#### Geotechnical Engineering

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**Materials Testing** 

**Building Science** 

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# **Phase II – Environmental Site Assessment**

78, 84, 86, 90 Beechwood Avenue & 69, 73, 77, 83, 85, 89, 93 Barrette Street Ottawa, Ontario

# **Prepared For**

Minto Communities

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# EXECUTIVE SUMMARY

# Assessment

A Phase II ESA was conducted for the properties addressed 78-90 Beechwood Avenue and 69-93 Barrette Street, in the City of Ottawa, Ontario. The purpose of the Phase II ESA was to address the potentially contaminating activities (PCAs) that were identified during the Phase I ESA and were considered to result in areas of potential environmental concern (APECs) on the subject site.

The subsurface investigation consisted of drilling fifteen (15) boreholes (BH1-BH14 and BH8B) on May 23-24, and July 15, 2019, as well as on August 24, 2020. Upon completion, ten (10) of the boreholes were instrumented with groundwater monitoring wells (BH2, BH3, BH4, BH7, BH8 BH8B, BH9, BH10, BH11, and BH13). The boreholes were drilled to depths ranging from approximately 3.25 m to 9.14 m below ground surface and terminated within weathered black shale bedrock.

Nineteen (19) soil samples were submitted for laboratory analysis of either: BTEX, PHCs F<sub>1</sub>-F<sub>4</sub>, VOCs, PAHs, metals, and/or pH parameters. According to the analytical results, the concentrations of some PHC, VOC, PAH, and/or metal parameters in the soil samples tested from BH1, BH3, BH7, BH8, BH10, and BH11 were in excess of the selected MECP Table 3 residential standards.

Twelve (12) groundwater samples were submitted for laboratory analysis of either: BTEX, PHCs  $F_1$ - $F_4$ , VOCs, and/or PAHs parameters. According to the analytical results, all detected parameter concentrations in the groundwater samples analyzed comply with the selected MECP Table 3 residential standards.

# Recommendations

Soil

Based on the findings of this assessment, PHC, VOC, PAH, and/or metal impacted soil was identified in the vicinity of BH1, BH3, BH7, BH8, BH10, and BH11, requiring some remedial work. It is our understanding that the subject site is to be redeveloped with a multi-storey residential high-rise building in the near future. Therefore, it is our recommendation that an environmental site remediation program be completed in conjunction with site redevelopment. This will require the segregation of clean soils from impacted soils, the latter of which will require disposal at an approved waste disposal facility.

It should be noted that several PHC, VOC, PAH, and metal parameters were detected in BH1, BH2, BH3, BH7 BH9, BH11, and/or BH14 at concentrations exceeding the MECP Table 1 background standards. These exceedances are not considered to pose an environmental concern to the subject site, however, if the soil is to be removed from the property for construction purposes, it should be classified as contaminated and disposed of at an approved waste disposal site.

Prior to off-site disposal at a licensed landfill, a leachate analysis of a representative sample of contaminated soil must be conducted in accordance with Ontario Regulation 347/558.

It is recommended that Paterson personnel be present on-site during remediation activities to direct the excavation and segregation of impacted soil, as well as to conduct confirmatory sampling as required.

#### Monitoring Wells

If the groundwater monitoring wells installed on-site are not going to be used in the future, or will be destroyed during future redevelopment activities, then they must be decommissioned according to Ontario Regulation Reg. 903 (Ontario Water Resources Act). The monitoring wells will be registered with the MECP under this regulation. Further information can be provided upon request in this regard.

# **1.0 INTRODUCTION**

At the request of Minto Communities, Paterson Group (Paterson) conducted a Phase II – Environmental Site Assessment for the properties addressed 78-90 Beechwood Avenue and 69-93 Barrette Street, in the City of Ottawa, Ontario. The purpose of this investigation was to assess the areas of potential environmental concern identified in the Phase I ESA report.

# 1.1 Site Description

Address:	78, 84, 86, 90 Beechwood Avenue, Ottawa, ON 69, 73, 77, 83, 85, 89, 93 Barrette Street Ottawa, ON.
Legal Description:	Part of Lot 4, Junction Gore Concession, Rideau Front, Formerly the Township of Gloucester, in the City of Ottawa, Ontario.
Location:	The subject site is situated between the south side of Beechwood Avenue and the north side of Barrette Street, approximately 40 m west of St. Charles Street, in the City of Ottawa, Ontario. Refer to Figure 1 - Key Plan, appended to this report.
Latitude and Longitude:	45° 26' 25" N, 75° 40' 27" W
Configuration:	Rectangular
Site Area:	4,170 m <sup>2</sup> (approximate)
Zoning:	TM – Traditional Mainstreet Zone R4T – Residential Fourth Density Zone
Current Use:	The subject site is currently occupied by twelve (12) buildings of mixed commercial and residential uses.
Services:	The subject site is located in a municipally serviced area.

# **1.2 Property Ownership**

The subject property is currently owned by Minto Communities. Paterson was engaged to conduct this Phase II ESA by Mr. Kevin Harper of Minto Communities. Minto's offices are located at 180 Kent Street, Suite 200, Ottawa, Ontario. Mr. Harper can be reached at 613-751-2857.

# **1.3 Current and Proposed Future Uses**

The subject site is currently occupied by twelve (12) buildings of mixed commercial and residential uses.

It is our understanding that the subject property is to be redeveloped with a multistorey residential building, with ground floor commercial units and underground parking.

# **1.4 Applicable Site Condition Standard**

The site condition standards for the subject property were obtained from Table 3 of the document entitled, *"Soil, Ground Water and Sediment Standards for Use Under Part XV.1 of the Environmental Protection Act"*, prepared by the Ministry of the Environment, Conservation and Parks (MECP), and dated April 15, 2011. The selected MECP standards are based on the following considerations:

- □ Coarse-grained soil conditions;
- □ Non-potable groundwater conditions;
- **Residential land use.**

The residential standards were selected based on the future land use of the subject site. Grain size analysis was not conducted as part of this assessment, however, the coarse-grained soil standards were chosen as a conservative approach.

The MECP Table 1 standards for Full Depth Background Site Conditions were also selected for additional consideration in order to assess the on-site soil conditions prior to future off-site disposal.

# 2.0 BACKGROUND INFORMATION

# 2.1 Physical Setting

The subject site is occupied by the twelve (12) aforementioned buildings which occupy the majority of the site area. The remainder of the site consists of asphaltic concrete laneways or gravel parking areas. The subject site is considered to be at grade with respect to the neighbouring properties and the adjacent streets.

The site topography appears to be relatively flat, whereas the regional topography appears to slope down to the south, in the general direction of the Rideau River.

Water drainage on the subject site occurs primarily via infiltration in the gravel and landscaped areas, as well as via surface run-off towards catch basins located on the adjacent streets.

# 2.2 Past Investigations

The following report was reviewed prior to conducting this assessment:

"Phase I - Environmental Site Assessment, 92 Beechwood Avenue, City of Ottawa, Ontario", prepared by Kollaard Associates and dated June 2009.

This Phase I ESA report discusses several landfill sites as posing a potential source of contamination with respect to the subject site. The report also makes reference to a geotechnical investigation completed by Fondex Ontario Limited in 2004 in which debris consisting of brick, ceramic, wood, plastic bags, cans, glass and asphalt was encountered during the borehole drilling program. The debris was considered to be representative of the types of materials described in the aforementioned Landfill Management Strategy as having been interred.

The report recommended that a soil and groundwater sampling program be carried out.

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# 3.0 SCOPE OF INVESTIGATION

### 3.1 Overview of Site Investigations

The subsurface investigation for this assessment was conducted on May 23-24, and July 15, 2019, as well as on August 24, 2020. The field program consisted of drilling fifteen (15) boreholes (BH1-BH14 and BH8B) throughout the subject site, of which ten (10) were instrumented with groundwater monitoring wells (BH2, BH3, BH4, BH7, BH8 BH8B, BH9, BH10, BH11, and BH13). The boreholes were drilled to depths ranging from approximately 3.25 m to 9.14 m below ground surface and terminated within weathered black shale bedrock.

#### 3.2 Media Investigated

During the subsurface investigation, soil and groundwater samples were obtained and submitted for laboratory analysis. The rationale for sampling and analyzing these media is based on the contaminants of potential concern identified in the Phase I ESA.

The contaminants of potential concern for the soil and groundwater on the subject site are considered to be:

- □ Volatile Organic Compounds (VOCs);
- Benzene, Toluene, Ethylbenzene, and Xylenes (BTEX);
- Petroleum Hydrocarbons, fractions 1 4 (PHCs F<sub>1</sub>-F<sub>4</sub>);
- □ Metals (including Mercury and Hexavalent Chromium);
- D Polycyclic Aromatic Hydrocarbons (PAHs).

# 3.3 Phase I Conceptual Site Model

#### **Existing Buildings and Structures**

The subject site is currently occupied by twelve (12) buildings of mixed residential and commercial uses.

#### Drinking Water Wells

Based on the availability of municipal services, no drinking water wells are expected to be present within the Phase I study area.

#### Geological and Hydrogeological Setting

Based on the available information, the bedrock in the area of the subject site consists of shale of the Billings Formation, whereas the surficial geology consists of fine-textured glaciomarine deposits, with an overburden thickness of approximately 3 m.

The site topography appears to be relatively flat, whereas the regional topography appears to slope down to the south, in the general direction of the Rideau River. Based on the regional topography, the groundwater in the area of the subject site is interpreted to be flowing in a southerly direction.

#### Neighbouring Land Use

Neighbouring land use in the Phase I study area consists mainly of residential and commercial properties.

#### Water Bodies and Areas of Natural and Scientific Interest

There are no waterbodies or areas of natural and scientific interest located on the subject site or within the Phase I study area. The nearest named waterbody with respect to the subject site is the Rideau River, approximately 315 m to the south.

# Potentially Contaminating Activities and Areas of Potential Environmental Concern

As per Section 7.1 of the Phase I ESA report, four (4) PCAs were determined to result in APECs with respect to the subject site. These PCAs include:

- □ A former autobody shop and service garage, located at 89 Barrette Street and situated in the central portion of the subject site;
- □ The presence of a former landfill site, located throughout the Phase I study area and situated beneath the entirety of the subject site;
- □ A dry cleaners, located at the property addressed 110 Beechwood Avenue and situated approximately 20 m to the east of the subject site;
- □ A former retail fuel outlet, located at the property addressed 64 Beechwood Avenue and situated immediately to the west of the subject site.

