	11

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-084  
Project: 0180-292-00

**TOTAL MERCURY  
EPA 7470A  
METHOD BLANK QUALITY CONTROL**

Date Extracted: 2-10-12

Date Analyzed: 2-10-12

Matrix: Water

Units: ug/L (ppb)

Lab ID: MB0210W1

Analyte	Method	Result	PQL
Mercury	7470A	<b>ND</b>	0.50

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8  
 DUPLICATE QUALITY CONTROL**

Date Extracted: 2-10-12

Date Analyzed: 2-10-12

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-084-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	11.8	11.4	3	3.3	
Barium	299	270	10	28	
Cadmium	ND	ND	NA	4.4	
Chromium	138	118	16	22	
Lead	12.0	11.6	4	1.1	
Selenium	ND	ND	NA	5.6	
Silver	ND	ND	NA	11	



Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-084  
Project: 0180-292-00

**TOTAL MERCURY  
EPA 7470A  
DUPLICATE QUALITY CONTROL**

Date Extracted: 2-10-12

Date Analyzed: 2-10-12

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-028-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Mercury	<b>ND</b>	<b>ND</b>	NA	0.50	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**TOTAL METALS  
 EPA 200.8  
 MS/MSD QUALITY CONTROL**

Date Extracted: 2-10-12

Date Analyzed: 2-10-12

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-084-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	222	<b>218</b>	93	<b>219</b>	93	0	
Barium	222	<b>493</b>	87	<b>482</b>	82	2	
Cadmium	222	<b>218</b>	98	<b>218</b>	98	0	
Chromium	222	<b>308</b>	77	<b>314</b>	79	2	
Lead	222	<b>238</b>	102	<b>229</b>	98	4	
Selenium	222	<b>200</b>	90	<b>202</b>	91	1	
Silver	222	<b>211</b>	95	<b>214</b>	97	2	

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-084  
Project: 0180-292-00

**TOTAL MERCURY  
EPA 7470A  
MS/MSD QUALITY CONTROL**

Date Extracted: 2-10-12

Date Analyzed: 2-10-12

Matrix: Water

Units: ug/L (ppb)

Lab ID: 02-028-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Mercury	12.5	<b>13.9</b>	111	<b>13.4</b>	107	4	

Date of Report: February 16, 2012  
Samples Submitted: February 9, 2012  
Laboratory Reference: 1202-084  
Project: 0180-292-00

**DISSOLVED METALS  
EPA 200.8/7470A  
METHOD BLANK QUALITY CONTROL**

Date Filtered: 2-9-12  
Date Analyzed: 2-10-12  
  
Matrix: Water  
Units: ug/L (ppb)  
  
Lab ID: MB0209F1

Analyte	Method	Result	PQL
Arsenic	200.8	ND	3.0
Barium	200.8	ND	25
Cadmium	200.8	ND	4.0
Chromium	200.8	ND	10
Lead	200.8	ND	1.0
Mercury	7470A	ND	0.50
Selenium	200.8	ND	5.0
Silver	200.8	ND	10

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**DISSOLVED METALS  
 EPA 200.8/7470A  
 DUPLICATE QUALITY CONTROL**

Date Filtered: 2-9-12  
 Date Analyzed: 2-10-12  
  
 Matrix: Water  
 Units: ug/L (ppb)  
  
 Lab ID: 02-084-01

Analyte	Sample Result	Duplicate Result	RPD	PQL	Flags
Arsenic	ND	ND	NA	3.0	
Barium	ND	ND	NA	25	
Cadmium	ND	ND	NA	4.0	
Chromium	ND	ND	NA	10	
Lead	ND	ND	NA	1.0	
Mercury	ND	ND	NA	0.50	
Selenium	ND	ND	NA	5.0	
Silver	ND	ND	NA	10	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**DISSOLVED METALS  
 EPA 200.8/7470A  
 MS/MSD QUALITY CONTROL**

Date Filtered: 2-9-12  
 Date Analyzed: 2-10-12

Matrix: Water  
 Units: ug/L (ppb)

