

**Kennedy/Jenks Consultants**  
**Engineers & Scientists**

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13 February 2017

Mr. Shane DeGross  
Manager Environmental Remediation  
BNSF Railway Company  
605 Puyallup Avenue  
Tacoma, Washington 98421

Subject: November 2016 Groundwater Compliance Monitoring Report  
BNSF D Street Oil Pipeline Site  
Tacoma, Washington  
K/J Project 1696039.00

Dear Mr. DeGross:

This report presents the results of the November 2016 groundwater compliance monitoring completed at the BNSF Railway Company (BNSF) D Street Oil Pipeline Site in Tacoma, Washington (Site). The Site is located along the former alignment of an oil pipeline that was used to transport bunker oil from bulk storage tanks located north of the intersection of East 15th Street and D Street, to what is now the Tacoma Rail Yard located south and east of the intersection of East 21st Street and East D Street (Figure 1).

### Background and Purpose

This groundwater sampling event was conducted as required by the Washington State Department of Ecology (Ecology) as part of a 5-year (Periodic) review to be conducted in 2016 in accordance with Washington Administrative Code (WAC) 173-340-420. The groundwater compliance monitoring is intended to monitor chemicals of concern (COCs) and natural attenuation parameters in groundwater following completion of the final remedial action in December 2009 (GeoEngineers 2010). The monitoring activities and this report are required by a Consent Decree (CD), dated 5 August 2008, between BNSF and Ecology (Pierce County Superior Court Action No. 08 2 11105 4).

The concentrations of COCs are evaluated at the Conditional Point of Compliance (CPOC) relative to Site cleanup levels (CULs) identified in the *Final Remedial Investigation (RI)/Feasibility Study (FS) Report* dated 4 April 2007 and the *Final Cleanup Action Plan (CAP)* dated 14 November 2007 (GeoEngineers 2007). Monitoring wells MW-2 through MW-5, MW-8, MW-9, and MW-11, collectively known as 'the D Street wells', serve to define the CPOC identified in the CAP (Figure 2).

Mr. Shane DeGross  
BNSF Railway Company  
13 February 2017  
Page 2

The RI/FS and CAP identified the COCs for groundwater at the Site as petroleum hydrocarbons [quantified as Bunker Range Petroleum Hydrocarbons (BRPH)], carcinogenic polycyclic aromatic hydrocarbons (cPAHs), and naphthalene. Model Toxics Control Act (MTCA) CULs for groundwater were selected for use at this Site, as defined in the CAP, and are summarized as:

- Petroleum hydrocarbons (quantified as BRPH): 0.5 milligrams per liter (mg/L)
- Total cPAHs (as determined using toxicity equivalency methodology as required in WAC 173-340-708(8)): 0.1 micrograms per liter ( $\mu\text{g/L}$ )
- Naphthalene: 160  $\mu\text{g/L}$ .

## Ecology Correspondence

In June 2011, the Ecology Site Manager, Mr. Marv Coleman, approved a request to discontinue annual groundwater monitoring at the Site. As part of that approval, Ecology noted that because the Site has institutional controls applied to it, a 5-year review will be required (Ecology 2011). According to Ecology, prior to the 5-year review, an additional round of groundwater sampling is required, along with an inspection of physical barriers put in place during remedial excavations. The planned installation locations of these physical barriers were provided in the *Engineering Design Report* (EDR) and were to be located surrounding the western ends of the Home Electric Building and Supervalu Building, along East D Street between East 18th Street and East 19th Street (GeoEngineers 2008).

Ecology indicated in a 2015 email that they would consider approving abandonment of the D Street wells after the 5-year review was completed (Ecology 2015a). This is consistent with previous correspondence with GeoEngineers' staff, where Ecology indicated they would consider terminating the CD after the 5-year review (GeoEngineers 2011).

In a follow-up letter to Mr. Bruce Sheppard (BNSF), dated 29 April 2015, Ecology indicated no further remedial action was necessary at the Site by some of the potentially liable parties (PLPs), including the Washington State Department of Transportation, the City of Tacoma, John Backus Home Electric, and 1815 D Street LLC. This NFA did not apply to BNSF, and the letter noted "*further remedial action is still necessary under MTCA and the consent decree to control and monitor remaining contamination and periodically review conditions at the site*" (Ecology 2015b). The letter noted that Ecology considers cleanup required at the Site under the CD to be complete, but post-cleanup actions are still required, including sampling of the D Street wells as part of the 5-year review, and inspecting physical barriers put in place during remedial excavations.

## Field Methods

Per Ecology's request, a Site inspection was conducted on 19 August 2016. This included a visual inspection for evidence of recent ground disturbance activities in areas where subsurface physical barriers were installed during remedial excavations. The condition of Site monitoring

Mr. Shane DeGross  
BNSF Railway Company  
13 February 2017  
Page 3

wells was also noted during the inspection. No evidence of recent ground disturbance was observed in areas surrounding the subsurface physical barriers.

Because the D Street wells are located in traffic lanes along East D Street, accessing the wells requires a Right-of-Way (ROW) permit from the City of Tacoma. Prior to conducting well development and sampling activities, Kennedy/Jenks Consultants obtained ROW Permit No. RCON16-0282 and subcontracted with a licensed traffic control services provider.

The D Street monitoring wells had not been accessed or sampled since 2011; as such, they were redeveloped prior to sampling on 21 November 2016. Development of the wells was accomplished by a combination of surging and over-pumping each well to remove accumulated sediment and clean the well screen. Wells MW-2, MW-4, MW-5, MW-8, MW-9, and MW-11 were purged using a 12-volt submersible well pump and dedicated polyethylene tubing. The submersible pump was decontaminated between each well using an Alconox® detergent wash and tap water rinse, followed by a distilled water rinse. Monitoring well MW-3 could not be purged using the submersible pump because the casing was partially blocked approximately 3 feet below the top of casing. The source of the obstruction could not be determined, but the casing was still open enough to allow passage of a water level meter and polyethylene tubing. As such, well MW-3 was surged as much as possible using the polyethylene tubing, and was purged using a peristaltic pump until the water was clear. Wells MW-2, MW-4, MW-5, MW-8, MW-9, and MW-11 were purged until each well was producing clear water after surging. Approximately 5 well volumes of water were purged from each well. Purge water was placed into 55-gallon drums, labeled, and temporarily stored at the nearby Tacoma Rail Yard for offsite disposal.

Groundwater samples were collected for laboratory analysis in general accordance with the 9 July 2008 Compliance Monitoring Plan (CMP) as presented in the EDR. The November 2016 groundwater compliance monitoring samples were collected on 30 November 2016, primarily during an outgoing tide, as measured by the National Oceanic and Atmospheric Administration's Commencement Bay Tidal Station 9446484. Although the CMP does not specify when groundwater samples will be collected in regards to tidal stage, samples were collected primarily during an outgoing tide to limit the potential for tidally-induced groundwater gradients to bias conditions observed at the CPOC.

The groundwater samples were collected following low-flow sampling guidelines to minimize the suspension of particulates in the samples. Groundwater samples were collected using a peristaltic pump with new, dedicated polyethylene tubing at each well. Groundwater samples were collected by pumping at flow rates of less than 0.5 liters per minute from the approximate mid-point of the screened interval in each well.

A YSI water quality meter was used with an inline flow-through cell to monitor electrical conductivity, dissolved oxygen (DO), pH, oxidation-reduction potential (ORP) and temperature. Turbidity was measured using a stand-alone Micro TPI turbidity meter. As noted in the CMP, ambient groundwater conditions were assumed to have been reached once three consecutive

Mr. Shane DeGross  
BNSF Railway Company  
13 February 2017  
Page 4

measurements of temperature, pH, and conductivity were within 10 percent of each other, and the turbidity is approximately 5 nephelometric turbidity units (NTUs) or less.

Once the water quality parameters had stabilized, the flow-through-cell was disconnected and the groundwater samples were collected in laboratory-provided containers. The samples were placed into a cooler with ice and shipped under chain-of-custody protocol to ESC Lab Sciences of Mt. Juliet, Tennessee, a Washington State accredited environmental laboratory. The groundwater samples were submitted for analysis of:

- Diesel Range Petroleum Hydrocarbons by Method NWTPH-Dx with silica gel cleanup
- PAHs by U.S. Environmental Protection Agency (EPA) Method 8270D-SIM
- Manganese by EPA Method 6020
- Nitrates by EPA Method 9056A
- Sulfates by EPA Method 9056A
- Total Alkalinity by EPA Method 2320 B-2011.

Natural attenuation parameter data were collected to monitor Site conditions as described in the CMP.

Purge water from well development and groundwater sampling was placed into labeled 55-gallon drums for future waste characterization and offsite disposal.

## Analytical Results

Monitoring wells MW-2 through MW-5 and MW-8, MW-9, and MW-11 were sampled on 30 November 2016. Analytical results are summarized in Table 1, and the laboratory analytical report is included as Attachment A. A data validation summary is included in Attachment B.

### **Petroleum Hydrocarbons**

Diesel-range and residual-range petroleum hydrocarbons were not detected at concentrations greater than the laboratory reporting limit in any of the groundwater samples collected during the November 2016 groundwater compliance monitoring.