Other off-site PCAs were identified within the Phase I study area but were deemed not to be of concern based on their separation distances as well as their down-gradient or cross-gradient orientations

#### **Contaminants of Potential Concern**

The contaminants of potential concern (CPCs) associated with the subject site are considered to be:

- □ Volatile Organic Compounds (VOCs);
- Benzene, Toluene, Ethylbenzene, and Xylenes (BTEX);
- D Petroleum Hydrocarbons, fractions 1 4 (PHCs F<sub>1</sub>-F<sub>4</sub>);
- □ Metals (including Mercury and Hexavalent Chromium);
- D Polycyclic Aromatic Hydrocarbons (PAHs).

These contaminants have the potential to be present in the soil/fill matrix as well as the groundwater on the subject site.

#### Assessment of Uncertainty and/or Absence of Information

The information available for review as part of the preparation of the Phase I ESA is considered to be sufficient to conclude that there are PCAs and APECs associated with the subject site. The presence of these PCAs were confirmed by a variety of independent sources, and as such, the conclusions of this report are not affected by uncertainty which may be present with respect to the individual sources.

#### 3.4 Deviations from Sampling and Analysis Plan

The Sampling and Analysis Plan for this project is included in Appendix 1 of this report. There were no deviations made from the Sampling and Analysis Plan.

#### 3.5 Impediments

No physical impediments or denial of access was encountered during this Phase II – Environmental Site Assessment.

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# 4.0 INVESTIGATION METHOD

# 4.1 Subsurface Investigation

The subsurface investigation for this assessment was conducted on May 23-24, and July 15, 2019, as well as on August 24, 2020. The field program consisted of drilling fifteen (15) boreholes (BH1-BH14 and BH8B) throughout the subject site, of which ten (10) were instrumented with groundwater monitoring wells (BH2, BH3, BH4, BH7, BH8 BH8B, BH9, BH10, BH11, and BH13). The boreholes were drilled to depths ranging from approximately 3.25 m to 9.14 m below ground surface and terminated within an underlying layer of weathered black shale bedrock.

Under the full-time supervision of Paterson personnel, the boreholes were drilled using a low-clearance drill rig provided by George Downing Estate Drilling of Hawkesbury, Ontario. The locations of the boreholes are illustrated on Drawing PE4614-4 – Test Hole Location Plan, appended to this report

# 4.2 Soil Sampling

Soil sampling protocols were followed using the MECP document entitled, "Guidance on Sampling and Analytical Methods for Use at Contaminated Sites in Ontario", dated May 1996. The samples were recovered using a stainless-steel split spoon while wearing protective gloves (changed after each sample), and immediately placed into plastic bags. If significant contamination was encountered, the samples were instead placed into glass jars. Sampling equipment was routinely washed in soapy water and rinsed with methylhydrate after each split spoon to prevent any cross contamination of the samples. The samples were also stored in coolers to reduce analyte volatilization during transportation.

One hundred (100) soil samples obtained as part of the current and previous subsurface investigations were recovered from the boreholes by means of auger and split spoon sampling, with samples taken at approximate 0.76 m intervals. The depths at which split spoon, auger flight, and rock core samples were obtained from the boreholes are shown as "SS", "AU", and "RC" on the Soil Profile and Test Data Sheets in Appendix 1.

The soil profile at the subject site generally consisted of gravel or asphaltic concrete over gravel, followed by shaley till or fill material over shaley till. The till graded into a weathered shale bedrock deposit at various depths. Specific details of the soil profile at each test hole location are presented on the Soil Profile and Test Data sheets appended to this report.

# 4.3 Field Screening Measurements

All soil samples collected were subjected to a preliminary screening procedure, which included a visual screening for colour and evidence of metals, as well as soil vapour screening with a Photo Ionization Detector. The soil vapours were measured by inserting the analyzer probe into the nominal headspace above the soil sample. The samples were then agitated/manipulated gently as the measurements were taken, and the peak reading registered within the first 15 seconds was recorded as the vapour measurement.

Samples with the highest vapour readings within a given borehole were generally selected as candidates for laboratory analysis. Additional samples were selected from different stratigraphic units to attempt to delineate the vertical extent of the contamination within a given borehole.

The measured vapour readings are depicted on the Soil Profile and Test Data Sheets in Appendix 1

#### 4.4 Groundwater Monitoring Well Installation

Ten (10) groundwater monitoring wells were installed on the subject site as part of this Phase II ESA investigation. These monitoring wells were constructed using 50 mm diameter Schedule 40 threaded PVC risers and screens. A sand pack consisting of silica sand was placed around the screen and a bentonite seal was placed above the screen to minimize cross-contamination. A summary of the monitoring well construction details are listed below in Table 1 as well as on the Soil Profile and Test Data Sheets provided in Appendix 1.

Upon completion, the groundwater monitoring wells were developed using a dedicated inertial lift pump, with a minimum of three (3) well volumes being removed from the wells at the time of installation. The wells were developed until the appearance of the water was noted to be stabilized. In addition, the ground surface elevations of each borehole were subsequently surveyed with respect to a known geodetic elevation.

	ig Well Cor	ISTRUCTION	Details			
Well ID	Ground Surface Elevation	Total Depth (m BGS)	Screened Interval (m BGS)	Sand Pack (m BGS)	Bentonite Seal (m BGS)	Casing Type
BH2	57.94	6.86	5.36 – 6.86	4.90 - 6.86	0.00 - 4.90	Flushmour
BH3	58.66	9.14	7.64 – 9.14	7.32 – 9.14	6.10 – 7.32	Flushmour
BH4	58.43	5.94	2.94 – 5.94	2.40 – 5.94	0.00 - 2.40	Flushmour
BH7	58.59	6.10	3.97 – 6.10	2.44 – 6.10	0.15 – 2.44	Flushmour
BH8	58.67	6.70	3.81 – 6.10	3.51 – 6.10	0.15 – 3.51	Flushmour
BH8B	58.67	8.84	7.34 – 8.84	6.71 – 8.84	0.15 – 6.71	Flushmour
BH9	58.71	6.86	3.71 – 6.71	3.35 – 6.71	0.15 – 3.35	Flushmour
BH10	58.66	6.70	3.81 – 6.10	3.35 – 6.71	0.15 – 3.35	Flushmour
BH11	58.11	5.84	2.84 – 5.84	2.52 – 5.84	0.23 – 2.52	Flushmour
BH13	58.73	5.89	2.89 - 5.89	2.46 - 5.89	0.10 – 2.46	Flushmour

# 4.5 Groundwater Sampling

Groundwater sampling protocols were followed using the MECP document entitled, "Guidance on Sampling and Analytical Methods for Use at Contaminated Sites in Ontario", dated May 1996. Groundwater samples were obtained from each monitoring well using dedicated sampling equipment. Standing water was purged from each well prior to sampling. Samples were stored in coolers to reduce analyte volatilization during transportation. Details of our standard operating procedure for groundwater sampling are provided in the Sampling and Analysis Plan in Appendix 1.

#### 4.6 Residue Management

Soil cuttings, purge water, and equipment cleaning fluids were retained on-site.

# 4.7 Analytical Testing

The following soil and groundwater samples were submitted for laboratory analysis as part of this Phase II ESA:

Table 2								
Testing	Parameters	for S	Subn	nitte	d So	il Sa	Impl	es
	Sample Depth		Parar	neter	s Ana	lyzed		
Sample ID	& Stratigraphic Unit	втех	PHCs	VOCs	PAHs	Metals <sup>1</sup>	Ηd	Rationale
BH1-SS2	0.76 – 1.37 m Fill Material	х	х		х	х	Х	Assess for potential impacts resulting from a former landfill site.
BH2-SS3	1.52 – 2.13 m Clayey Silt Till	х	х				х	Assess for potential impacts resulting from a former off-site retail fuel outlet.
BH3-AU1	0.00 – 0.61 m Fill Material	х				х		Assess for potential impacts resulting from a former landfill site.
BH3- SS9/10	6.10 – 7.47 m Silty Sand Till			х				Assess for potential impacts resulting from an off-site dry-cleaners.
BH7-AU1	0.00 – 0.61 m Fill Material					х		Assess for potential impacts resulting from a former landfill site.
BH7-SS3	1.52 – 2.13 m Shale Bedrock			х				Assess for potential impacts resulting from an off-site dry-cleaners.
BH8-SS4	2.29 – 2.90 m Silty Sand Till		х	х				Assess for potential impacts resulting from an off-site dry-cleaners
BH8-SS9	6.10 – 6.71 m Shale Bedrock			х				Assess for potential impacts resulting from an off-site dry-cleaners.
BH9-AU1	0.00 – 0.61 m Fill Material		х					Assess for potential impacts resulting from a former landfill site.
BH9-SS6	3.81 – 4.42 m Silty Sand Till			х				Assess for potential impacts resulting from an off-site dry-cleaners.
BH9-SS9	6.10 – 6.71 m Silty Sand Till			х				Assess for potential impacts resulting from an off-site dry-cleaners.
BH10- SS2	0.76 – 1.37 m Fill Material				х	х		Assess for potential impacts resulting from a former autobody shop.
BH11- AU1	0.00 – 0.61 m Fill Material		х					Assess for potential impacts resulting from a former landfill site.
BH11- SS2	0.76 – 1.37 m Silty Sand Till				х	х		Assess for potential impacts resulting from a former landfill site.
BH11- SS4	2.29 – 2.90 m Silty Sand Till	х	х					Assess for potential impacts resulting from a former off-site retail fuel outlet.
BH13- AU1	0.00 – 0.61 m Fill Material		х					Assess for potential impacts resulting from a former autobody shop.
BH13- SS3	1.52 – 2.13 m Silty Sand Till		х			х		Assess for potential impacts resulting from a former autobody shop.
BH13- SS7	4.57 – 4.83 m Shale Bedrock			х				Assess for potential impacts resulting from an off-site dry-cleaners.
BH14- SS2	0.76 – 1.37 m Fill Material					х		Assess for potential impacts resulting from a former landfill site.