Lab ID: 02-084-01

Analyte	Spike Level	MS	Percent Recovery	MSD	Percent Recovery	RPD	Flags
Arsenic	200	<b>203</b>	101	<b>203</b>	101	0	
Barium	200	<b>191</b>	95	<b>190</b>	95	1	
Cadmium	200	<b>206</b>	103	<b>203</b>	101	2	
Chromium	200	<b>203</b>	101	<b>203</b>	101	0	
Lead	200	<b>207</b>	103	<b>206</b>	103	0	
Mercury	12.5	<b>12.4</b>	99	<b>12.4</b>	100	0	
Selenium	200	<b>208</b>	104	<b>204</b>	102	2	
Silver	200	<b>200</b>	100	<b>198</b>	99	1	

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**NWTPH-Gx  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0210W1					
Gasoline	<b>ND</b>	100	NWTPH-Gx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Fluorobenzene</i>	90	73-121				

Analyte	Result	Spike Level	Source Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>								
Laboratory ID:	02-084-01							
	ORIG	DUP						
Gasoline	<b>ND</b>	<b>ND</b>	NA	NA	NA	NA	NA	30
<i>Surrogate:</i>								
<i>Fluorobenzene</i>				90	90	73-121		

Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**NWTPH-Dx  
 QUALITY CONTROL  
 (with acid/silica gel clean-up)**

Matrix: Water  
 Units: mg/L (ppm)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0210W1					
Diesel Range Organics	<b>ND</b>	0.25	NWTPH-Dx	2-10-12	2-10-12	
Lube Oil Range Organics	<b>ND</b>	0.40	NWTPH-Dx	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>o-Terphenyl</i>	92	50-150				

Analyte	Result	Percent Recovery	Recovery Limits	RPD	RPD Limit	Flags
<b>DUPLICATE</b>						
Laboratory ID:	02-073-01					
	ORIG	DUP				
Diesel Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
Lube Oil Range Organics	<b>ND</b>	<b>ND</b>		NA	NA	
<i>Surrogate:</i>						
<i>o-Terphenyl</i>			97 98	50-150		



Date of Report: February 16, 2012  
 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**

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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0210W1					
Dichlorodifluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloromethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Vinyl Chloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Trichlorofluoromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Acetone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Iodomethane	ND	1.0	EPA 8260	2-10-12	2-10-12	
Carbon Disulfide	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methylene Chloride	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl t-Butyl Ether	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Vinyl Acetate	ND	2.0	EPA 8260	2-10-12	2-10-12	
2,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
(cis) 1,2-Dichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Butanone	ND	5.0	EPA 8260	2-10-12	2-10-12	
Bromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chloroform	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Carbon Tetrachloride	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Benzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Trichloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Dibromomethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromodichloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chloroethyl Vinyl Ether	ND	1.0	EPA 8260	2-10-12	2-10-12	
(cis) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Methyl Isobutyl Ketone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Toluene	ND	1.0	EPA 8260	2-10-12	2-10-12	
(trans) 1,3-Dichloropropene	ND	0.20	EPA 8260	2-10-12	2-10-12	

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**VOLATILES by EPA 8260B**  
**METHOD BLANK QUALITY CONTROL**  
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Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0210W1					
1,1,2-Trichloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Tetrachloroethene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Hexanone	ND	2.0	EPA 8260	2-10-12	2-10-12	
Dibromochloromethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromoethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Chlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,1,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
Ethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
m,p-Xylene	ND	0.40	EPA 8260	2-10-12	2-10-12	
o-Xylene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Styrene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromoform	ND	1.0	EPA 8260	2-10-12	2-10-12	
Isopropylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Bromobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,1,2,2-Tetrachloroethane	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichloropropane	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Propylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
2-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
4-Chlorotoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3,5-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
tert-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2,4-Trimethylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
sec-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,3-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
p-Isopropyltoluene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,4-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
n-Butylbenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
1,2-Dibromo-3-chloropropane	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,4-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Hexachlorobutadiene	ND	0.20	EPA 8260	2-10-12	2-10-12	
Naphthalene	ND	1.0	EPA 8260	2-10-12	2-10-12	
1,2,3-Trichlorobenzene	ND	0.20	EPA 8260	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>Dibromofluoromethane</i>	<i>89</i>	<i>68-120</i>				
<i>Toluene-d8</i>	<i>91</i>	<i>73-120</i>				
<i>4-Bromofluorobenzene</i>	<i>88</i>	<i>65-120</i>				