### **cPAHs**

cPAHs were not detected at concentrations greater than the laboratory reporting limit in any of the groundwater samples collected during the November 2016 groundwater compliance monitoring.

Mr. Shane DeGross  
BNSF Railway Company  
13 February 2017  
Page 5

## Naphthalene

Naphthalene, 1-methyl-naphthalene, and 2-methyl-naphthalene were not detected at concentrations greater than the laboratory reporting limit in any of the groundwater samples collected during the November 2016 groundwater compliance monitoring.

## Natural Attenuation Parameters

The primary natural attenuation parameters (pH, conductivity, temperature, DO, and ORP) collected in the field during the November 2016 groundwater compliance monitoring were variable across the Site and similar to results from the last monitoring event in February 2011 (GeoEngineers 2011). The measured parameters continue to indicate conditions are generally favorable for natural attenuation of any residual hydrocarbons that may potentially remain at the Site.

Secondary natural attenuation parameters (manganese, total alkalinity, nitrate, and sulfate) are also variable and similar to those observed during the last groundwater compliance monitoring event in 2011. These parameters also continue to indicate conditions are favorable for natural attenuation of any residual BRPHs that potentially remain at the Site.

## Conclusion

Diesel and heavy oil range hydrocarbons, PAHs, and cPAHs were not detected above the laboratory reporting limits in any of the groundwater samples collected from the D Street wells during the November 2016 groundwater compliance monitoring event. The results of the November 2016 groundwater compliance monitoring are consistent with the results of the 2010 and 2011 events. The November 2016 groundwater compliance monitoring results indicate that the Site is in compliance with the MTCA CULs established in the CD.

Based on our evaluation of the historical compliance monitoring results, it is our opinion that the conditions for site cleanup as described in both the CD and the CMP have been met. As it appears the goals identified in the CD and the CMP have been achieved, we recommend discontinuing the groundwater compliance monitoring program, finalizing the 5-year review, permanently abandoning the D Street monitoring wells, and terminating the CD.

Very truly yours,

KENNEDY/JENKS CONSULTANTS



Nathan Moxley, LG, LHG  
Project Manager



Ty C. Schreiner  
Vice President

cc: Marv Coleman, Washington State Department of Ecology

Mr. Shane DeGross  
BNSF Railway Company  
13 February 2017  
Page 6

**Attachments:**

Table 1 Summary of Groundwater Compliance Monitoring Analytical Results  
Table 2 Summary of Groundwater Compliance Monitoring Natural Attenuation Parameters  
Figure 1: Vicinity Map  
Figure 2: Site Map  
Attachment A: Laboratory Analytical Report and Chain-of-Custody Documentation  
Attachment B: Data Quality Assessment Summary

**References**

GeoEngineers. 2007. Draft Final Cleanup Action Plan; The BNSF Oil Pipeline Site, Tacoma, Washington. 14 November.

GeoEngineers. 2008. Engineering Design Report, The BNSF Oil Pipeline, Tacoma, Washington. 9 July.

GeoEngineers. 2010. Final Remedial Action Report; The BNSF Oil Pipeline Site, Tacoma, Washington. 28 May.

GeoEngineers. 2011. Email from Nick Rohrbach (GeoEngineers) to Matt Wells (Tupper Mack Wells PLLC) and Bruce Sheppard (BNSF) summarizing conversation with Marv Coleman (Ecology). 7 July 2011.

Washington State Department of Ecology. 2011. Letter to Bruce Sheppard re: February 2011 Groundwater Compliance Monitoring Report. BNSF Oil Pipeline Site, Tacoma, WA, 15 June 2011. BNSF D Street Pipeline FSID# 2923141.

Washington State Department of Ecology. 2015a. Email from Marv Coleman (Ecology) to Laura Himes (Kennedy/Jenks Consultants). 20 May 2015.

Washington State Department of Ecology. 2015b. Letter to Bruce Sheppard re: Status of Consent Decree No 08 2 1115 4 and No Further Action to complete Cleanup of the BNSF D Street Pipeline Site. 29 April 2015.

## Tables

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**TABLE 1**

**SUMMARY OF GROUNDWATER COMPLIANCE MONITORING ANALYTICAL RESULTS  
November 2016 Compliance Monitoring  
November 30, 2016  
BNSF D Street Pipeline, Tacoma, Washington**

Monitoring Well ID	Total Petroleum Hydrocarbons <sup>(a)</sup> (mg/L)		Carcinogenic Polycyclic Aromatic Hydrocarbons (cPAHs) <sup>(b)</sup> (µg/L)							PAHs <sup>(b)</sup> (µg/L)		
	Diesel Range	Heavy Oil Range	Benzo(a)anthracene	Benzo(a)pyrene	Benzo(b)fluoranthene	Benzo(k)fluoranthene	Chrysene	Dibenz(a,h)anthracene	Indeno (1,2,3-cd)pyrene	1-Methyl-naphthalene	2-Methyl-naphthalene	Naphthalene
MW-2	0.250 U	0.500 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.500 U	0.500 U	0.500 U
D-1	0.250 U	0.500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.250 U	0.250 U	0.250 U
MW-3	0.250 U	0.500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.250 U	0.250 U	0.250 U
MW-4	0.250 U	0.500 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.500 U	0.500 U	0.500 U
MW-5	0.250 U	0.500 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.500 U	0.500 U	0.500 U
MW-8	0.250 U	0.500 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.500 U	0.500 U	0.500 U
MW-9	0.250 U	0.500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.0500 U	0.250 U	0.250 U	0.250 U
MW-11	0.250 U	0.500 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.100 U	0.500 U	0.500 U	0.500 U
MTCA Method A Cleanup Levels	0.5	0.5	0.1 <sup>(c)</sup>							160 <sup>(d)</sup>		

Notes:

(a) Analyzed by Washington State Department of Ecology Method NWTPH-Dx.

(b) Analyzed by U.S. Environmental Protection Agency (EPA) Method 8270-SIM.

(c) Cleanup level for cPAHs is defined as the total toxic equivalent concentration of benzo(a)pyrene, as defined in WAC 173-340-708(8)(e).

(d) Cleanup level is a total value for naphthalene, 1-methyl naphthalene, and 2-methyl naphthalene.

Duplicate sample 'D-1' was collected from MW-2.

mg/L = milligrams per liter

µg/L = micrograms per liter

MTCA = Washington State Department of Ecology Model Toxics Control Act (WAC 173-340)

U = Not detected at or greater than the laboratory reporting limits

**TABLE 2**

**SUMMARY OF GROUNDWATER COMPLIANCE MONITORING NATURAL ATTENUATION PARAMETERS  
November 2016 Compliance Monitoring  
November 30, 2016  
BNSF D Street Pipeline, Tacoma, Washington**

Monitoring Well ID	Water Quality Parameters						Natural Attenuation Parameters (mg/L)				
	pH	Conductivity (mS/cm)	Turbidity (NTU)	Temperature (°C)	Dissolved Oxygen (mg/L)	ORP (mV)	Manganese (total)	Manganese (dissolved)	Alkalinity	Nitrate	Sulfate
MW-2	6.7	556.8	20.86	13.36	0.41	-102.8	1.170	1.260	314	0.102	77.3
D-1	--	--	--	--	--	--	1.140	1.260	258	0.104	77.6
MW-3	6.6	4991	6.5	12.85	0.98	-108.1	2.090	2.400	332	0.100 U	59.6
MW-4	6.3	2708	4.45	12.29	0.90	-185.0	1.460	1.390	323	0.100 U	5 U
MW-5	6.6	4870	0.76	12.58	1.04	-239.2	0.643	0.648	391	0.100 U	18.8
MW-8	6.4	3318	0.82	12.38	0.64	-201.6	0.304	0.332	252	0.100 U	1,360
MW-9	7.0	1451	6.9	13.22	0.93	-154.8	0.260	0.251	349	0.100 U	5 U
MW-11	6.4	2884	0.22	11.46	0.89	-282.0	0.569	0.586	311	11.500	1,100

Notes:

Duplicate sample 'D-1' was collected from MW-2.