1 – Including Mercury and Chromium VI

Table 3						
Testing	Parameters 1	for Su	bmitte	d Soil	Samp	les
		Pa	rameter	s Analyz	ed .	
Sample ID	Sample Depth & Stratigraphic Unit	ВТЕХ	PHCs	VOCs	PAHs	Rationale
BH2- GW1	5.36 – 6.86 m Shale Bedrock	х	Х	Х	х	Assess for potential impacts resulting from a former off-site retail fuel outlet.
BH3- GW1	7.64 – 9.14 m Shale Bedrock	Х	х	Х	х	Assess for potential impacts resulting from an off-site dry-cleaners.
BH3- GW2	7.64 – 9.14 m Shale Bedrock			Х		Assess for potential impacts resulting from an off-site dry-cleaners.
BH3- GW3	7.64 – 9.14 m Shale Bedrock			Х		Assess for potential impacts resulting from an off-site dry-cleaners.
BH4- GW1	2.94 – 5.94 m Shale Bedrock		Х	Х	х	Assess for potential impacts resulting from a former landfill site.
BH7- GW1	3.97 – 6.10 m Shale Bedrock		Х	Х		Assess for potential impacts resulting from a former landfill site.
BH8B- GW1	7.34 – 8.84 m Shale Bedrock			х		Assess for potential impacts resulting from an off-site dry-cleaners.
BH8B- GW2	7.34 – 8.84 m Shale Bedrock			Х		Assess for potential impacts resulting from an off-site dry-cleaners.
BH9- GW1	3.71 – 6.71 m Silty Sand Till			Х		Assess for potential impacts resulting from an off-site dry-cleaners.
BH10- GW1	3.81 – 6.10 m Shale Bedrock		х	Х		Assess for potential impacts resulting from a former autobody shop.
BH11- GW1	2.84 – 5.84 m Shale Bedrock		Х	Х		Assess for potential impacts resulting from a former off-site retail fuel outlet.
BH13- GW1	2.89 – 5.89 m Shale Bedrock		х	х		Assess for potential impacts resulting from a former autobody shop.
1 – Including	g Mercury and Chrom	nium VI				

Paracel Laboratories (Paracel), of Ottawa, Ontario, performed the laboratory analysis on the samples submitted for analytical testing. Paracel is a member of the Standards Council of Canada/Canadian Association for Environmental Analytical Laboratories (SCC/CAEAL). Paracel is accredited and certified by SCC/CAEAL for specific tests registered with the association.

# 4.8 Elevation Surveying

Monitoring well locations were surveyed using a laser level device, with the elevations surveyed relative to a magnetic nail, located on the southern sidewalk of Beechwood Avenue, which contains a known geodetic elevation. The elevations of the monitoring wells are shown on Drawing PE4614-4 – Test Hole Location Plan.

# 4.9 Quality Assurance and Quality Control Measures

A summary of quality assurance and quality control (QA/QC) measures, including sampling containers, preservation, labelling, handling, and custody, equipment cleaning procedures, and field quality control measurements is provided in the Sampling and Analysis Plan in Appendix 1.

# 5.0 REVIEW AND EVALUATION

# 5.1 Geology

Site geology details are provided in the Soil Profile and Test Data Sheets provided in Appendix 1. Site soils generally consist of fill material, consisting of either crushed stone and/or brown silty sand and gravel, underlain by either brown clayey silt till or brown silty sand till. Bedrock, consisting of weathered shale, was generally encountered at depths ranging from approximately 1.52 m to 6.86 m below the existing ground surface.

# 5.2 Groundwater Elevations, Flow Direction, and Hydraulic Gradient

Groundwater levels were measured at the subject site during the September 3, 2020 sampling event, using an electronic water level meter. It should be noted that BH7 and BH9 could not be located at the time of the sampling event. The groundwater levels are summarized below in Table 4.

Table 4 Groundwa	ater Level Meas	urements		
Borehole Location	Ground Surface Elevation	Water Level Depth (m below grade)	Water Level Elevation (m ASL)	Date of Measurement
BH2	57.94	3.61	54.33	
BH3	58.66	5.79	52.87	
BH4	58.43	4.06	54.37	
BH7	58.59	Could Not Locate	-	
BH8	58.67	6.34	52.33	O antanah an O 0000
BH8B	58.67	6.37	52.30	September 3, 2020
BH9	58.71	Could Not Locate	-	1
BH10	58.66	4.72	53.94	1
BH11	58.11	4.28	53.83	1
BH13	58.73	3.20	55.53	1

The groundwater on-site was typically encountered within the underlying shale bedrock unit at depths ranging from approximately 3.20 m to 6.37 m below ground surface. No unusual visual or olfactory observations were noted in the groundwater samples obtained from the boreholes.

Using the groundwater elevations recorded during the September 3, 2020 sampling event, groundwater contour mapping was completed as part of this assessment. According to the mapped contour data, illustrated on Drawing PE4614-4 Test Hole Location Plan in the appendix, the groundwater flow on the subject site is interpreted to be in a northerly direction. A horizontal hydraulic gradient of approximately 0.07 m/m was also calculated as part of this assessment.

It should be noted that groundwater levels fluctuate throughout the year with seasonal variations.

### 5.3 Field Screening

Field screening of the soil samples collected during the drilling programs resulted in organic vapour readings ranging from 0.0 ppm to 51.8 ppm. These vapour readings indicate that there is a negligible potential for the presence of volatile substances. Heavier substances, such as heavy oil, may not be detected by the above field screening method, however, visual observations did not suggest the presence of hydrocarbon contamination. Field screening results of each individual soil sample are provided on the Soil Profile and Test Data Sheets appended to this report.

# 5.4 Fine/Coarse Soil Texture

Grain size analysis was not completed as part of this investigation. Coarse grained soil standards were chosen based on the nature of the recovered soil samples.

# 5.5 Soil Quality

As part of this assessment, nineteen (19) soil samples were submitted for laboratory analysis of either: BTEX, PHCs F<sub>1</sub>-F<sub>4</sub>, VOCs, PAHs, Metals, and/or pH parameters. The results of the analytical testing are presented below in Tables 5 to 9. The laboratory certificates of analysis are provided in Appendix 1.

Analytical BTEX & PH							
			Soil Samp	oles (µg/g)		MECP	MECP
			May 2	3, 2019		Table 1	Table 3
Parameter	MDL (µg/g)	BH1- SS2	BH2- SS3	BH3- AU1	BH3- SS9/10	Residential Soil Standards (µg/g)	Residential Soil Standards (µg/g)
Benzene	0.05	nd	nd	nd	nd	0.02	0.21
Ethylbenzene	0.05	nd	nd	nd	nd	0.05	2
Toluene	0.05	0.05	nd	nd	nd	0.2	2.3
Xylenes	0.05	0.73	nd	nd	0.29	0.05	3.1
PHCs F <sub>1</sub>	7	44	nd	15	nt	25	55
PHCs F <sub>2</sub>	4	22	80	1,040	nt	10	98
PHCs F <sub>3</sub>	8	42	29	2,370	nt	240	300
PHCs F <sub>4</sub>	6	12	nd	2,590	nt	120	2,800
PHCs F <sub>4G</sub>	50	nt	nt	9,490	nt	120	2,800

nt – not tested for this parameter

□ nd – not detected above the MDL

<u>Underlined</u> – Value exceeds MECP Table 1 standards <u>Bold and Underlined</u> – value exceeds selected MECP standards

Table 5 (continued)
Analytical Test Results - Soil
BTEX & PHCs (F1-F4)

			So	oil Samp July 1	MECP Table 1	MECP Table 3			
Parameter	MDL (µg/g)	BH7- SS3	BH8- SS4	BH8- SS9	BH9- AU1	BH9- SS6	BH9- SS9	Residential Soil Standards (µg/g)	Residential Soil Standards (µg/g)
Benzene	0.05	nd	nd	nd	nt	nd	nd	0.02	0.21
Ethylbenzene	0.05	nd	nd	nd	nt	nd	nd	0.05	2
Toluene	0.05	nd	nd	nd	nt	nd	nd	0.2	2.3
Xylenes	0.05	nd	nd	nd	nt	0.87	nd	0.05	3.1
PHCs F1	7	nt	<u>60</u>	nt	nt	nt	nt	25	55
PHCs F <sub>2</sub>	4	nt	2,980	nt	nd	nt	nt	10	98
PHCs F <sub>3</sub>	8	nt	888	nt	98	nt	nt	240	300
PHCs F <sub>4</sub>	6	nt	nd	nt	171	nt	nt	120	2,800
PHCs F <sub>4G</sub>	50	nt	nt	nt	nt	nt	nt	120	2,800
<ul> <li>nt – not</li> <li>nd – not</li> </ul>	Method Dete tested for th detected at <u>ted</u> – Value	is parame pove the I	eter MDL	able 1 sta	ndards				

Bold and Underlined – value exceeds selected MECP standards

			Soil Samp	MECP	MECP		
			August	Table 1	Table 3		
Parameter	MDL (µg/g)	BH11- AU1	BH11- SS4	BH13- AU1	BH13- SS3	Residential Soil Standards (µg/g)	Residential Soil Standards (µg/g)
Benzene	0.05	nt	nd	nt	nt	0.02	0.21
Ethylbenzene	0.05	nt	nd	nt	nt	0.05	2
Toluene	0.05	nt	nd	nt	nt	0.2	2.3
Xylenes	0.05	nt	nd	nt	nt	0.05	3.1
PHCs F1	7	nt	nd	nt	nt	25	55
PHCs F <sub>2</sub>	4	nd	<u>35</u>	5	nd	10	98
PHCs F <sub>3</sub>	8	131	83	54	nd	240	300
PHCs F <sub>4</sub>	6	877	36	57	nd	120	2,800
	50	1,730	nt	nt	nt	120	2,800

The concentrations of PHCs  $F_2$ ,  $F_3$ , and  $F_{4G}$  in sample BH3-AU1, as well as the concentrations of PHCs  $F_1$ ,  $F_2$ , and  $F_3$  in sample BH8-SS4 are in excess of the MECP Table 3 residential standards and MECP Table 1 background standards.

While in compliance with the selected MECP Table 3 standards, several BTEX and/or PHC parameters were detected in samples: BH1-SS2, BH2-SS3, BH3-SS9/10, BH9-AU1, BH9-SS6, BH11-AU1, and BH11-SS4 at concentrations exceeding the MECP Table 1 standards. These exceedances are not considered to pose an environmental concern to the subject site, however, if this soil is ever to be removed from the property, it should be classified as contaminated and disposed of at an approved waste disposal site.