Date of Report: February 16, 2012  
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 Laboratory Reference: 1202-084  
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**VOLATILES by EPA 8260B**  
**SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Limit	Flags
					Recovery	Limits	RPD				
<b>SPIKE BLANKS</b>											
Laboratory ID:	SB0210W1										
	SB	SBD	SB	SBD	SB	SBD					
1,1-Dichloroethene	11.2	10.8	10.0	10.0	112	108	70-130	4		11	
Benzene	10.3	10.2	10.0	10.0	103	102	75-123	1		8	
Trichloroethene	10.4	10.5	10.0	10.0	104	105	80-113	1		9	
Toluene	10.4	10.5	10.0	10.0	104	105	80-113	1		8	
Chlorobenzene	11.2	11.0	10.0	10.0	112	110	80-115	2		8	
<i>Surrogate:</i>											
Dibromofluoromethane					86	89	68-120				
Toluene-d8					87	91	73-120				
4-Bromofluorobenzene					87	91	65-120				

Date of Report: February 16, 2012  
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**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
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Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0210W1					
n-Nitrosodimethylamine	ND	1.0	EPA 8270	2-10-12	2-10-12	
Pyridine	ND	1.0	EPA 8270	2-10-12	2-10-12	
Phenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
Aniline	ND	5.0	EPA 8270	2-10-12	2-10-12	
bis(2-Chloroethyl)ether	ND	1.0	EPA 8270	2-10-12	2-10-12	
2-Chlorophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,3-Dichlorobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,4-Dichlorobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
Benzyl alcohol	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,2-Dichlorobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
2-Methylphenol (o-Cresol)	ND	1.0	EPA 8270	2-10-12	2-10-12	
bis(2-Chloroisopropyl)ether	ND	1.0	EPA 8270	2-10-12	2-10-12	
(3+4)-Methylphenol (m,p-Cresol)	ND	1.0	EPA 8270	2-10-12	2-10-12	
n-Nitroso-di-n-propylamine	ND	1.0	EPA 8270	2-10-12	2-10-12	
Hexachloroethane	ND	1.0	EPA 8270	2-10-12	2-10-12	
Nitrobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
Isophorone	ND	1.0	EPA 8270	2-10-12	2-10-12	
2-Nitrophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,4-Dimethylphenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
bis(2-Chloroethoxy)methane	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,4-Dichlorophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,2,4-Trichlorobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
Naphthalene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
4-Chloroaniline	ND	1.0	EPA 8270	2-10-12	2-10-12	
Hexachlorobutadiene	ND	1.0	EPA 8270	2-10-12	2-10-12	
4-Chloro-3-methylphenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
2-Methylnaphthalene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
1-Methylnaphthalene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
Hexachlorocyclopentadiene	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,4,6-Trichlorophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,3-Dichloroaniline	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,4,5-Trichlorophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
2-Chloronaphthalene	ND	1.0	EPA 8270	2-10-12	2-10-12	
2-Nitroaniline	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,4-Dinitrobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
Dimethylphthalate	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,3-Dinitrobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,6-Dinitrotoluene	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,2-Dinitrobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
Acenaphthylene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
3-Nitroaniline	ND	1.0	EPA 8270	2-10-12	2-10-12	

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**SEMIVOLATILES by EPA 8270D/SIM**  
**METHOD BLANK QUALITY CONTROL**  
 page 2 of 2