mS/cm = microsiemens per centimeter

NTU = Nephelometric Turbidity Unit

°C = degrees celcius

ORP = oxidation-reduction potential

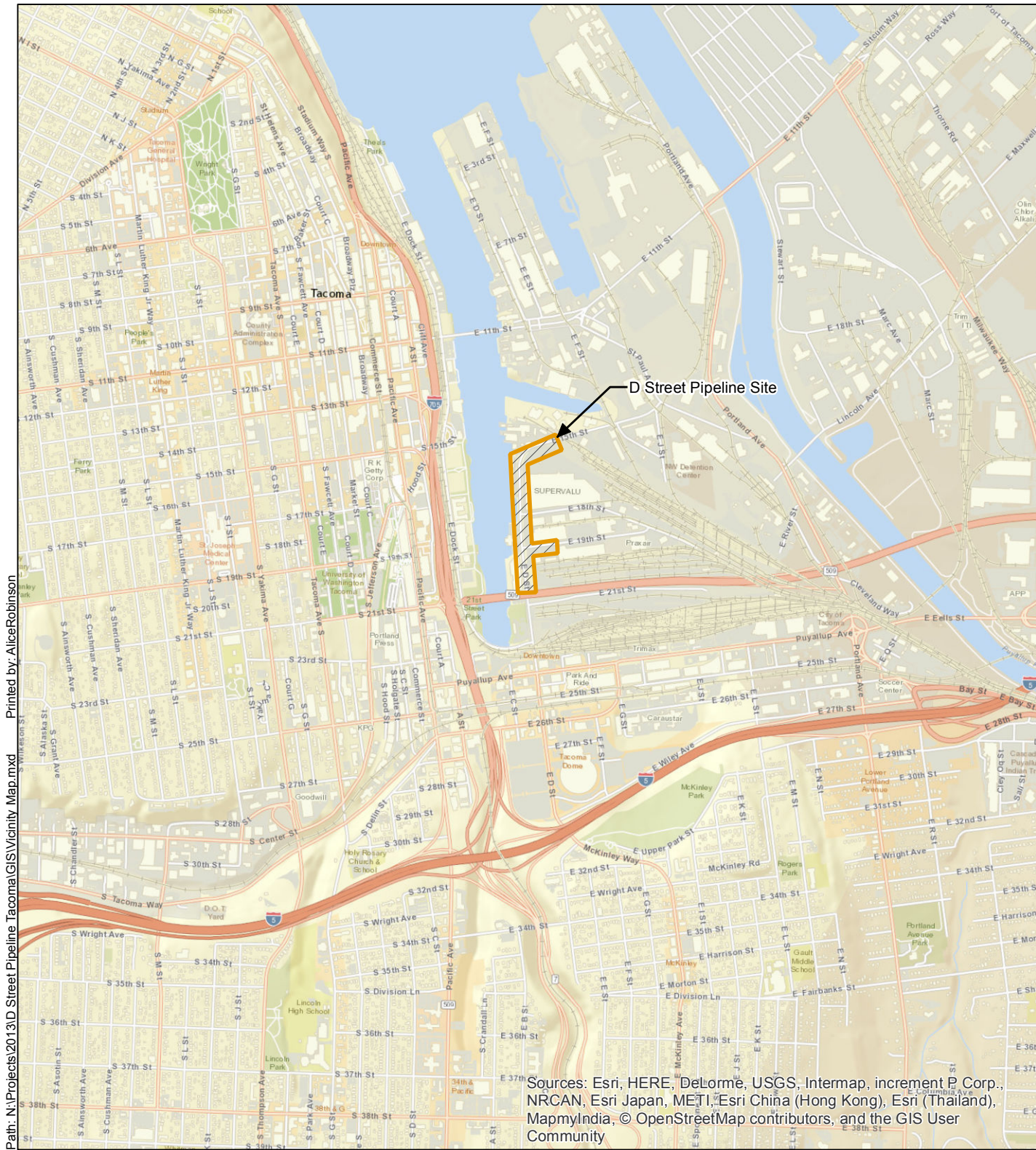
mg/L = milligrams per liter

mV = millivolts


U = Not detected at or above laboratory reporting limits

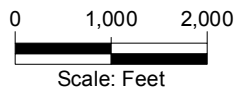
## Figures

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**Legend**

-  Approximate Site Extent



**Kennedy/Jenks Consultants**

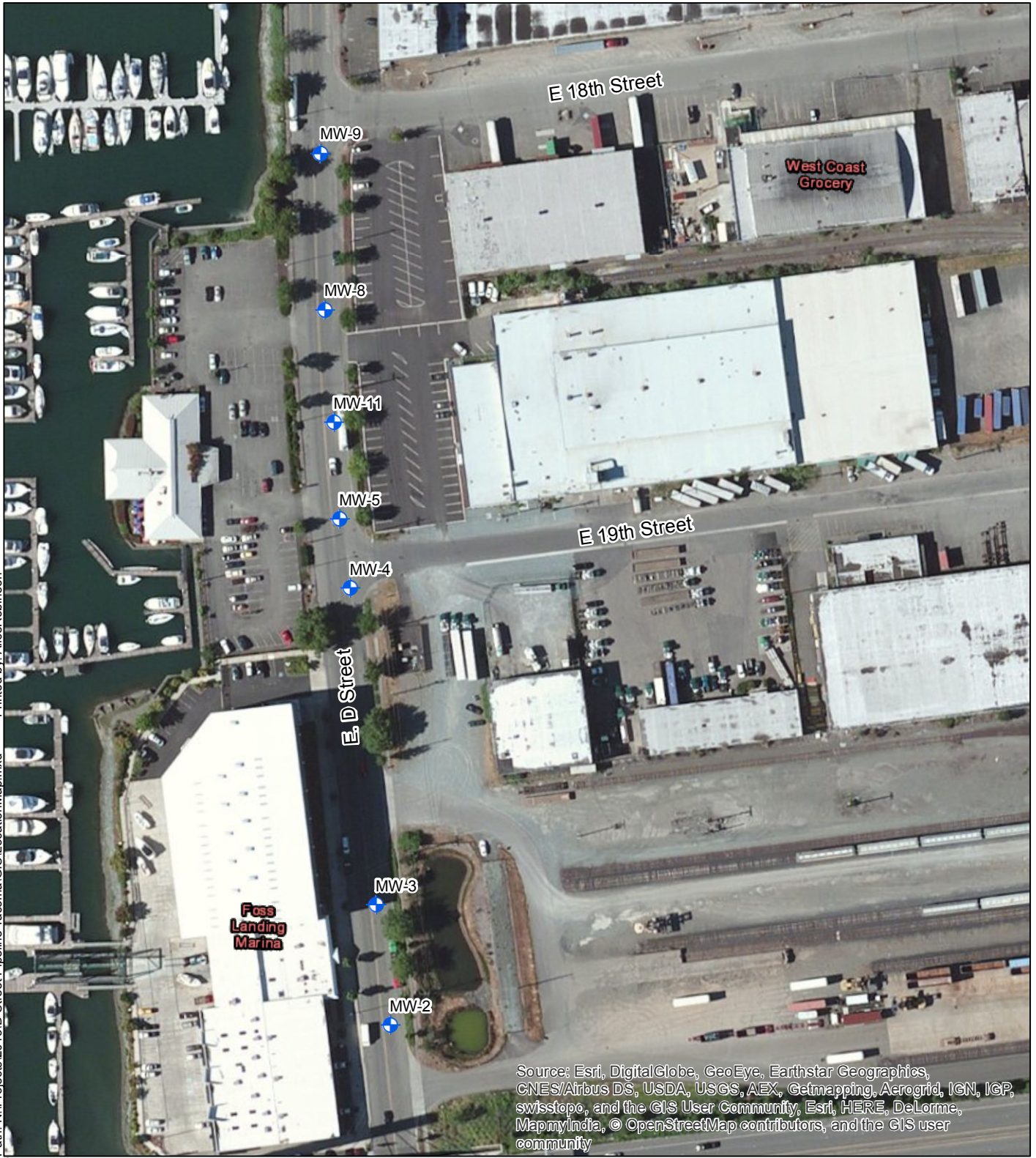
BNSF D Street Oil Pipeline Site  
Tacoma, Washington

**Vicinity Map**

1696039°00  
December 2016

**Figure 1**

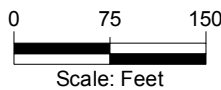
Path: N:\Projects\2013\D\_Street\_Pipeline\_Tacoma\GIS\LocationMap.mxd Printed by: AliceRobinson



Source: Esri, DigitalGlobe, GeoEye, Earthstar Geographics, CNES/Airbus DS, USDA, USGS, AEX, Getmapping, Aerogrid, IGN, IGP, swisstopo, and the GIS User Community, Esri, HERE, DeLorme, MapmyIndia, © OpenStreetMap contributors, and the GIS user community

**Legend**

◆ Monitoring Well Location and Designation



**Kennedy/Jenks Consultants**

BNSF D Street Oil Pipeline Site  
Tacoma, Washington

**Site Map**

1696039°00  
December 2016

**Figure 2**

## Attachment A

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Laboratory Analytical Report and Chain-of-Custody Documentation

December 09, 2016

## Kennedy/Jenks Con-BNSF Region 1

Sample Delivery Group: L875470  
Samples Received: 12/01/2016  
Project Number: 1696039.00  
Description: BNSF - D St. Wells  
Site: D STREET PIPELINE  
Report To: Nathan Moxley  
32001 32nd Avenue South, Ste 100  
Federal Way, WA 98001

Entire Report Reviewed By:



Mark W. Beasley  
Technical Service Representative

Results relate only to the items tested or calibrated and are reported as rounded values. This test report shall not be reproduced, except in full, without written approval of the laboratory. Where applicable, sampling conducted by ESC is performed per guidance provided in laboratory standard operating procedures: 060302, 060303, and 060304.