# Table 6 Analytical Test Results - Soil

			Soil Sam	oles (ug/g)		MECP	MECP	
Parameter	MDL	May 23, 2019	J	uly 15, 201	19	Table 1 Residential Soil	Table 3 Residential Soil	
	(µg/g)	BH3-	BH7-	BH8-	BH8-	Standards	Soll Standards	
		SS9/10	SS3	SS4	SS9	(µg/g)	(µg/g)	
Acetone	0.50	nd	nd	nd	nd	0.5	16	
Benzene	0.02	nd	nd	nd	nd	0.02	0.21	
Bromodichloromethane	0.05	nd	nd	nd	nd	0.05	13	
Bromoform	0.05	nd	nd	nd	nd	0.05	0.27	
Bromomethane	0.05	nd	nd	nd	nd	0.05	0.05	
Carbon Tetrachloride	0.05	nd	nd	nd	nd	0.05	0.05	
Chlorobenzene	0.05	nd	nd	nd	nd	0.05	2.4	
Chloroform	0.05	nd	nd	nd	nd	0.05	0.05	
Dibromochloromethane	0.05	nd	nd	nd	nd	0.05	9.4	
Dichlorodifluoromethane	0.05	nd	nd	nd	nd	0.05	16	
1,2-Dichlorobenzene	0.05	nd	nd	nd	nd	0.05	3.4	
1,3-Dichlorobenzene	0.05	nd	nd	nd	nd	0.05	4.8	
1,4-Dichlorobenzene	0.05	nd	nd	nd	nd	0.05	0.083	
1,1-Dichloroethane	0.05	nd	nd	nd	nd	0.05	3.5	
1,2-Dichloroethane	0.05	nd	nd	nd	nd	0.05	0.05	
1,1-Dichloroethylene	0.05	nd	nd	nd	nd	0.05	0.05	
cis-1,2-Dichloroethylene	0.05	nd	nd	nd	nd	0.05	3.4	
trans-1,2-Dichloroethylene	0.05	nd	nd	nd	nd	0.05	0.084	
1,2-Dichloropropane	0.05	nd	nd	nd	nd	0.05	0.05	
1,3-Dichloropropene	0.05	nd	nd	nd	nd	0.05	0.05	
Ethylbenzene	0.05	nd	nd	nd	nd	0.05	2	
Ethylene Dibromide	0.05	nd	nd	nd	nd	0.05	0.05	
Hexane	0.05	nd	nd	nd	nd	0.05	2.8	
Methyl Ethyl Ketone	0.50	nd	nd	nd	nd	0.5	16	
Methyl Isobutyl Ketone	0.50	nd	nd	nd	nd	0.5	1.7	
Methyl tert-butyl ether	0.05	nd	nd	nd	nd	0.05	0.75	
Methylene Chloride	0.05	nd	nd	nd	nd	0.05	0.1	
Styrene	0.05	nd	nd	nd	nd	0.05	0.7	
1,1,1,2-Tetrachloroethane	0.05	nd	nd	nd	nd	0.05	0.058	
1,1,2,2-Tetrachloroethane	0.05	nd	nd	nd	nd	0.05	0.05	
Tetrachloroethylene	0.05	<u>1.79</u>	nd	nd	nd	0.05	0.28	
Toluene	0.05	nd	nd	nd	nd	0.2	2.3	
1,1,1-Trichloroethane	0.05	nd	nd	nd	nd	0.05	0.38	
1,1,2-Trichloroethane	0.05	nd	nd	nd	nd	0.05	0.05	
Trichloroethylene	0.05	nd	nd	nd	nd	0.05	0.061	
Trichlorofluoromethane	0.05	nd	nd	nd	nd	0.25	4	
Vinyl Chloride	0.02	nd	nd	nd	nd	0.02	0.02	
Xylenes	0.05	<u>0.29</u>	nd	nd	nd	0.05	3.1	
Notes: MDL – Method De nt – not tested for nd – not detected <u>Underlined</u> – Valu Bold and Underli	this parame above the N e exceeds N	ter IDL IECP Table						

Table 6 (continued)
Analytical Test Results – Soil
VOCs

		So	il Samples (u	g/g)	MECP	MECP
Parameter	MDL (µg/g)	July 1	5, 2019	August 24, 2020	Table 1 Residential Soil	Table 3 Residentia Soil
	(1-9, 9)	BH9-SS6	BH9-SS9	BH13-SS7	Standards (µg/g)	Standards (µg/g)
Acetone	0.50	nd	nd	nd	0.5	16
Benzene	0.02	nd	nd	nd	0.02	0.21
Bromodichloromethane	0.05	nd	nd	nd	0.05	13
Bromoform	0.05	nd	nd	nd	0.05	0.27
Bromomethane	0.05	nd	nd	nd	0.05	0.05
Carbon Tetrachloride	0.05	nd	nd	nd	0.05	0.05
Chlorobenzene	0.05	nd	nd	nd	0.05	2.4
Chloroform	0.05	nd	nd	nd	0.05	0.05
Dibromochloromethane	0.05	nd	nd	nd	0.05	9.4
Dichlorodifluoromethane	0.05	nd	nd	nd	0.05	16
1,2-Dichlorobenzene	0.05	nd	nd	nd	0.05	3.4
1,3-Dichlorobenzene	0.05	nd	nd	nd	0.05	4.8
1,4-Dichlorobenzene	0.05	nd	nd	nd	0.05	0.083
1,1-Dichloroethane	0.05	nd	nd	nd	0.05	3.5
1,2-Dichloroethane	0.05	nd	nd	nd	0.05	0.05
1,1-Dichloroethylene	0.05	nd	nd	nd	0.05	0.05
cis-1,2-Dichloroethylene	0.05	nd	nd	nd	0.05	3.4
trans-1,2-Dichloroethylene	0.05	nd	nd	nd	0.05	0.084
1,2-Dichloropropane	0.05	nd	nd	nd	0.05	0.05
1,3-Dichloropropene	0.05	nd	nd	nd	0.05	0.05
Ethylbenzene	0.05	nd	nd	nd	0.05	2
Ethylene Dibromide	0.05	nd	nd	nd	0.05	0.05
Hexane	0.05	nd	nd	nd	0.05	2.8
Methyl Ethyl Ketone	0.50	nd	nd	nd	0.5	16
Methyl Isobutyl Ketone	0.50	nd	nd	nd	0.5	1.7
Methyl tert-butyl ether	0.05	nd	nd	nd	0.05	0.75
Methylene Chloride	0.05	nd	nd	nd	0.05	0.1
Styrene	0.05	nd	nd	nd	0.05	0.7
1,1,1,2-Tetrachloroethane	0.05	nd	nd	nd	0.05	0.058
1,1,2,2-Tetrachloroethane	0.05	nd	nd	nd	0.05	0.05
Tetrachloroethylene	0.05	nd	nd	nd	0.05	0.28
Toluene	0.05	nd	nd	nd	0.2	2.3
1,1,1-Trichloroethane	0.05	nd	nd	nd	0.05	0.38
1,1,2-Trichloroethane	0.05	nd	nd	nd	0.05	0.05
Trichloroethylene	0.05	nd	nd	nd	0.05	0.061
Trichlorofluoromethane	0.05	nd	nd	nd	0.25	4
Vinyl Chloride	0.02	nd	nd	nd	0.02	0.02
Xylenes	0.05	0.87	nd	nd	0.05	3.1

INDE - Method Detection Faint
 Int - not tested for this parameter
 Ind - not detected above the MDL
 <u>Underlined</u> - Value exceeds MECP Table 1 standards
 <u>Bold and Underlined</u> - value exceeds selected MECP standards

All detected VOC parameter concentrations are in compliance with the selected MECP Table 3 residential standards, with one (1) exception. The concentration of tetrachloroethylene in sample BH3-SS9/10 is in excess of the MECP Table 3 residential standards and the MECP Table 1 background standards.

While in compliance with the selected MECP Table 3 standards, the concentrations of xylenes detected in samples: BH3-SS9/10 and BH9-SS6 are in excess of the MECP Table 1 standards. These exceedances are not considered to pose an environmental concern to the subject site, however, if the soil is ever to be removed from the property, it should be classified as contaminated and disposed of at an approved waste disposal site.

		Soil	Samples (µ	ıq/q)	MECP	MECP
Parameter	MDL	May 23, 2019	July 15, 2019	August 24, 2020	Table 1 Residential	Table 3 Residential
i ulullotoi	(µg/g)	BH1- SS2	BH10- SS2	BH11- SS2	Soil Standards (µg/g)	Soil Standards (µg/g)
Acenaphthene	0.02	0.06	nd	nd	0.072	7.9
Acenaphthylene	0.02	0.05	nd	0.05	0.093	0.15
Anthracene	0.02	0.13	nd	0.05	0.16	0.67
Benzo[a]anthracene	0.02	<u>0.37</u>	0.04	0.13	0.36	0.5
Benzo[a]pyrene	0.02	0.29	0.03	0.14	0.3	0.3
Benzo[b]fluoranthene	0.02	0.45	0.04	0.19	0.47	0.78
Benzo[g,h,i]perylene	0.02	0.2	0.03	0.10	0.68	6.6
Benzo[k]fluoranthene	0.02	0.31	0.03	0.10	0.48	0.78
Chrysene	0.02	0.38	0.06	0.13	2.8	7
Dibenzo[a,h]anthracene	0.02	0.07	nd	0.03	0.1	0.1
Fluoranthene	0.02	<u>0.78</u>	0.09	0.27	0.56	0.69
Fluorene	0.02	0.06	nd	nd	0.12	62
Indeno[1,2,3-cd]pyrene	0.02	0.16	nd	0.09	0.23	0.38
1-Methylnaphthalene	0.02	nd	nd	nd	0.59	0.99
2-Methylnaphthalene	0.02	0.03	nd	nd	0.59	0.99
Methylnaphthalene (1&2)	0.04	0.03	nd	nd	0.59	0.99
Naphthalene	0.01	0.04	nd	nd	0.09	0.6
Phenanthrene	0.02	0.58	0.07	0.15	0.69	6.2
Pyrene	0.02	0.62	0.09	0.24	1	78

All detected PAH parameter concentrations are in compliance with the selected MECP Table 3 residential standards, with one (1) exception. The concentration of fluoranthene in sample BH1-SS2 is in excess of the MECP Table 3 residential standards and the MECP Table 1 background standards.