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0210W1					
2,4-Dinitrophenol	ND	5.0	EPA 8270	2-10-12	2-10-12	
Acenaphthene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
4-Nitrophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,4-Dinitrotoluene	ND	1.0	EPA 8270	2-10-12	2-10-12	
Dibenzofuran	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,3,5,6-Tetrachlorophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
2,3,4,6-Tetrachlorophenol	ND	1.0	EPA 8270	2-10-12	2-10-12	
Diethylphthalate	ND	1.0	EPA 8270	2-10-12	2-10-12	
4-Chlorophenyl-phenylether	ND	1.0	EPA 8270	2-10-12	2-10-12	
4-Nitroaniline	ND	1.0	EPA 8270	2-10-12	2-10-12	
Fluorene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
4,6-Dinitro-2-methylphenol	ND	5.0	EPA 8270	2-10-12	2-10-12	
n-Nitrosodiphenylamine	ND	1.0	EPA 8270	2-10-12	2-10-12	
1,2-Diphenylhydrazine	ND	1.0	EPA 8270	2-10-12	2-10-12	
4-Bromophenyl-phenylether	ND	1.0	EPA 8270	2-10-12	2-10-12	
Hexachlorobenzene	ND	1.0	EPA 8270	2-10-12	2-10-12	
Pentachlorophenol	ND	5.0	EPA 8270	2-10-12	2-10-12	
Phenanthrene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
Anthracene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
Carbazole	ND	1.0	EPA 8270	2-10-12	2-10-12	
Di-n-butylphthalate	ND	1.0	EPA 8270	2-10-12	2-10-12	
Fluoranthene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
Benzidine	ND	5.0	EPA 8270	2-10-12	2-10-12	
Pyrene	ND	0.10	EPA 8270/SIM	2-10-12	2-10-12	
Butylbenzylphthalate	ND	1.0	EPA 8270	2-10-12	2-10-12	
bis-2-Ethylhexyladipate	ND	5.0	EPA 8270	2-10-12	2-10-12	
3,3'-Dichlorobenzidine	ND	1.0	EPA 8270	2-10-12	2-10-12	
Benzo[a]anthracene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
Chrysene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
bis(2-Ethylhexyl)phthalate	ND	1.0	EPA 8270	2-10-12	2-10-12	
Di-n-octylphthalate	ND	1.0	EPA 8270	2-10-12	2-10-12	
Benzo[b]fluoranthene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
Benzo[a]pyrene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
Indeno[1,2,3-cd]pyrene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270/SIM	2-10-12	2-10-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorophenol</i>	<i>49</i>	<i>18 - 86</i>				
<i>Phenol-d6</i>	<i>38</i>	<i>10 - 88</i>				
<i>Nitrobenzene-d5</i>	<i>70</i>	<i>37 - 112</i>				
<i>2-Fluorobiphenyl</i>	<i>73</i>	<i>42 - 108</i>				
<i>2,4,6-Tribromophenol</i>	<i>89</i>	<i>39 - 118</i>				
<i>Terphenyl-d14</i>	<i>80</i>	<i>49 - 122</i>				

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**SEMIVOLATILES by EPA 8270D/SIM  
 SB/SBD QUALITY CONTROL**