<b><sup>1</sup>Cp: Cover Page</b>	<b>1</b>	
<b><sup>2</sup>Tc: Table of Contents</b>	<b>2</b>	
<b><sup>3</sup>Ss: Sample Summary</b>	<b>3</b>	
<b><sup>4</sup>Cn: Case Narrative</b>	<b>5</b>	
<b><sup>5</sup>Sr: Sample Results</b>	<b>6</b>	
MW-9-20161130 L875470-01	6	
MW-8-20161130 L875470-02	7	
MW-11-20161130 L875470-03	8	
MW-5-20161130 L875470-04	9	
MW-4-20161130 L875470-05	10	
MW-3-20161130 L875470-06	11	
MW-2-20161130 L875470-07	12	
D-1-20161130 L875470-08	13	
<b><sup>6</sup>Qc: Quality Control Summary</b>	<b>14</b>	
Wet Chemistry by Method 2320 B-2011	14	
Wet Chemistry by Method 9056A	15	
Metals (ICPMS) by Method 6020	18	
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	20	
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	21	
<b><sup>7</sup>Gl: Glossary of Terms</b>	<b>23</b>	
<b><sup>8</sup>Al: Accreditations &amp; Locations</b>	<b>24</b>	
<b><sup>9</sup>Sc: Chain of Custody</b>	<b>25</b>	

# SAMPLE SUMMARY



## MW-9-20161130 L875470-01 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 08:55  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 14:58	LAT
Metals (ICPMS) by Method 6020	WG931186	1	12/01/16 20:20	12/02/16 12:19	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	1	12/04/16 07:13	12/05/16 02:59	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 11:54	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 07:48	12/02/16 07:48	AMC
Wet Chemistry by Method 9056A	WG931112	1	12/01/16 14:17	12/01/16 14:17	SAM

1  
Cp

2  
Tc

3  
Ss

4  
Cn

## MW-8-20161130 L875470-02 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 10:00  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 15:02	LAT
Metals (ICPMS) by Method 6020	WG931186	1	12/01/16 20:20	12/02/16 10:42	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	2	12/04/16 07:13	12/05/16 03:20	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 12:10	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 07:57	12/02/16 07:57	AMC
Wet Chemistry by Method 9056A	WG931112	1	12/01/16 14:47	12/01/16 14:47	SAM
Wet Chemistry by Method 9056A	WG931112	50	12/01/16 15:03	12/01/16 15:03	SAM

5  
Sr

6  
Qc

7  
Gl

8  
Al

9  
Sc

## MW-11-20161130 L875470-03 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 10:40  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 15:12	LAT
Metals (ICPMS) by Method 6020	WG931186	1	12/01/16 20:20	12/02/16 12:23	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	2	12/04/16 07:13	12/05/16 03:42	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 12:27	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 08:05	12/02/16 08:05	AMC
Wet Chemistry by Method 9056A	WG931112	5	12/01/16 15:18	12/01/16 15:18	SAM
Wet Chemistry by Method 9056A	WG931536	20	12/03/16 12:57	12/03/16 12:57	SAM

## MW-5-20161130 L875470-04 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 11:15  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 14:40	LAT
Metals (ICPMS) by Method 6020	WG931186	1	12/01/16 20:20	12/02/16 12:26	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	2	12/04/16 07:13	12/05/16 04:04	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 12:44	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 08:14	12/02/16 08:14	AMC
Wet Chemistry by Method 9056A	WG931112	1	12/01/16 16:04	12/01/16 16:04	SAM

## MW-4-20161130 L875470-05 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 11:55  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 15:16	LAT
Metals (ICPMS) by Method 6020	WG931186	1	12/01/16 20:20	12/02/16 12:30	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	2	12/04/16 07:13	12/05/16 04:26	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 13:00	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 08:23	12/02/16 08:23	AMC
Wet Chemistry by Method 9056A	WG931112	1	12/01/16 16:35	12/01/16 16:35	SAM

# SAMPLE SUMMARY



## MW-3-20161130 L875470-06 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 12:30  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 15:19	LAT
Metals (ICPMS) by Method 6020	WG931186	5	12/01/16 20:20	12/02/16 14:23	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	1	12/04/16 07:13	12/05/16 04:48	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 13:17	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 08:31	12/02/16 08:31	AMC
Wet Chemistry by Method 9056A	WG931112	1	12/01/16 16:50	12/01/16 16:50	SAM

1  
Cp

2  
Tc

3  
Ss

4  
Cn

## MW-2-20161130 L875470-07 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 13:05  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 15:23	LAT
Metals (ICPMS) by Method 6020	WG931186	1	12/01/16 20:20	12/02/16 12:37	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	2	12/04/16 07:13	12/05/16 05:09	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 13:33	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 10:27	12/02/16 10:27	AMC
Wet Chemistry by Method 9056A	WG931112	1	12/01/16 17:06	12/01/16 17:06	SAM

5  
Sr

6  
Qc

7  
Gl

8  
Al

9  
Sc

## D-1-20161130 L875470-08 GW

Collected by  
Alexander Leshner  
Collected date/time  
11/30/16 00:00  
Received date/time  
12/01/16 09:00

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Metals (ICPMS) by Method 6020	WG931185	1	12/02/16 12:44	12/03/16 15:27	LAT
Metals (ICPMS) by Method 6020	WG931186	1	12/01/16 20:20	12/02/16 13:39	LAT
Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM	WG931662	1	12/04/16 07:13	12/05/16 05:31	FMB
Semi-Volatile Organic Compounds (GC) by Method NWTPHDX	WG931238	1	12/02/16 08:12	12/03/16 13:50	TRF
Wet Chemistry by Method 2320 B-2011	WG931302	1	12/02/16 10:37	12/02/16 10:37	AMC
Wet Chemistry by Method 9056A	WG931112	1	12/01/16 17:21	12/01/16 17:21	SAM



All sample aliquots were received at the correct temperature, in the proper containers, with the appropriate preservatives, and within method specified holding times. All MDL (LOD) and RDL (LOQ) values reported for environmental samples have been corrected for the dilution factor used in the analysis. All Method and Batch Quality Control are within established criteria except where addressed in this case narrative, a non-conformance form or properly qualified within the sample results. By my digital signature below, I affirm to the best of my knowledge, all problems/anomalies observed by the laboratory as having the potential to affect the quality of the data have been identified by the laboratory, and no information or data have been knowingly withheld that would affect the quality of the data.

Mark W. Beasley  
Technical Service Representative

- <sup>1</sup> Cp
- <sup>2</sup> Tc
- <sup>3</sup> Ss
- <sup>4</sup> Cn
- <sup>5</sup> Sr
- <sup>6</sup> Qc
- <sup>7</sup> Gl
- <sup>8</sup> Al
- <sup>9</sup> Sc



## Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Alkalinity	349000		20000	1	12/02/2016 07:48	<a href="#">WG931302</a>

## Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Nitrate	ND		100	1	12/01/2016 14:17	<a href="#">WG931112</a>
Sulfate	ND		5000	1	12/01/2016 14:17	<a href="#">WG931112</a>

## Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Manganese	260		5.00	1	12/03/2016 14:58	<a href="#">WG931185</a>
Manganese,Dissolved	251		5.00	1	12/02/2016 12:19	<a href="#">WG931186</a>

## Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 11:54	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 11:54	<a href="#">WG931238</a>
(S) o-Terphenyl	127		50.0-150		12/03/2016 11:54	<a href="#">WG931238</a>

## Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Anthracene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Acenaphthene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Acenaphthylene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Chrysene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Fluoranthene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Fluorene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Naphthalene	ND		0.250	1	12/05/2016 02:59	<a href="#">WG931662</a>
Phenanthrene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
Pyrene	ND		0.0500	1	12/05/2016 02:59	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.250	1	12/05/2016 02:59	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.250	1	12/05/2016 02:59	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.250	1	12/05/2016 02:59	<a href="#">WG931662</a>
(S) Nitrobenzene-d5	98.4		45.1-170		12/05/2016 02:59	<a href="#">WG931662</a>
(S) 2-Fluorobiphenyl	103		57.7-153		12/05/2016 02:59	<a href="#">WG931662</a>
(S) p-Terphenyl-d14	103		53.2-156		12/05/2016 02:59	<a href="#">WG931662</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Alkalinity	252000		20000	1	12/02/2016 07:57	<a href="#">WG931302</a>

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Nitrate	ND		100	1	12/01/2016 14:47	<a href="#">WG931112</a>
Sulfate	1360000		250000	50	12/01/2016 15:03	<a href="#">WG931112</a>

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Manganese	304		5.00	1	12/03/2016 15:02	<a href="#">WG931185</a>
Manganese,Dissolved	332	<u>Q1</u>	5.00	1	12/02/2016 10:42	<a href="#">WG931186</a>

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 12:10	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 12:10	<a href="#">WG931238</a>
(S) o-Terphenyl	130		50.0-150		12/03/2016 12:10	<a href="#">WG931238</a>

Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Anthracene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Acenaphthene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Acenaphthylene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Chrysene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Fluoranthene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Fluorene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Naphthalene	ND		0.500	2	12/05/2016 03:20	<a href="#">WG931662</a>
Phenanthrene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
Pyrene	ND		0.100	2	12/05/2016 03:20	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.500	2	12/05/2016 03:20	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.500	2	12/05/2016 03:20	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.500	2	12/05/2016 03:20	<a href="#">WG931662</a>
(S) Nitrobenzene-d5	90.2		45.1-170		12/05/2016 03:20	<a href="#">WG931662</a>
(S) 2-Fluorobiphenyl	93.0		57.7-153		12/05/2016 03:20	<a href="#">WG931662</a>
(S) p-Terphenyl-d14	101		53.2-156		12/05/2016 03:20	<a href="#">WG931662</a>