While in compliance with the selected MECP Table 3 standards, the concentration of benzo[a]anthracene detected in samples: BH1-SS2 is in excess of the MECP Table 1 standards. This exceedances is not considered to pose an environmental concern to the subject site, however, if the soil is ever to be removed from the property, it should be classified as contaminated and disposed of at an approved waste disposal site.

				oles (µg/g)		MECP	MECP
		May 2	3, 2019	July 1	5, 2019	Table 1	Table 3
Parameter	MDL (µg/g)	BH1- SS2	BH3- AU1	BH7- AU1	BH10- SS2	Residential Soil Standards (µg/g)	Residential Soil Standards (µg/g)
Antimony	1.0	nd	nd	<u>1.5</u>	nd	1.3	7.5
Arsenic	1.0	7.7	4.0	10.8	5.8	18	18
Barium	1.0	<u>256</u>	54.9	293	171	220	390
Beryllium	0.5	0.6	nd	0.7	0.6	2.5	4
Boron	5.0	15.8	6.8	7.8	6.7	36	120
Cadmium	0.5	nd	nd	0.7	nd	1.2	1.2
Chromium	5.0	30.1	10.1	26.8	19.7	70	160
Chromium (VI)	0.2	nd	nd	nd	nd	0.66	8
Cobalt	1.0	11.7	4.7	9.1	8.1	21	22
Copper	5.0	37.5	13.5	77.6	43.8	92	140
Lead	1.0	<u>181</u>	44.7	<u>295</u>	<u>255</u>	120	120
Mercury	0.1	0.3	nd	0.2	0.3	0.27	0.27
Molybdenum	1.0	<u>3.3</u>	<u>2.2</u>	<u>2.7</u>	1.5	2	6.9
Nickel	5.0	31.6	16.0	27.8	21.0	82	100
Selenium	1.0	nd	nd	1.0	nd	1.5	2.4
Silver	0.3	0.5	nd	0.3	nd	0.5	20
Thallium	1.0	nd	nd	nd	nd	1	1
Uranium	1.0	1.4	nd	nd	nd	2.5	23
Vanadium	10.0	27.5	24.7	28.0	23.0	86	86
Zinc	20.0	143	41.6	279	186	290	340

		Sc	oil Samples (µg	/g)	MECP	MECP
Parameter	MDL (µg/g)	BH11-SS2	August 24, 2020 BH13-SS3	BH14-SS2	Table 1 Residential Soil Standards (µg/g)	Table 3 Residentia Soil Standards (µg/g)
Antimony	1.0	nd	nd	<u>3.9</u>	1.3	7.5
Arsenic	1.0	8.2	2.0	2.3	18	18
Barium	1.0	180	14.8	99.5	220	390
Beryllium	0.5	0.8	nd	nd	2.5	4
Boron	5.0	9.0	nd	7.6	36	120
Cadmium	0.5	nd	nd	nd	1.2	1.2
Chromium	5.0	30.4	7.9	13.7	70	160
Chromium (VI)	0.2	nd	nd	0.3	0.66	8
Cobalt	1.0	9.5	3.5	5.4	21	22
Copper	5.0	63.1	8.0	18.4	92	140
Lead	1.0	226	3.6	22.7	120	120
Mercury	0.1	<u>1.4</u>	nd	0.2	0.27	0.27
Molybdenum	1.0	<u>2.2</u>	nd	nd	2	6.9
Nickel	5.0	34.8	10.1	11.6	82	100
Selenium	1.0	nd	nd	nd	1.5	2.4
Silver	0.3	<u>1.1</u>	nd	nd	0.5	20
Thallium	1.0	nd	nd	nd	1	1
Uranium	1.0	2.4	nd	nd	2.5	23
Vanadium	10.0	30.5	14.6	16.5	86	86
Zinc	20.0	175	nd	36.6	290	340

The concentration of lead in sample BH7-AU1, as well as the concentrations of lead and mercury in samples BH1-SS2, BH10-SS2, and BH11-SS2 are in excess of the selected MECP Table 3 residential standards and the MECP Table 1 background standards.

While in compliance with the selected MECP Table 3 standards, several metal parameters were detected in samples: BH1-SS2, BH3-AU1, BH7-AU1, BH11-SS2, and BH14-SS2 at concentrations exceeding the MECP Table 1 standards. These exceedances are not considered to pose an environmental concern to the subject site, however, if the soil is ever to be removed from the property, it should be classified as contaminated and disposed of at an approved waste disposal site.

Table 9 Analytical Test Results – Soil pH					
			oles (µg/g)	MECP Table 3	
Parameter	MDL (units)	May 23	3, 2019	Residential Standards	
	(units)	BH1-SS2	BH2-SS3	(units)	
рН	0.05	7.94	7.39	5.00 - 9.00	
<ul> <li>nt – not</li> <li>nd – not</li> <li><u>Underlin</u></li> </ul>		rameter	9 standards		

The pH levels detected in the soil samples analyzed are in compliance with the selected MECP Table 3 residential standards.

# DatersongroupOttawaKingstonNorth Bay

Parameter	Maximum Concentration (µg/g)	Sample ID	Depth Interval (m BGS)
Toluene	0.05	BH1-SS2	0.76 – 1.37 m
Kylenes	<u>0.87</u>	BH9-SS6	3.81 – 4.42 m
PHCs F1	<u>60</u>	BH8-SS4	2.29 – 2.90 m
HCs F <sub>2</sub>	<u>2,980</u>	BH8-SS4	2.29 – 2.90 m
PHCs F <sub>3</sub>	<u>2,370</u>	BH3-AU1	0.00 – 0.61 m
HCs F4	2,590	BH3-AU1	0.00 – 0.61 m
PHCs F <sub>4G</sub>	<u>9,490</u>	BH3-AU1	0.00 – 0.61 m
etrachloroethylene	<u>1.79</u>	BH3-SS9/10	6.10 – 7.47 m
Acenaphthene	0.06	BH1-SS2	0.76 – 1.37 m
cenaphthylene	0.05	BH1-SS2 / BH11-SS2	0.76 – 1.37 m / 0.76 – 1.37 m
Anthracene	0.13	BH1-SS2	0.76 – 1.37 m
Benzo[a]anthracene	<u>0.37</u>	BH1-SS2	0.76 – 1.37 m
Benzo[a]pyrene	0.29	BH1-SS2	0.76 – 1.37 m
enzo[b]fluoranthene	0.45	BH1-SS2	0.76 – 1.37 m
enzo[g,h,i]perylene	0.2	BH1-SS2	0.76 – 1.37 m
Benzo[k]fluoranthene	0.31	BH1-SS2	0.76 – 1.37 m
Chrysene	0.38	BH1-SS2	0.76 – 1.37 m
bibenzo[a,h]anthracene	0.07	BH1-SS2	0.76 – 1.37 m
luoranthene	<u>0.78</u>	BH1-SS2	0.76 – 1.37 m
luorene	0.06	BH1-SS2	0.76 – 1.37 m
ndeno[1,2,3-cd]pyrene	0.16	BH1-SS2	0.76 – 1.37 m
-Methylnaphthalene	0.03	BH1-SS2	0.76 – 1.37 m
lethylnaphthalene (1&2)	0.03	BH1-SS2	0.76 – 1.37 m
laphthalene	0.04	BH1-SS2	0.76 – 1.37 m
henanthrene	0.58	BH1-SS2	0.76 – 1.37 m
Pyrene	0.62	BH1-SS2	0.76 – 1.37 m
Intimony	<u>3.9</u>	BH14-SS2	0.76 – 1.37 m
rsenic	10.8	BH7-AU1	0.00 – 0.61 m
Barium	<u>293</u>	BH7-AU1	<u>0.00 – 0.61 m</u>
eryllium	0.8	BH11-SS2	0.76 – 1.37 m
Boron	15.8	BH1-SS2	0.76 – 1.37 m
admium	0.7	BH7-AU1	0.00 - 0.61 m
hromium	30.4	BH11-SS2	0.76 – 1.37 m
Chromium (VI)	0.3	BH14-SS2	0.76 – 1.37 m
obalt	11.7 77.6	BH1-SS2 BH7-AU1	<u>0.76 – 1.37 m</u> 0.00 – 0.61 m
Copper			
ead Iercury	<u>295</u>	BH7-AU1	<u>0.00 – 0.61 m</u>
lercury lolybdenum	<u>1.4</u> 3.3	BH11-SS2 BH1-SS2	<u>0.76 – 1.37 m</u> 0.76 – 1.37 m
ickel	<u>3.3</u> 34.8	BH11-SS2	0.76 – 1.37 m
elenium	1.0	BH7-AU1	0.00 – 0.61 m
ilver	1.1	BH11-SS2	0.76 – 1.37 m
Iranium	2.4	BH11-SS2 BH11-SS2	0.76 – 1.37 m
anadium	30.5	BH11-SS2 BH11-SS2	0.76 – 1.37 m
inc	279	BH7-AU1	0.00 – 0.61 m
H	7.94	BH1-SS2	0.76 – 1.37 m

All other parameter concentrations analyzed were below the laboratory detection limits.

# 5.6 Groundwater Quality

As part of this assessment, twelve (12) groundwater samples, were submitted for laboratory analysis of either: BTEX, PHCs, VOCs, and/or PAHs parameters. The results of the analytical testing are presented below in Tables 11 to 13. The laboratory certificates of analysis are provided in Appendix 1.

			Gr	oundwa	ter Sam	ples (µg	/L)		MECP Table 3
Parameter	MDL	Ju	ne 3, 20	19		July 19	9, 2019		Residential Groundwater
i urumeter	(µg/L)	BH2- GW1	BH3- GW1	BH4- GW1	BH3- GW2	BH7- GW1	BH9- GW1	BH10 -GW1	Standards (µg/L)
Benzene	0.5	nd	44						
Ethylbenzene	0.5	nd	2,300						
Toluene	0.5	nd	18,000						
Xylenes	0.5	nd	4,200						
PHC F1	25	nd	nd	nd	nt	nd	nt	nd	750
PHC F <sub>2</sub>	100	nd	nd	nd	nt	nd	nt	nd	150
PHC F <sub>3</sub>	100	nd	nd	nd	nt	nd	nt	nd	500
PHC F4	100	nd	nd	nd	nt	nd	nt	nd	500

Bold and Underlined – value exceeds selected MECP standards

	MDL		Groundwater Septemb	Samples (µg/L er 3, 2020	)	MECP Table 3 Residential
Parameter	(µg/L)	BH3-GW3	BH8B-GW1	BH11-GW1	BH13-GW1	Groundwater Standards (µg/L)
Benzene	0.5	nd	nd	nd	nd	44
Ethylbenzene	0.5	nd	nd	0.6	nd	2,300
Toluene	0.5	nd	nd	nd	nd	18,000
Xylenes	0.5	nd	nd	0.7	nd	4,200
PHC F1	25	nt	nt	nd	nd	750
PHC F <sub>2</sub>	100	nt	nt	nd	nd	150
PHC F <sub>3</sub>	100	nt	nt	nd	nd	500
PHC F <sub>4</sub>	100	nt	nt	nd	nd	500

All detected BTEX and PHC parameters are in compliance with the selected MECP Table 3 residential standards.