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					Recovery	Limits	Limits	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0210W1									
	SB	SBD	SB	SBD	SB	SBD				
Phenol	16.9	19.1	40.0	40.0	42	48	26 - 60	12	29	
2-Chlorophenol	32.2	35.2	40.0	40.0	81	88	46 - 104	9	34	
1,4-Dichlorobenzene	15.2	16.5	20.0	20.0	76	83	46 - 92	8	29	
n-Nitroso-di-n-propylamine	15.9	18.0	20.0	20.0	80	90	30 - 102	12	25	
1,2,4-Trichlorobenzene	15.7	17.1	20.0	20.0	79	86	45 - 92	9	25	
4-Chloro-3-methylphenol	35.2	39.6	40.0	40.0	88	99	53 - 108	12	18	
Acenaphthene	16.7	19.1	20.0	20.0	84	96	57 - 103	13	15	
4-Nitrophenol	24.3	27.0	40.0	40.0	61	68	21 - 76	11	33	
2,4-Dinitrotoluene	18.1	21.7	20.0	20.0	91	109	60 - 118	18	20	
Pentachlorophenol	35.9	43.6	40.0	40.0	90	109	48 - 119	19	31	
Pyrene	17.6	20.4	20.0	20.0	88	102	62 - 111	15	19	
<i>Surrogate:</i>										
2-Fluorophenol					51	57	18 - 86			
Phenol-d6					40	44	10 - 88			
Nitrobenzene-d5					70	80	37 - 112			
2-Fluorobiphenyl					70	82	42 - 108			
2,4,6-Tribromophenol					80	94	39 - 118			
Terphenyl-d14					75	86	49 - 122			

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**PCBs by EPA 8082  
 QUALITY CONTROL**

Matrix: Water  
 Units: ug/L (ppb)

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
<b>METHOD BLANK</b>						
Laboratory ID:	MB0210W1					
Aroclor 1016	ND	0.050	EPA 8082	2-10-12	2-10-12	
Aroclor 1221	ND	0.050	EPA 8082	2-10-12	2-10-12	
Aroclor 1232	ND	0.050	EPA 8082	2-10-12	2-10-12	
Aroclor 1242	ND	0.050	EPA 8082	2-10-12	2-10-12	
Aroclor 1248	ND	0.050	EPA 8082	2-10-12	2-10-12	
Aroclor 1254	ND	0.050	EPA 8082	2-10-12	2-10-12	
Aroclor 1260	ND	0.050	EPA 8082	2-10-12	2-10-12	
Surrogate:	<i>Percent Recovery</i>		<i>Control Limits</i>			
DCB	80		36-127			

Analyte	Result		Spike Level		Source Result	Percent Recovery		Recovery Limits	RPD	RPD Limit	Flags
<b>SPIKE BLANKS</b>											
Laboratory ID:	SB0210W1										
	SB	SBD	SB	SBD		SB	SBD				
Aroclor 1260	0.389	0.431	0.500	0.500	N/A	78	86	57-122	10	11	
Surrogate:											
DCB						83	85	36-127			

Date of Report: February 16, 2012  
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 Project: 0180-292-00

**PAHs by EPA 8270D/SIM  
 METHOD BLANK QUALITY CONTROL  
 (with silica gel clean-up)**

Matrix: Water  
 Units: ug/L

Analyte	Result	PQL	Method	Date Prepared	Date Analyzed	Flags
Laboratory ID:	MB0213W1					
Naphthalene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
2-Methylnaphthalene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
1-Methylnaphthalene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Acenaphthylene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Acenaphthene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Fluorene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Phenanthrene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Anthracene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Fluoranthene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Pyrene	ND	0.10	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[a]anthracene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Chrysene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[b]fluoranthene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo(j,k)fluoranthene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[a]pyrene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Indeno(1,2,3-c,d)pyrene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Dibenz[a,h]anthracene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
Benzo[g,h,i]perylene	ND	0.010	EPA 8270/SIM	2-13-12	2-13-12	
<i>Surrogate:</i>	<i>Percent Recovery</i>	<i>Control Limits</i>				
<i>2-Fluorobiphenyl</i>	<i>78</i>	<i>38 - 105</i>				
<i>Pyrene-d10</i>	<i>90</i>	<i>37 - 121</i>				
<i>Terphenyl-d14</i>	<i>85</i>	<i>32 - 112</i>				



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 Samples Submitted: February 9, 2012  
 Laboratory Reference: 1202-084  
 Project: 0180-292-00