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



## Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Alkalinity	311000		20000	1	12/02/2016 08:05	<a href="#">WG931302</a>

## Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Nitrate	11500		500	5	12/01/2016 15:18	<a href="#">WG931112</a>
Sulfate	1100000		100000	20	12/03/2016 12:57	<a href="#">WG931536</a>

## Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Manganese	569		5.00	1	12/03/2016 15:12	<a href="#">WG931185</a>
Manganese,Dissolved	586		5.00	1	12/02/2016 12:23	<a href="#">WG931186</a>

## Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 12:27	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 12:27	<a href="#">WG931238</a>
<i>(S) o-Terphenyl</i>	128		50.0-150		12/03/2016 12:27	<a href="#">WG931238</a>

## Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Anthracene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Acenaphthene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Acenaphthylene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Chrysene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Fluoranthene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Fluorene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Naphthalene	ND		0.500	2	12/05/2016 03:42	<a href="#">WG931662</a>
Phenanthrene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
Pyrene	ND		0.100	2	12/05/2016 03:42	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.500	2	12/05/2016 03:42	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.500	2	12/05/2016 03:42	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.500	2	12/05/2016 03:42	<a href="#">WG931662</a>
<i>(S) Nitrobenzene-d5</i>	88.5		45.1-170		12/05/2016 03:42	<a href="#">WG931662</a>
<i>(S) 2-Fluorobiphenyl</i>	92.7		57.7-153		12/05/2016 03:42	<a href="#">WG931662</a>
<i>(S) p-Terphenyl-d14</i>	99.9		53.2-156		12/05/2016 03:42	<a href="#">WG931662</a>

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Alkalinity	391000		20000	1	12/02/2016 08:14	<a href="#">WG931302</a>

1 Cp

2 Tc

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Nitrate	ND		100	1	12/01/2016 16:04	<a href="#">WG931112</a>
Sulfate	18800		5000	1	12/01/2016 16:04	<a href="#">WG931112</a>

3 Ss

4 Cn

5 Sr

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Manganese	643	V	5.00	1	12/03/2016 14:40	<a href="#">WG931185</a>
Manganese,Dissolved	648		5.00	1	12/02/2016 12:26	<a href="#">WG931186</a>

6 Qc

7 Gl

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 12:44	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 12:44	<a href="#">WG931238</a>
(S) o-Terphenyl	133		50.0-150		12/03/2016 12:44	<a href="#">WG931238</a>

8 Al

9 Sc

Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Anthracene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Acenaphthene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Acenaphthylene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Chrysene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Fluoranthene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Fluorene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Naphthalene	ND		0.500	2	12/05/2016 04:04	<a href="#">WG931662</a>
Phenanthrene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
Pyrene	ND		0.100	2	12/05/2016 04:04	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.500	2	12/05/2016 04:04	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.500	2	12/05/2016 04:04	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.500	2	12/05/2016 04:04	<a href="#">WG931662</a>
(S) Nitrobenzene-d5	87.1		45.1-170		12/05/2016 04:04	<a href="#">WG931662</a>
(S) 2-Fluorobiphenyl	84.7		57.7-153		12/05/2016 04:04	<a href="#">WG931662</a>
(S) p-Terphenyl-d14	81.4		53.2-156		12/05/2016 04:04	<a href="#">WG931662</a>



## Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
Alkalinity	ug/l		ug/l		date / time	
Alkalinity	323000		20000	1	12/02/2016 08:23	<a href="#">WG931302</a>

## Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
Nitrate	ug/l		ug/l		date / time	
Nitrate	ND		100	1	12/01/2016 16:35	<a href="#">WG931112</a>
Sulfate	ND		5000	1	12/01/2016 16:35	<a href="#">WG931112</a>

## Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
Manganese	ug/l		ug/l		date / time	
Manganese	1460		5.00	1	12/03/2016 15:16	<a href="#">WG931185</a>
Manganese,Dissolved	1390		5.00	1	12/02/2016 12:30	<a href="#">WG931186</a>

## Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
Diesel Range Organics (DRO)	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 13:00	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 13:00	<a href="#">WG931238</a>
(S) o-Terphenyl	123		50.0-150		12/03/2016 13:00	<a href="#">WG931238</a>

## Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
Anthracene	ug/l		ug/l		date / time	
Anthracene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Acenaphthene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Acenaphthylene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Chrysene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Fluoranthene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Fluorene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Naphthalene	ND		0.500	2	12/05/2016 04:26	<a href="#">WG931662</a>
Phenanthrene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
Pyrene	ND		0.100	2	12/05/2016 04:26	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.500	2	12/05/2016 04:26	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.500	2	12/05/2016 04:26	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.500	2	12/05/2016 04:26	<a href="#">WG931662</a>
(S) Nitrobenzene-d5	88.3		45.1-170		12/05/2016 04:26	<a href="#">WG931662</a>
(S) 2-Fluorobiphenyl	95.1		57.7-153		12/05/2016 04:26	<a href="#">WG931662</a>
(S) p-Terphenyl-d14	90.7		53.2-156		12/05/2016 04:26	<a href="#">WG931662</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



## Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Alkalinity	332000		20000	1	12/02/2016 08:31	<a href="#">WG931302</a>

## Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Nitrate	ND		100	1	12/01/2016 16:50	<a href="#">WG931112</a>
Sulfate	59600		5000	1	12/01/2016 16:50	<a href="#">WG931112</a>

## Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Manganese	2090		5.00	1	12/03/2016 15:19	<a href="#">WG931185</a>
Manganese,Dissolved	2400		25.0	5	12/02/2016 14:23	<a href="#">WG931186</a>

## Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 13:17	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 13:17	<a href="#">WG931238</a>
<i>(S) o-Terphenyl</i>	122		50.0-150		12/03/2016 13:17	<a href="#">WG931238</a>

## Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Anthracene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Acenaphthene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Acenaphthylene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Chrysene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Fluoranthene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Fluorene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Naphthalene	ND		0.250	1	12/05/2016 04:48	<a href="#">WG931662</a>
Phenanthrene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
Pyrene	ND		0.0500	1	12/05/2016 04:48	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.250	1	12/05/2016 04:48	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.250	1	12/05/2016 04:48	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.250	1	12/05/2016 04:48	<a href="#">WG931662</a>
<i>(S) Nitrobenzene-d5</i>	97.9		45.1-170		12/05/2016 04:48	<a href="#">WG931662</a>
<i>(S) 2-Fluorobiphenyl</i>	106		57.7-153		12/05/2016 04:48	<a href="#">WG931662</a>
<i>(S) p-Terphenyl-d14</i>	102		53.2-156		12/05/2016 04:48	<a href="#">WG931662</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Alkalinity	314000		20000	1	12/02/2016 10:27	<a href="#">WG931302</a>

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Nitrate	102		100	1	12/01/2016 17:06	<a href="#">WG931112</a>
Sulfate	77300		5000	1	12/01/2016 17:06	<a href="#">WG931112</a>

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Manganese	1170		5.00	1	12/03/2016 15:23	<a href="#">WG931185</a>
Manganese,Dissolved	1260		5.00	1	12/02/2016 12:37	<a href="#">WG931186</a>

Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 13:33	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 13:33	<a href="#">WG931238</a>
<i>(S) o-Terphenyl</i>	126		50.0-150		12/03/2016 13:33	<a href="#">WG931238</a>

Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Anthracene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Acenaphthene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Acenaphthylene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Chrysene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Fluoranthene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Fluorene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Naphthalene	ND		0.500	2	12/05/2016 05:09	<a href="#">WG931662</a>
Phenanthrene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
Pyrene	ND		0.100	2	12/05/2016 05:09	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.500	2	12/05/2016 05:09	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.500	2	12/05/2016 05:09	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.500	2	12/05/2016 05:09	<a href="#">WG931662</a>
<i>(S) Nitrobenzene-d5</i>	93.4		45.1-170		12/05/2016 05:09	<a href="#">WG931662</a>
<i>(S) 2-Fluorobiphenyl</i>	99.8		57.7-153		12/05/2016 05:09	<a href="#">WG931662</a>
<i>(S) p-Terphenyl-d14</i>	90.4		53.2-156		12/05/2016 05:09	<a href="#">WG931662</a>

- 1 Cp
- 2 Tc
- 3 Ss
- 4 Cn
- 5 Sr
- 6 Qc
- 7 Gl
- 8 Al
- 9 Sc



## Wet Chemistry by Method 2320 B-2011

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Alkalinity	258000		20000	1	12/02/2016 10:37	<a href="#">WG931302</a>

## Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Nitrate	104		100	1	12/01/2016 17:21	<a href="#">WG931112</a>
Sulfate	77600		5000	1	12/01/2016 17:21	<a href="#">WG931112</a>

## Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Manganese	1140		5.00	1	12/03/2016 15:27	<a href="#">WG931185</a>
Manganese,Dissolved	1260		5.00	1	12/02/2016 13:39	<a href="#">WG931186</a>