# Table 12 Analytical Test Results – Groundwater

			Grou	ndwater	Sample (	ug/L)		MECP
	MDL	Ju	une 3, 20	19	Ju	ily 19, 20	19	Table 3 Residential
Parameter	(μg/L)	BH2- GW1	BH3- GW1	BH4- GW1	BH3- GW2	BH7- GW1	BH9- GW1	Groundwater Standards (µg/L)
Acetone	5.0	nd	nd	nd	nd	nd	nd	130,000
Benzene	0.5	nd	nd	nd	nd	nd	nd	44
Bromodichloromethane	0.5	nd	nd	nd	nd	nd	nd	85,000
Bromoform	0.5	nd	nd	nd	nd	nd	nd	380
Bromomethane	0.5	nd	nd	nd	nd	nd	nd	5.6
Carbon Tetrachloride	0.2	nd	nd	nd	nd	nd	nd	0.79
Chlorobenzene	0.5	nd	nd	nd	nd	nd	nd	630
Chloroform	0.5	nd	nd	nd	nd	nd	nd	2.4
Dibromochloromethane	0.5	nd	nd	nd	nd	nd	nd	82,000
Dichlorodifluoromethane	1.0	nd	nd	nd	nd	nd	nd	4,400
1,2-Dichlorobenzene	0.5	nd	nd	nd	nd	nd	nd	4,600
1,3-Dichlorobenzene	0.5	nd	nd	nd	nd	nd	nd	9,600
1,4-Dichlorobenzene	0.5	nd	nd	nd	nd	nd	nd	8
1,1-Dichloroethane	0.5	nd	nd	nd	nd	nd	nd	320
1,2-Dichloroethane	0.5	nd	nd	nd	nd	nd	nd	1.6
1,1-Dichloroethylene	0.5	nd	nd	nd	nd	nd	nd	1.6
cis-1,2-Dichloroethylene	0.5	nd	nd	nd	nd	nd	nd	1.6
trans-1,2-Dichloroethylene	0.5	nd	nd	nd	nd	nd	nd	1.6
1,2-Dichloropropane	0.5	nd	nd	nd	nd	nd	nd	16
1,3-Dichloropropene	0.5	nd	nd	nd	nd	nd	nd	5.2
Ethylbenzene	0.5	nd	nd	nd	nd	nd	nd	2,300
Ethylene Dibromide	0.2	nd	nd	nd	nd	nd	nd	0.25
Hexane	1.0	nd	nd	nd	nd	nd	nd	51
Methyl Ethyl Ketone	5.0	nd	nd	nd	nd	nd	nd	470,000
Methyl Isobutyl Ketone	5.0	nd	nd	nd	nd	nd	nd	140,000
Methyl tert-butyl ether	2.0	nd	nd	nd	nd	nd	nd	190
Methylene Chloride	5.0	nd	nd	nd	nd	nd	nd	610
Styrene	0.5	nd	nd	nd	nd	nd	nd	1,300
1,1,1,2-Tetrachloroethane	0.5	nd	nd	nd	nd	nd	nd	3.3
1,1,2,2-Tetrachloroethane	0.5	nd	nd	nd	nd	nd	nd	3.2
Tetrachloroethylene	0.5	nd	nd	nd	nd	nd	nd	1.6
Toluene	0.5	nd	nd	nd	nd	nd	nd	18,000
1,1,1-Trichloroethane	0.5	nd	nd	nd	nd	nd	nd	640
1,1,2-Trichloroethane	0.5	nd	nd	nd	nd	nd	nd	4.7
Trichloroethylene	0.5	nd	nd	nd	nd	nd	nd	1.6
Trichlorofluoromethane	1.0	nd	nd	nd	nd	nd	nd	2,500
Vinyl Chloride	0.5	nd	nd	nd	nd	nd	nd	0.5
Xylenes	0.5	nd	nd	nd	nd	nd	nd	4,200

nd – not detected above the MDL <u>Bold and Underlined</u> – value exceeds selected MECP standards 

# Table 12 (continued) Analytical Test Results - Groundwater

North Bay

Ditawa Kingston North Bay

			Ground	water Samp	ole (ug/L)		MECP
Parameter	MDL (µg/L)	July 19, 2019		Septemb	er 3, 2020		Table 3 Residential Groundwater
	(µg/⊏)	BH10- GW1	BH3- GW3	BH8B- GW1	BH11- GW1	BH13- GW1	Standards (µg/L)
Acetone	5.0	nd	nd	nd	nd	nd	130,000
Benzene	0.5	nd	nd	nd	nd	nd	44
Bromodichloromethane	0.5	nd	nd	nd	nd	nd	85,000
Bromoform	0.5	nd	nd	nd	nd	nd	380
Bromomethane	0.5	nd	nd	nd	nd	nd	5.6
Carbon Tetrachloride	0.2	nd	nd	nd	nd	nd	0.79
Chlorobenzene	0.5	nd	nd	nd	nd	nd	630
Chloroform	0.5	nd	nd	nd	nd	nd	2.4
Dibromochloromethane	0.5	nd	nd	nd	nd	nd	82,000
Dichlorodifluoromethane	1.0	nd	nd	nd	nd	nd	4,400
1,2-Dichlorobenzene	0.5	nd	nd	nd	nd	nd	4,600
1,3-Dichlorobenzene	0.5	nd	nd	nd	nd	nd	9,600
1,4-Dichlorobenzene	0.5	nd	nd	nd	nd	nd	8
1,1-Dichloroethane	0.5	nd	nd	nd	nd	nd	320
1,2-Dichloroethane	0.5	nd	nd	nd	nd	nd	1.6
1,1-Dichloroethylene	0.5	nd	nd	nd	nd	nd	1.6
cis-1,2-Dichloroethylene	0.5	nd	nd	nd	nd	nd	1.6
trans-1,2-Dichloroethylene	0.5	nd	nd	nd	nd	nd	1.6
1,2-Dichloropropane	0.5	nd	nd	nd	nd	nd	16
1,3-Dichloropropene	0.5	nd	nd	nd	nd	nd	5.2
Ethylbenzene	0.5	nd	nd	nd	0.6	nd	2,300
Ethylene Dibromide	0.2	nd	nd	nd	nd	nd	0.25
Hexane	1.0	nd	nd	nd	nd	nd	51
Methyl Ethyl Ketone	5.0	nd	nd	nd	nd	nd	470,000
Methyl Isobutyl Ketone	5.0	nd	nd	nd	nd	nd	140,000
Methyl tert-butyl ether	2.0	nd	nd	nd	nd	nd	190
Methylene Chloride	5.0	nd	nd	nd	nd	nd	610
Styrene	0.5	nd	nd	nd	nd	nd	1,300
1,1,1,2-Tetrachloroethane	0.5	nd	nd	nd	nd	nd	3.3
1,1,2,2-Tetrachloroethane	0.5	nd	nd	nd	nd	nd	3.2
Tetrachloroethylene	0.5	nd	nd	nd	nd	nd	1.6
Toluene	0.5	nd	nd	nd	nd	nd	18,000
1,1,1-Trichloroethane	0.5	nd	nd	nd	nd	nd	640
1,1,2-Trichloroethane	0.5	nd	nd	nd	nd	nd	4.7
Trichloroethylene	0.5	nd	nd	nd	nd	nd	1.6
Trichlorofluoromethane	1.0	nd	nd	nd	nd	nd	2,500
Vinyl Chloride	0.5	nd	nd	nd	nd	nd	0.5
Xylenes	0.5	nd	nd	nd	0.7	nd	4,200
Notes: MDL – Method De nt – not tested for nd – not detected Bold and Underli	this param above the	neter MDL	elected ME0	CP standards			

All detected VOC parameter concentrations are in compliance with the selected MECP Table 3 residential standards.

#### Table 13 Analytical Test Results – Groundwater PAHs

	-	Grou	ndwater Sample	(µg/L)	MECP Table 3	
Parameter	MDL (µg/L)	BH2-GW1	May 23, 2019 BH3-GW1	BH4-GW1	Residential Groundwater Standards (µg/L)	
Acenaphthene	0.05	nd	nd	nd	600	
Acenaphthylene	0.05	nd	nd	nd	1.8	
Anthracene	0.01	0.05	nd	nd	2.4	
Benzo[a]anthracene	0.01	nd	nd	nd	4.7	
Benzo[a]pyrene	0.01	0.03	nd	nd	0.81	
Benzo[b]fluoranthene	0.05	nd	nd	nd	0.75	
Benzo[g,h,i]perylene	0.05	nd	nd	nd	0.2	
Benzo[k]fluoranthene	0.05	nd	nd	nd	0.4	
Chrysene	0.05	nd	nd	nd	1	
Dibenzo[a,h]anthracene	0.05	nd	nd	nd	0.52	
Fluoranthene	0.01	0.07	nd	nd	130	
Fluorene	0.05	nd	nd	nd	400	
Indeno[1,2,3-cd]pyrene	0.05	nd	nd	nd	0.2	
1-Methylnaphthalene	0.05	nd	nd	nd	1,800	
2-Methylnaphthalene	0.05	nd	nd	nd	1,800	
Methylnaphthalene (1&2)	0.10	nd	nd	nd	1,800	
Naphthalene	0.05	nd	nd	nd	1,400	
Phenanthrene	0.05	nd	nd	nd	580	
Pyrene	0.01	0.06	nd	nd	60	

Ind – not detected above the MDL
Bold and Underlined – value exceeds selected MEC

Bold and Underlined – value exceeds selected MECP standards

All detected PAH parameter concentrations are in compliance with the selected MECP Table 3 residential standards.