**PAHs by EPA 8270D/SIM**  
**SB/SBD QUALITY CONTROL**  
 (with silica gel clean-up)

Matrix: Water  
 Units: ug/L

Analyte	Result		Spike Level		Percent Recovery		Recovery	RPD	RPD	Flags
					Recovery	Limits	Limits	Limit		
<b>SPIKE BLANKS</b>										
Laboratory ID:	SB0213W1									
	SB	SBD	SB	SBD	SB	SBD				
Naphthalene	<b>0.352</b>	<b>0.360</b>	0.500	0.500	70	72	38 - 110	2	35	
Acenaphthylene	<b>0.350</b>	<b>0.395</b>	0.500	0.500	70	79	47 - 120	12	30	
Acenaphthene	<b>0.349</b>	<b>0.374</b>	0.500	0.500	70	75	46 - 113	7	26	
Fluorene	<b>0.373</b>	<b>0.404</b>	0.500	0.500	75	81	58 - 104	8	25	
Phenanthrene	<b>0.370</b>	<b>0.390</b>	0.500	0.500	74	78	61 - 99	5	19	
Anthracene	<b>0.378</b>	<b>0.404</b>	0.500	0.500	76	81	55 - 122	7	19	
Fluoranthene	<b>0.434</b>	<b>0.459</b>	0.500	0.500	87	92	58 - 129	6	18	
Pyrene	<b>0.423</b>	<b>0.442</b>	0.500	0.500	85	88	57 - 126	4	25	
Benzo[a]anthracene	<b>0.408</b>	<b>0.447</b>	0.500	0.500	82	89	51 - 124	9	18	
Chrysene	<b>0.398</b>	<b>0.412</b>	0.500	0.500	80	82	53 - 123	3	20	
Benzo[b]fluoranthene	<b>0.497</b>	<b>0.526</b>	0.500	0.500	99	105	53 - 126	6	18	
Benzo(j,k)fluoranthene	<b>0.461</b>	<b>0.489</b>	0.500	0.500	92	98	51 - 126	6	23	
Benzo[a]pyrene	<b>0.392</b>	<b>0.410</b>	0.500	0.500	78	82	52 - 127	4	21	
Indeno(1,2,3-c,d)pyrene	<b>0.416</b>	<b>0.428</b>	0.500	0.500	83	86	49 - 123	3	26	
Dibenz[a,h]anthracene	<b>0.444</b>	<b>0.465</b>	0.500	0.500	89	93	39 - 125	5	31	
Benzo[g,h,i]perylene	<b>0.441</b>	<b>0.466</b>	0.500	0.500	88	93	40 - 125	6	30	
<i>Surrogate:</i>										
<i>2-Fluorobiphenyl</i>					<i>63</i>	<i>69</i>	<i>38 - 105</i>			
<i>Pyrene-d10</i>					<i>84</i>	<i>86</i>	<i>37 - 121</i>			
<i>Terphenyl-d14</i>					<i>78</i>	<i>77</i>	<i>32 - 112</i>			



### Data Qualifiers and Abbreviations

- A - Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B - The analyte indicated was also found in the blank sample.
- C - The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E - The value reported exceeds the quantitation range and is an estimate.
- F - Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H - The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I - Compound recovery is outside of the control limits.
- J - The value reported was below the practical quantitation limit. The value is an estimate.
- K - Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L - The RPD is outside of the control limits.
- M - Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 - Hydrocarbons in the gasoline range (toluene-naphthalene) are present in the sample.
- N - Hydrocarbons in the lube oil range are impacting the diesel range result.
- N1 - Hydrocarbons in diesel range are impacting lube oil range results.
- O - Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P - The RPD of the detected concentrations between the two columns is greater than 40.
- Q - Surrogate recovery is outside of the control limits.
- S - Surrogate recovery data is not available due to the necessary dilution of the sample.
- T - The sample chromatogram is not similar to a typical \_\_\_\_\_.
- U - The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 - The practical quantitation limit is elevated due to interferences present in the sample.
- V - Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W - Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X - Sample extract treated with a mercury cleanup procedure.
- Y - Sample extract treated with an acid/silica gel cleanup procedure.
- Z -
- ND - Not Detected at PQL
- PQL - Practical Quantitation Limit
- RPD - Relative Percent Difference