## Semi-Volatile Organic Compounds (GC) by Method NWTPHDX

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Diesel Range Organics (DRO)	ND		250	1	12/03/2016 13:50	<a href="#">WG931238</a>
Residual Range Organics (RRO)	ND		500	1	12/03/2016 13:50	<a href="#">WG931238</a>
<i>(S) o-Terphenyl</i>	128		50.0-150		12/03/2016 13:50	<a href="#">WG931238</a>

## Semi Volatile Organic Compounds (GC/MS) by Method 8270D-SIM

Analyte	Result	Qualifier	RDL	Dilution	Analysis	Batch
	ug/l		ug/l		date / time	
Anthracene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Acenaphthene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Acenaphthylene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Benzo(a)anthracene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Benzo(a)pyrene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Benzo(b)fluoranthene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Benzo(g,h,i)perylene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Benzo(k)fluoranthene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Chrysene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Dibenz(a,h)anthracene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Fluoranthene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Fluorene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Indeno(1,2,3-cd)pyrene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Naphthalene	ND		0.250	1	12/05/2016 05:31	<a href="#">WG931662</a>
Phenanthrene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
Pyrene	ND		0.0500	1	12/05/2016 05:31	<a href="#">WG931662</a>
1-Methylnaphthalene	ND		0.250	1	12/05/2016 05:31	<a href="#">WG931662</a>
2-Methylnaphthalene	ND		0.250	1	12/05/2016 05:31	<a href="#">WG931662</a>
2-Chloronaphthalene	ND		0.250	1	12/05/2016 05:31	<a href="#">WG931662</a>
<i>(S) Nitrobenzene-d5</i>	93.1		45.1-170		12/05/2016 05:31	<a href="#">WG931662</a>
<i>(S) 2-Fluorobiphenyl</i>	99.0		57.7-153		12/05/2016 05:31	<a href="#">WG931662</a>
<i>(S) p-Terphenyl-d14</i>	98.7		53.2-156		12/05/2016 05:31	<a href="#">WG931662</a>

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Method Blank (MB)

(MB) R3182049-2 12/02/16 07:20

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
Alkalinity	U		2710	20000

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

L875488-02 Original Sample (OS) • Duplicate (DUP)

(OS) L875488-02 12/02/16 09:17 • (DUP) R3182049-7 12/02/16 09:26

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
Alkalinity	406000	384000	1	6.00		20

L875488-01 Original Sample (OS) • Duplicate (DUP)

(OS) L875488-01 12/02/16 09:34 • (DUP) R3182049-8 12/02/16 09:41

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
Alkalinity	432000	419000	1	3.00		20

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3182049-3 12/02/16 08:39 • (LCSD) R3182049-9 12/02/16 09:55

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
Alkalinity	100000	102000	103000	102	103	85.0-115			2.00	20



Method Blank (MB)

(MB) R3181971-1 12/01/16 11:35

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
Nitrate	U		22.7	100
Sulfate	U		77.4	5000

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc

L875470-01 Original Sample (OS) • Duplicate (DUP)

(OS) L875470-01 12/01/16 14:17 • (DUP) R3181971-4 12/01/16 14:32

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
Nitrate	ND	66.7	1	12	J	15
Sulfate	ND	296	1	19	J P1	15

L875474-02 Original Sample (OS) • Duplicate (DUP)

(OS) L875474-02 12/01/16 17:52 • (DUP) R3181971-6 12/01/16 19:55

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
Nitrate	1670	1650	1	1		15

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3181971-2 12/01/16 11:51 • (LCSD) R3181971-3 12/01/16 12:06

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
Nitrate	8000	7940	7940	99	99	80-120			0	15
Sulfate	40000	39000	39000	97	97	80-120			0	15

L875470-04 Original Sample (OS) • Matrix Spike (MS)

(OS) L875470-04 12/01/16 16:04 • (MS) R3181971-5 12/01/16 16:20

Analyte	Spike Amount	Original Result	MS Result	MS Rec.	Dilution	Rec. Limits	MS Qualifier
Nitrate	5000	ND	5120	101	1	80-120	
Sulfate	50000	18800	69400	101	1	80-120	



L875474-08 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L875474-08 12/01/16 20:42 • (MS) R3181971-7 12/01/16 20:57 • (MSD) R3181971-8 12/01/16 21:43

Analyte	Spike Amount ug/l	Original Result ug/l	MS Result ug/l	MSD Result ug/l	MS Rec. %	MSD Rec. %	Dilution	Rec. Limits %	MS Qualifier	MSD Qualifier	RPD %	RPD Limits %
Nitrate	5000	ND	5060	5200	101	104	1	80-120			3	15
Sulfate	50000	ND	52600	51500	105	103	1	80-120			2	15

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



Method Blank (MB)

(MB) R3182304-1 12/03/16 11:43

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
Sulfate	U		77.4	5000

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

L875736-01 Original Sample (OS) • Duplicate (DUP)

(OS) L875736-01 12/03/16 13:12 • (DUP) R3182304-4 12/03/16 13:27

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
Sulfate	5460	5180	1	5		15

L875736-10 Original Sample (OS) • Duplicate (DUP)

(OS) L875736-10 12/03/16 16:26 • (DUP) R3182304-6 12/03/16 16:41

Analyte	Original Result	DUP Result	Dilution	DUP RPD	DUP Qualifier	DUP RPD Limits
Sulfate	10200	10200	1	0		15

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3182304-2 12/03/16 11:58 • (LCSD) R3182304-3 12/03/16 12:12

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
Sulfate	40000	38100	39500	95	99	80-120			4	15

L875736-02 Original Sample (OS) • Matrix Spike (MS)

(OS) L875736-02 12/03/16 13:42 • (MS) R3182304-5 12/03/16 13:57

Analyte	Spike Amount	Original Result	MS Result	MS Rec.	Dilution	Rec. Limits	MS Qualifier
Sulfate	50000	8280	60600	105	1	80-120	

L875741-02 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L875741-02 12/03/16 17:41 • (MS) R3182304-7 12/03/16 17:55 • (MSD) R3182304-8 12/03/16 18:10

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
Sulfate	50000	8020	60400	60900	105	106	1	80-120			1	15



Method Blank (MB)

(MB) R3182334-1 12/03/16 14:30

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
Manganese	U		0.250	5.00

<sup>1</sup>Cp

<sup>2</sup>Tc

<sup>3</sup>Ss

<sup>4</sup>Cn

<sup>5</sup>Sr

<sup>6</sup>Qc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3182334-2 12/03/16 14:33 • (LCSD) R3182334-3 12/03/16 14:37

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
Manganese	50.0	48.8	51.9	98	104	80-120			6	20

L875470-04 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L875470-04 12/03/16 14:40 • (MS) R3182334-5 12/03/16 14:48 • (MSD) R3182334-6 12/03/16 14:51

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
Manganese	50.0	643	700	706	114	126	1	75-125		V	1	20

<sup>7</sup>Gl

<sup>8</sup>Al

<sup>9</sup>Sc



Method Blank (MB)

(MB) R3182109-1 12/02/16 10:31

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
Manganese,Dissolved	U		0.250	5.00

<sup>1</sup>Cp

<sup>2</sup>Tc

<sup>3</sup>Ss

<sup>4</sup>Cn

<sup>5</sup>Sr

<sup>6</sup>Qc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3182109-2 12/02/16 10:35 • (LCSD) R3182109-3 12/02/16 10:38

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
Manganese,Dissolved	50.0	49.9	50.2	100	100	80-120			1	20

L875470-02 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L875470-02 12/02/16 10:42 • (MS) R3182109-5 12/02/16 10:49 • (MSD) R3182109-6 12/02/16 10:52

Analyte	Spike Amount	Original Result	MS Result	MSD Result	MS Rec.	MSD Rec.	Dilution	Rec. Limits	MS Qualifier	MSD Qualifier	RPD	RPD Limits
Manganese,Dissolved	50.0	332	387	385	111	106	1	75-125			1	20

<sup>7</sup>Gl

<sup>8</sup>Al

<sup>9</sup>Sc



Method Blank (MB)

(MB) R3182235-1 12/02/16 15:19

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Diesel Range Organics (DRO)	U		83.3	250
Residual Range Organics (RRO)	U		167	500
<i>(S) o-Terphenyl</i>	126			64.0-146

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3182235-2 12/02/16 15:36 • (LCSD) R3182235-3 12/02/16 15:52

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	%	%	%			%	%
Diesel Range Organics (DRO)	750	937	999	125	133	50.0-150			6.43	20
Residual Range Organics (RRO)	750	828	874	110	117	50.0-150			5.35	20
<i>(S) o-Terphenyl</i>				120	128	64.0-146				

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



Method Blank (MB)