Parameter	Maximum Concentration (μg/L)	Sample ID	Depth Interval (m BGS)	
Ethylbenzene	0.6	BH11-GW1	2.84 – 5.84 m	
Xylenes	0.7	BH11-GW1	2.84 – 5.84 m	
Anthracene	0.05	BH2-GW1	5.36 – 6.86 m	
Benzo[a]pyrene	0.03	BH2-GW1	5.36 – 6.86 m	
Fluoranthene	0.07	BH2-GW1	5.36 – 6.86 m	
Pyrene	0.06	BH2-GW1	5.36 – 6.86 m	

# 5.7 Quality Assurance and Quality Control Results

As per the Sampling and Analysis Plan, a duplicate groundwater sample was obtained from BH3 during the July 19, 2019 sampling event as well as from BH13 during the September 3, 2020 sampling event, and analyzed for VOCs. Based on the analytical test results, no VOC parameters were detected above the laboratory method detection limits in either the original sample or the field duplicate sample. As a result, the quality of the field data collected during this Phase II ESA is considered to be sufficient to meet the overall objectives of this assessment.

All samples submitted as part of this Phase II ESA were handled in accordance with the analytical protocols with respect to holding time, preservation method, storage requirement, and container type. As per Subsection 47(3) of O.Reg. 153/04, as amended by the Environmental Protection Act, the certificates of analysis have been received for each sample submitted for laboratory analysis and have been appended to this report.

# 5.8 Updated Phase II Conceptual Site Model

The following section has been prepared in accordance with the requirements of O.Reg. 269/11, amending O.Reg. 153/04 - Record of Site Condition regulation, made under the Environmental Protection Act. Conclusions and recommendations are discussed in a subsequent section.

# Site Description

# Potentially Contaminating Activities (PCAs) and Areas of Potential Environmental Concern (APECs)

As described in Section 6.1 of the Phase I ESA report, as well as in Section 2.2 of this report, the following PCAs, as described by Table 2 of O.Reg. 153/04, are considered to result in APECs on the subject site:

#### "Item 10: Commercial Autobody Shops"

This PCA was identified as a result of the presence of a former autobody shop and service garage located at 89 Barrette Street.

"Item 28: Gasoline and Associated Products Storage in Fixed Tanks"

This PCA was identified as a result of the presence of a former retail fuel outlet located at 64 Beechwood Avenue.

"Item 37: Operation of Dry-Cleaning Equipment (Where Chemicals Are Used)"

This PCA was identified as a result of the presence of a dry cleaners located at 110 Beechwood Avenue.

"Item 52: Storage, Maintenance, Fuelling, and Repair of Equipment, Vehicles, and Material Used to Maintain Transportation Systems"

This PCA was identified as a result of the presence of a former autobody shop and service garage located at 89 Barrette Street.

"Item 58: Waste Disposal and Waste Management, including thermal treatment, landfilling and transfer of waste, other than use of biosoils as soil conditioners"

This PCA was identified as a result of the presence of a former landfill site, located throughout the Phase I study area and situated beneath the entirety of the subject site.

#### **Contaminants of Potential Concern**

The contaminants of potential concern (CPCs) associated with the aforementioned APECs are considered to be:

- □ Volatile Organic Compounds (VOCs);
- Benzene, Toluene, Ethylbenzene, and Xylenes (BTEX);
- Petroleum Hydrocarbons, fractions 1 4 (PHCs F<sub>1</sub>-F<sub>4</sub>);
- □ Metals (including Mercury and Hexavalent Chromium);
- D Polycyclic Aromatic Hydrocarbons (PAHs).

These contaminants have the potential to be present in the soil/fill matrix as well as the groundwater on the subject site.

#### Subsurface Structures and Utilities

Underground service locates were completed as part of a Phase II ESA conducted for the subject site in tandem with this assessment. According to the locates, underground gas lines, electrical cables, as well as water and sewer pipes are present on the subject site.

# **Physical Setting**

#### Site Stratigraphy

The stratigraphy of the subject site generally consists of:

- Fill material, consisting of either crushed stone and/or brown silty sand and gravel, extending to depths ranging from approximately 0.28 m to 3.20 m below ground surface.
- Glacial till, consisting of either brown clayey silt or brown silty sand and gravel, extending to depths of approximately 1.52 m to 6.86 m below ground surface.
- □ Weathered shale bedrock; encountered at depths ranging from approximately 1.52 m to 6.86 m below the existing ground surface.

#### Hydrogeological Characteristics

Groundwater levels at the subject site were measured at each monitoring well, with the exception of BH7 and BH9, on September 3, 2020. The water table at the subject site was encountered within the underlying bedrock unit, at depths ranging from approximately 3.20 m to 6.37 m below the existing ground surface.

Based on the measured groundwater levels, the groundwater flow in the vicinity of the subject site is towards the north.

#### Approximate Depth to Bedrock

Bedrock was encountered in all boreholes at depths ranging from approximately 1.52 m to 6.86 m below ground surface, at the time of the drilling program.

#### Approximate Depth to Water Table

Based on the findings of the current groundwater sampling event, the depth to water table at the subject property is approximately 3.20 m to 6.37 m below the existing ground surface.

#### Fill Placement

Fill material, consisting of either crushed stone and/or brown silty sand and gravel with construction/demolition debris, was encountered throughout the entire subject site area and extending to depths ranging from approximately 0.28 m to 3.20 m below ground surface.

#### Sections 41 and 43.1 of the Regulation

Section 41 of the Regulation does not apply to the subject site as there are no areas of natural significance or bodies of water located on the subject site or within 30 m of the subject site. The subject site is not considered to be environmentally sensitive.

Section 43.1 of the Regulation does not apply to the subject site since the bedrock is located at a depth of greater than 2 m below the ground surface, and thus is not considered to be a Shallow Soil Property.

#### **Existing Buildings and Structures**

The subject site is currently occupied by twelve (12) buildings of mixed residential and commercial uses.

#### Proposed Buildings and Other Structures

It is our understanding that the subject site is to be redeveloped with a multistorey residential building with ground floor commercial units and underground parking. A Record of Site Condition (RSC) will be required due to the conversion to a more sensitive land use.

#### Water Bodies and Areas of Natural Significance

There are no waterbodies or areas of natural and scientific interest located on the subject site or within the Phase I study area. The nearest named waterbody with respect to the subject site is the Rideau River, located approximately 315 m to the south.

### **Environmental Condition**

### Areas Where Contaminants are Present

Based on the findings of this assessment, the soil/fill within the following areas were identified as being contaminated with various contaminants:

- Petroleum Hydrocarbons: identified within the soil at BH3 and BH8, located within the northeastern portion of the subject site;
- □ Volatile Organic Compounds: identified within the soil at BH3, located in the northeastern portion of the subject site;
- Polycyclic Aromatic Hydrocarbons: identified within the soil at BH1, located in the north-central portion of the subject site;
- Metals: identified within the soil at BH1, BH7, BH10, and BH11, located in the central, and western portions of the subject site.

No impacted groundwater was identified on the subject site. The analytical test results for all soil and groundwater samples tested are shown on the Analytical Testing Plans, appended to this report.

#### Types of Contaminants

Based on the findings of this assessment, the contaminants of concern identified on the subject site are summarized below in Table 15:

Table 15         Contaminants of Concern (CPCs)					
SOIL					
Contaminant Group	Parameters	Contaminant Locations			
Petroleum Hydrocarbons (PHCs)	PHCs F1           PHCs F2           PHCs F3           PHCs F4           PHCs F4G	BH3 / BH8			
Volatile Organic Compounds (VOCs)	Tetrachloroethylene	ВНЗ			
Polycyclic Aromatic Hydrocarbons (PAHs)	Fluoranthene	BH1			
Metals	Lead Mercury	BH1 / BH7/ BH10 / BH11			

### Contaminated Media

As noted above in Table 15, PHC, VOC, PAH, and/or metal impacted soil/fill was identified at BH1, BH3, BH7, BH8, BH10, and BH11. These boreholes are located within the northeastern, central, and western portions of the subject site.

Based on the analytical results, the groundwater is not contaminated.

### What Is Known About Areas Where Contaminants Are Present

Based on the findings of this assessment, PHC, VOC, PAH, and/or metal impacted soil was identified in the vicinity of BH1, BH3, BH7, BH8, BH10, and BH11.

The metal and PAH contaminants are likely a result of historical landfilling on the subject site. The PHC contaminants are likely the result of an unknown discharge on the subject site. The VOC contaminants are likely the result of the neighbouring dry cleaners.

### **Distribution and Migration of Contaminants**

PHC impacted soil was identified in the vicinity of BH3 and BH8, located in the northeastern portion of the subject site. Based on their low mobility, as well as the clean groundwater results, it is anticipated that these contaminants are primarily contained within the fill material in this portion of the subject site. Some limited vertical migration of these contaminants is expected to have occurred, based on their detection within a deeper native soil sample (BH8-SS4).

VOC impacted soil was identified in the vicinity of BH3, located in the northeastern portion of the subject site. Based on its depth, as well as the clean groundwater results, it is anticipated that this contaminant likely originated from the neighbouring dry-cleaners at 110 Beechwood Avenue.

PAH impacted soil was identified in the vicinity of BH1, located in the northcentral portion of the subject site. Based on their low mobility, it is anticipated that these contaminants are primarily contained within the fill material in this portion of the subject site.

Metal impacted soil was identified in the vicinity of BH1, BH7, BH10, and BH11, located in the central and western portions of the subject site. Based on their low mobility, it is anticipated that these contaminants are primarily contained within the fill material in these portions of the subject site.

### Discharge of Contaminants

The PHC impacted soil in the vicinity of BH3 and BH8, located in the northeastern portion of the subject site, is considered to have resulted from the importation of poor-quality fill material and/or the presence of building debris.

The VOC impacted soil in the vicinity of BH3, located in the northeastern portion of the subject site, is considered to have resulted from the neighbouring drycleaners at 110 Beechwood Avenue.

The PAH impacted soil in the vicinity of BH1, located in the north-central portion of the subject site, is considered to have resulted from the importation of poorquality fill material and/or the presence of building debris.

The metal impacted soil in the vicinity of BH1, BH7, BH10, and BH11, located in the central, and western portions of the subject site, is considered to have resulted from the importation of poor-quality fill material and/or the presence of building debris.

### Climatic and Meteorological Conditions

In general, climatic and meteorological conditions have the potential to affect contaminant distribution. Two (2) ways by which climatic and meteorological conditions may affect contaminant distribution include the downward leaching of contaminants via the infiltration of precipitation, and the migration of contaminants via groundwater levels and/or flow, which may fluctuate seasonally.