(MB) R3182439-3 12/05/16 01:53

Analyte	MB Result	MB Qualifier	MB MDL	MB RDL
	ug/l		ug/l	ug/l
Anthracene	U		0.0140	0.0500
Acenaphthene	U		0.0100	0.0500
Acenaphthylene	U		0.0120	0.0500
Benzo(a)anthracene	U		0.00410	0.0500
Benzo(a)pyrene	U		0.0116	0.0500
Benzo(b)fluoranthene	U		0.00212	0.0500
Benzo(g,h,i)perylene	U		0.00227	0.0500
Benzo(k)fluoranthene	U		0.0136	0.0500
Chrysene	U		0.0108	0.0500
Dibenz(a,h)anthracene	U		0.00396	0.0500
Fluoranthene	U		0.0157	0.0500
Fluorene	U		0.00850	0.0500
Indeno(1,2,3-cd)pyrene	U		0.0148	0.0500
Naphthalene	0.0419	J	0.0198	0.250
Phenanthrene	U		0.00820	0.0500
Pyrene	U		0.0117	0.0500
1-Methylnaphthalene	U		0.00821	0.250
2-Methylnaphthalene	U		0.00902	0.250
2-Chloronaphthalene	U		0.00647	0.250
(S) Nitrobenzene-d5	97.3			33.8-179
(S) 2-Fluorobiphenyl	103			55.5-150
(S) p-Terphenyl-d14	104			46.2-163

1 Cp

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3182439-1 12/05/16 01:10 • (LCSD) R3182439-2 12/05/16 01:31

Analyte	Spike Amount	LCS Result	LCSD Result	LCS Rec.	LCSD Rec.	Rec. Limits	LCS Qualifier	LCSD Qualifier	RPD	RPD Limits
	ug/l	ug/l	ug/l	%	%	%			%	%
Anthracene	2.00	2.48	2.33	124	117	68.9-153			6.10	20
Acenaphthene	2.00	2.19	2.05	110	103	67.7-141			6.38	20
Acenaphthylene	2.00	2.20	2.06	110	103	66.9-141			6.65	20
Benzo(a)anthracene	2.00	2.15	1.99	108	99.6	63.1-147			7.71	20
Benzo(a)pyrene	2.00	2.25	2.13	112	107	62.2-150			5.24	20
Benzo(b)fluoranthene	2.00	2.09	2.02	105	101	58.4-148			3.71	20
Benzo(g,h,i)perylene	2.00	2.30	2.20	115	110	57.4-152			4.55	20
Benzo(k)fluoranthene	2.00	2.23	2.15	112	107	60.5-154			3.94	20
Chrysene	2.00	2.36	2.24	118	112	64.8-155			5.61	20
Dibenz(a,h)anthracene	2.00	2.22	2.12	111	106	53.5-153			4.83	20
Fluoranthene	2.00	2.26	2.11	113	105	68.6-153			7.00	20



Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3182439-1 12/05/16 01:10 • (LCSD) R3182439-2 12/05/16 01:31

Analyte	Spike Amount ug/l	LCS Result ug/l	LCSD Result ug/l	LCS Rec. %	LCSD Rec. %	Rec. Limits %	<u>LCS Qualifier</u>	<u>LCSD Qualifier</u>	RPD %	RPD Limits %
Fluorene	2.00	2.13	1.98	106	98.9	67.3-141			7.37	20
Indeno(1,2,3-cd)pyrene	2.00	2.28	2.17	114	109	57.0-155			4.73	20
Naphthalene	2.00	2.44	2.28	122	114	66.7-135			6.70	20
Phenanthrene	2.00	2.09	1.97	105	98.3	64.3-143			6.33	20
Pyrene	2.00	2.38	2.21	119	111	60.2-154			7.07	20
1-Methylnaphthalene	2.00	2.44	2.27	122	114	68.3-144			6.96	20
2-Methylnaphthalene	2.00	2.30	2.14	115	107	67.6-143			7.51	20
2-Chloronaphthalene	2.00	2.27	2.12	113	106	69.7-144			6.72	20
<i>(S) Nitrobenzene-d5</i>				104	96.6	33.8-179				
<i>(S) 2-Fluorobiphenyl</i>				110	103	55.5-150				
<i>(S) p-Terphenyl-d14</i>				106	98.7	46.2-163				

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



## Abbreviations and Definitions

SDG	Sample Delivery Group.
MDL	Method Detection Limit.
RDL	Reported Detection Limit.
ND	Not detected at the Reporting Limit (or MDL where applicable).
U	Not detected at the Reporting Limit (or MDL where applicable).
RPD	Relative Percent Difference.
Original Sample	The non-spiked sample in the prep batch used to determine the Relative Percent Difference (RPD) from a quality control sample. The Original Sample may not be included within the reported SDG.
(S)	Surrogate (Surrogate Standard) - Analytes added to every blank, sample, Laboratory Control Sample/Duplicate and Matrix Spike/Duplicate; used to evaluate analytical efficiency by measuring recovery. Surrogates are not expected to be detected in all environmental media.
Rec.	Recovery.

Qualifier	Description
J	The identification of the analyte is acceptable; the reported value is an estimate.
O1	The analyte failed the method required serial dilution test and/or subsequent post-spike criteria. These failures indicate matrix interference.
P1	RPD value not applicable for sample concentrations less than 5 times the reporting limit.
V	The sample concentration is too high to evaluate accurate spike recoveries.

<sup>1</sup> Cp

<sup>2</sup> Tc

<sup>3</sup> Ss

<sup>4</sup> Cn

<sup>5</sup> Sr

<sup>6</sup> Qc

<sup>7</sup> Gl

<sup>8</sup> Al

<sup>9</sup> Sc



ESC Lab Sciences is the only environmental laboratory accredited/certified to support your work nationwide from one location. One phone call, one point of contact, one laboratory. No other lab is as accessible or prepared to handle your needs throughout the country. Our capacity and capability from our single location laboratory is comparable to the collective totals of the network laboratories in our industry. The most significant benefit to our "one location" design is the design of our laboratory campus. The model is conducive to accelerated productivity, decreasing turn-around time, and preventing cross contamination, thus protecting sample integrity. Our focus on premium quality and prompt service allows us to be **YOUR LAB OF CHOICE**.  
 \* Not all certifications held by the laboratory are applicable to the results reported in the attached report.



## State Accreditations

Alabama	40660	Nevada	TN-03-2002-34
Alaska	UST-080	New Hampshire	2975
Arizona	AZ0612	New Jersey–NELAP	TN002
Arkansas	88-0469	New Mexico	TN00003
California	01157CA	New York	11742
Colorado	TN00003	North Carolina	Env375
Connecticut	PH-0197	North Carolina <sup>1</sup>	DW21704
Florida	E87487	North Carolina <sup>2</sup>	41
Georgia	NELAP	North Dakota	R-140
Georgia <sup>1</sup>	923	Ohio–VAP	CL0069
Idaho	TN00003	Oklahoma	9915
Illinois	200008	Oregon	TN200002
Indiana	C-TN-01	Pennsylvania	68-02979
Iowa	364	Rhode Island	221
Kansas	E-10277	South Carolina	84004
Kentucky <sup>1</sup>	90010	South Dakota	n/a
Kentucky <sup>2</sup>	16	Tennessee <sup>14</sup>	2006
Louisiana	AI30792	Texas	T 104704245-07-TX
Maine	TN0002	Texas <sup>5</sup>	LAB0152
Maryland	324	Utah	6157585858
Massachusetts	M-TN003	Vermont	VT2006
Michigan	9958	Virginia	109
Minnesota	047-999-395	Washington	C1915
Mississippi	TN00003	West Virginia	233
Missouri	340	Wisconsin	9980939910
Montana	CERT0086	Wyoming	A2LA
Nebraska	NE-OS-15-05		

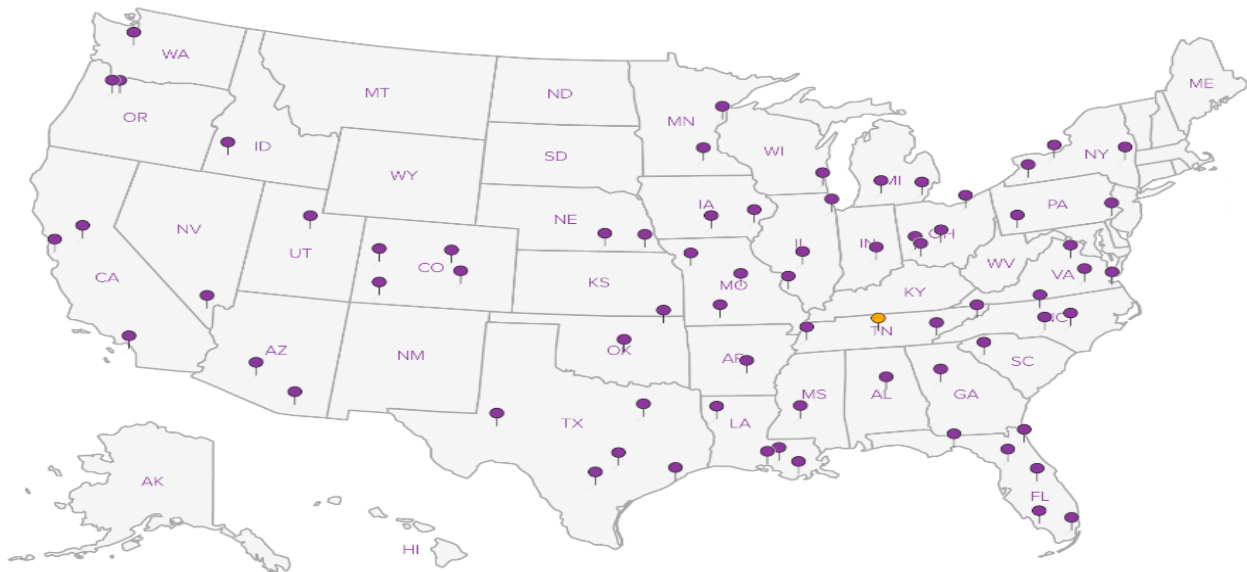
## Third Party & Federal Accreditations

A2LA – ISO 17025	1461.01	AIHA	100789
A2LA – ISO 17025 <sup>5</sup>	1461.02	DOD	1461.01
Canada	1461.01	USDA	S-67674
EPA–Crypto	TN00003		

<sup>1</sup> Drinking Water <sup>2</sup> Underground Storage Tanks <sup>3</sup> Aquatic Toxicity <sup>4</sup> Chemical/Microbiological <sup>5</sup> Mold <sup>n/a</sup> Accreditation not applicable

## Our Locations

ESC Lab Sciences has sixty-four client support centers that provide sample pickup and/or the delivery of sampling supplies. If you would like assistance from one of our support offices, please contact our main office. **ESC Lab Sciences performs all testing at our central laboratory.**



TRK: 7066 9126 3501

**Kennedy/Jenks Con-BNSF Region 1**  
 32001 32nd Avenue South, Ste 100  
 Federal Way, WA 98001

Billing Information & Quote Number:  
 Shane DeGross  
 605 Puyallup Ave S  
 Tacoma, WA 98421

Report to:  
**Nathan Moxley**

Email To: [nathanmoxley@kennedyjenks.com](mailto:nathanmoxley@kennedyjenks.com),  
[tyschreiner@kennedyjenks.com](mailto:tyschreiner@kennedyjenks.com)

Project Description: **BNSF - D St. Wells**

City/State Collected: **Tacoma, WA**

Phone: **253-835-6400**  
 Fax:

Client Project #  
**169603900**

Lab Project #  
**BNSF1KEN-D ST**

Collected by (print):  
**Alexander Leshner**

Site/Facility ID #  
**D Street Pipeline**

P.O. #

Collected by (signature):  
*Alexander Leshner*

**Rush?** (Lab MUST Be Notified)  
 \_\_\_ Same Day .....200%  
 \_\_\_ Next Day .....100%  
 \_\_\_ Two Day .....50%  
 \_\_\_ Three Day .....25%

Date Results Needed  
 Email? \_\_\_ No  Yes  
 FAX? \_\_\_ No \_\_\_ Yes

Packed on Ice N \_\_\_ Y

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs
MW-9-20161130	G	GW		11-30-16	0855	7
MW-8-20161130		GW			1000	7
MW-11-20161130		GW			1040	7
MW-5-20161130		GW			1115	7
MW-4-20161130		GW			1155	7
MW-3-20161130		GW			1230	7
MW-2-20161130		GW			1305	7
TB-1		GW			-	7
D-1-20161130	G	GW			-	7
		GW				7

Analysis / Container / Preservative										
**ALK, NO3, SO4**	125ml	HDPE-NoPres								
Disolved	250ml	HDPE-HNO3	Manganese							
NWTPHDXLVI-	w/	SGT	40ml	Amb-HCl-BT						
PAHSIMLVID	40ml	Amb-NoPres-WT								
Total Pb	250ml	HDPE-HNO3								
Total Mn										

Chain of Custody Page 1 of 1



YOUR LAB OF CHOICE

12055 Lebanon Rd  
 Mount Juliet, TN 37122  
 Phone: 615-758-5858  
 Phone: 800-767-5859  
 Fax: 615-758-5859



L# **L875V70**  
**D060**

Acctnum: **BNSF1KEN**  
 Template: **T117672**  
 Prelogin: **P576038**  
 TSR: **134 - Mark W. Beasley**  
 PB: **11-9-16**

Shipped Via: **FedEX Ground**

\* Matrix: SS - Soil GW - Groundwater WW - WasteWater DW - Drinking Water OT - Other

Remarks: **Manganese not Lead**

pH \_\_\_\_\_ Temp \_\_\_\_\_  
 Flow \_\_\_\_\_ Other \_\_\_\_\_

Relinquished by: (Signature) <i>Alexander Leshner</i>	Date: <b>11-30-16</b>	Time: <b>1500</b>	Received by: (Signature) <i>[Signature]</i>	Samples returned via: <input type="checkbox"/> UPS <input type="checkbox"/> FedEx <input type="checkbox"/> Courier <input type="checkbox"/> _____	Condition: (lab use only) <b>Jun 7</b>
Relinquished by: (Signature)	Date:	Time:	Received by: (Signature)	Temp: <b>3.4</b> °C Bottles Received: <b>57</b>	COC Seal Intact: ___ Y ___ N ___ NA
Relinquished by: (Signature)	Date:	Time:	Received by: (Signature) <i>[Signature]</i>	Date: <b>12/01/16</b> Time: <b>9:00</b>	pH Checked: NCF:

Rem./Contaminant	Sample # (lab only)
	-01
	02
	03
	04
	05
	06
	07
	08
	09



### Cooler Receipt Form

Client: <i>BNSF IKEN</i>	SDG#	<i>1875470</i>		
Cooler Received/Opened On: <i>12/01/16</i>	Temperature Upon Receipt:	<i>3.4 °C</i>		
Received By: Timiesha Scott				
Signature: <i>[Handwritten Signature]</i>				
Receipt Check List		Yes	No	N/A
Were custody seals on outside of cooler and intact?		<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were custody papers properly filled out?		<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Did all bottles arrive in good condition?		<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were correct bottles used for the analyses requested?		<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Was sufficient amount of sample sent in each bottle?		<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
Were all applicable sample containers correctly preserved and checked for preservation? (Any not in accepted range noted on COC)		<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>
If applicable, was an observable VOA headspace present?		<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Non Conformance Generated. (If yes see attached NCF)		<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>

# Attachment B

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## Data Quality Assessment Summary

**DATA VALIDATION SUMMARY – L875470  
NOVEMBER 2016 SAMPLING EVENT  
BNSF D STREET**

Laboratory Reports included in Data Validation	Report Date	Sample IDs
ESC L875470 Includes: NWTPH-Dx, PAHs, Alkalinity, Nitrate, Sulfate, Manganese	9 December 2016	Aqueous Samples: MW-9-20161130, MW-8-20161130, MW-11-20161130, MW-5-20161130, MW-4-20161130, MW-3-20161130, MW-2-20161130  Field Duplicate: D-1-20161130 collected at MW-2.

Criteria	(Yes or No)	Comment
<u>Chain-of-Custody</u> – Chain-of-custody protocol followed?	Yes	
<u>Temperature Blank</u> – Sample temperature criteria met?	Yes	Samples arrived at a temperature of 3.4 degrees Celsius (°C), which was within the recommended temperature of 4°C ± 2°.
<u>Holding times</u> – Samples analyzed within specified holding time?	Yes	
<u>Laboratory method blank samples</u> – Analytes present in method blank samples?	Yes	See Note 1 below.
<u>Field/Equipment blank samples</u> – Analytes present in field/equipment blank samples?	Not applicable	Field/equipment blank samples were not necessary and were not submitted with the batch of samples.
<u>Trip blank samples</u> – Analytes present in trip blank samples?	Not applicable	Trip blank samples were not submitted with the batch of samples.
<u>Matrix spikes/matrix spike duplicate samples</u> – Control limits met?	No	See Note 2 below.
<u>Surrogate percent recoveries</u> – Control limits met?	Yes	
<u>Laboratory control sample</u> – Control limits met?	Yes	
<u>Laboratory duplicate samples (if applicable)</u> – Control limits met?	No	See Note 3 below.
<u>Field duplicate samples (if submitted)</u> – Relative percent differences within control limits?	Yes	Primary and field duplicate samples did not contain detectable concentrations of cPAHs or NWTPH-Dx. Results for alkalinity, nitrate, sulfate, and manganese were within 20% RPD.

**NOTES**

- Naphthalene was detected at 0.0419 J micrograms per liter (µg/L) in the method blank for batch R3182439. Because naphthalene was not detected above the laboratory method reporting limit (0.05 µg/L) in the field samples, no action was taken.
- Percent recovery of manganese in the MSD for batch R3182334 was slightly higher than the laboratory control limit. Because the percent recovery was within the laboratory control limit for the MS and LCS/LCSD pair, no action was taken.
- Laboratory primary/duplicate samples for sulfate (19%) were greater than the 15% RPD limit. However, the reported concentration of 296 µg/L in the laboratory duplicate sample was below the laboratory reporting limit for the primary field sample (5,000 µg/L). Because the primary and laboratory duplicate sample results were below the 5,000 µg/L reporting limit, no action was taken.

**DATA VALIDATION SUMMARY – L875470  
NOVEMBER 2016 SAMPLING EVENT  
BNSF D STREET**

**SUMMARY**

Laboratory report L875470 was reviewed and compared to data quality objectives presented in the *Compliance Monitoring Plan, Remedial Excavation, BNSF Oil Pipeline Site, Tacoma, Washington*, dated 9 July 2008, prepared by GeoEngineers. Overall, the findings with respect to the quality assurance/quality control (QA/QC) data do not adversely affect the use of the analytical results.