Downward leaching is not considered to have affected any contaminant distribution at the subject site, since the groundwater test results comply with the MECP Table 3 residential standards. Fluctuations in the groundwater level and flow are also not considered to have affected any contaminant distribution, due to the depth of the water table within the bedrock, well below the shallow fill material.

### Potential for Vapour Intrusion

Given the location of impacted soil outside of the building footprints, as well as their relatively low-volatility, the potential for vapours to be present within the subject structures is considered to be low and does not pose a safety hazard to the current occupants. During redevelopment, all soil exceeding the MECP Table 3 residential standards will be removed and disposed off-site. As such, there is no anticipated potential for future vapour intrusion at the subject site.

## 6.0 CONCLUSION

A Phase II ESA was conducted for the properties addressed 78-90 Beechwood Avenue and 69-93 Barrette Street, in the City of Ottawa, Ontario. The purpose of the Phase II ESA was to address the potentially contaminating activities (PCAs) that were identified during the Phase I ESA and were considered to result in areas of potential environmental concern (APECs) on the subject site.

The subsurface investigation consisted of drilling fifteen (15) boreholes (BH1-BH14 and BH8B) on May 23-24, and July 15, 2019, as well as on August 24, 2020. Upon completion, ten (10) of the boreholes were instrumented with groundwater monitoring wells (BH2, BH3, BH4, BH7, BH8 BH8B, BH9, BH10, BH11, and BH13). The boreholes were drilled to depths ranging from approximately 3.25 m to 9.14 m below ground surface and terminated within weathered black shale bedrock.

Nineteen (19) soil samples were submitted for laboratory analysis of either: BTEX, PHCs F<sub>1</sub>-F<sub>4</sub>, VOCs, PAHs, metals, and/or pH parameters. According to the analytical results, the concentrations of some PHC, VOC, PAH, and/or metal parameters in the soil samples tested from BH1, BH3, BH7, BH8, BH10, and BH11 were in excess of the selected MECP Table 3 residential standards.

Twelve (12) groundwater samples were submitted for laboratory analysis of either: BTEX, PHCs F<sub>1</sub>-F<sub>4</sub>, VOCs, and/or PAHs parameters. According to the analytical results, all detected parameter concentrations in the groundwater samples analyzed comply with the selected MECP Table 3 residential standards.

### Recommendations

### Soil

Based on the findings of this assessment, PHC, VOC, PAH, and/or metal impacted soil was identified in the vicinity of BH1, BH3, BH7, BH8, BH10, and BH11, requiring some remedial work. It is our understanding that the subject site is to be redeveloped with a multi-storey residential high-rise building in the near future. Therefore, it is our recommendation that an environmental site remediation program be completed in conjunction with site redevelopment. This will require the segregation of clean soils from impacted soils, the latter of which will require disposal at an approved waste disposal facility.

It should be noted that several PHC, VOC, PAH, and metal parameters were detected in BH1, BH2, BH3, BH7 BH9, BH11, and/or BH14 at concentrations exceeding the MECP Table 1 background standards. These exceedances are not considered to pose an environmental concern to the subject site, however, if the soil is to be removed from the property for construction purposes, it should be classified as contaminated and disposed of at an approved waste disposal site.

Prior to off-site disposal at a licensed landfill, a leachate analysis of a representative sample of contaminated soil must be conducted in accordance with Ontario Regulation 347/558.

It is recommended that Paterson personnel be present on-site during remediation activities to direct the excavation and segregation of impacted soil, as well as to conduct confirmatory sampling as required.

### Monitoring Wells

If the groundwater monitoring wells installed on-site are not going to be used in the future, or will be destroyed during future redevelopment activities, then they must be decommissioned according to Ontario Regulation Reg. 903 (Ontario Water Resources Act). The monitoring wells will be registered with the MECP under this regulation. Further information can be provided upon request in this regard.

## 7.0 STATEMENT OF LIMITATIONS

This Phase II - Environmental Site Assessment report has been prepared as per the agreed scope-of-work, in general accordance with O.Reg. 153/04, as amended by O.Reg. 269/11, and meets the requirements of CSA Z769-00. The conclusions presented herein are based on information gathered from a limited sampling and testing program. The test results represent conditions at specific test locations at the time of the field program.

The client should be aware that any information pertaining to soils and all test hole logs are furnished as a matter of general information only and test hole descriptions or logs are not to be interpreted as descriptive of conditions at locations other than those of the test holes themselves.

Should any conditions be encountered at the subject site and/or historical information that differ from our findings, we request that we be notified immediately in order to allow for a reassessment.

This report was prepared for the sole use of Minto Communities. Permission and notification from Minto Communities and Paterson will be required to release this report to any other party.

### Paterson Group Inc.

N. Sullin

Nick Sullivan, B.Sc.

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Mark S. D'Arcy, P.Eng., QPESA

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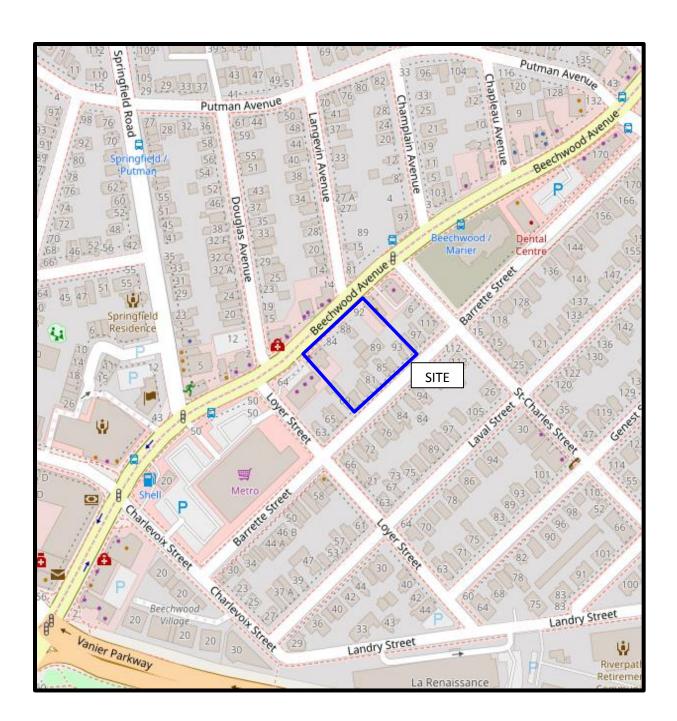
## **FIGURES**

#### FIGURE 1 – KEY PLAN

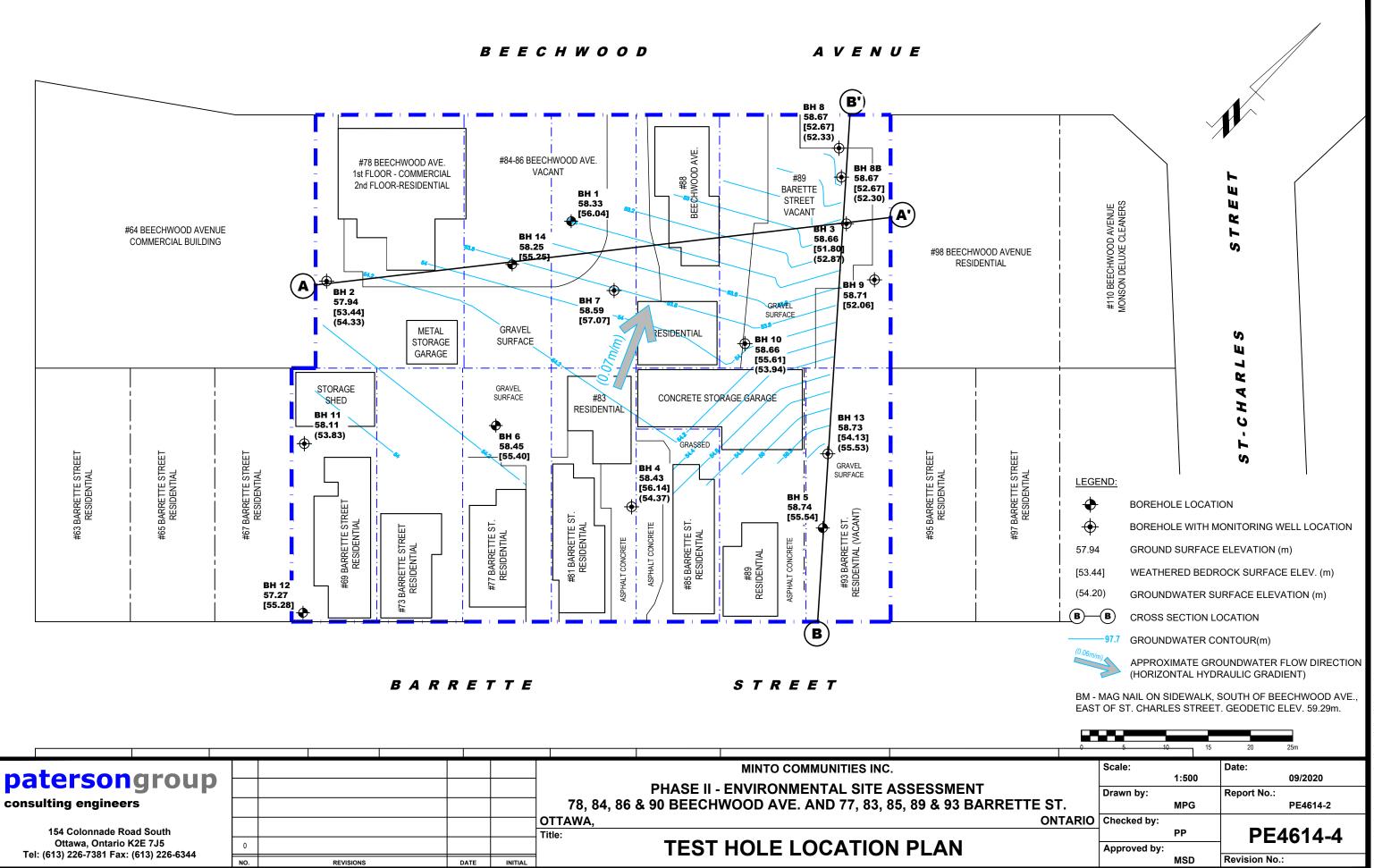
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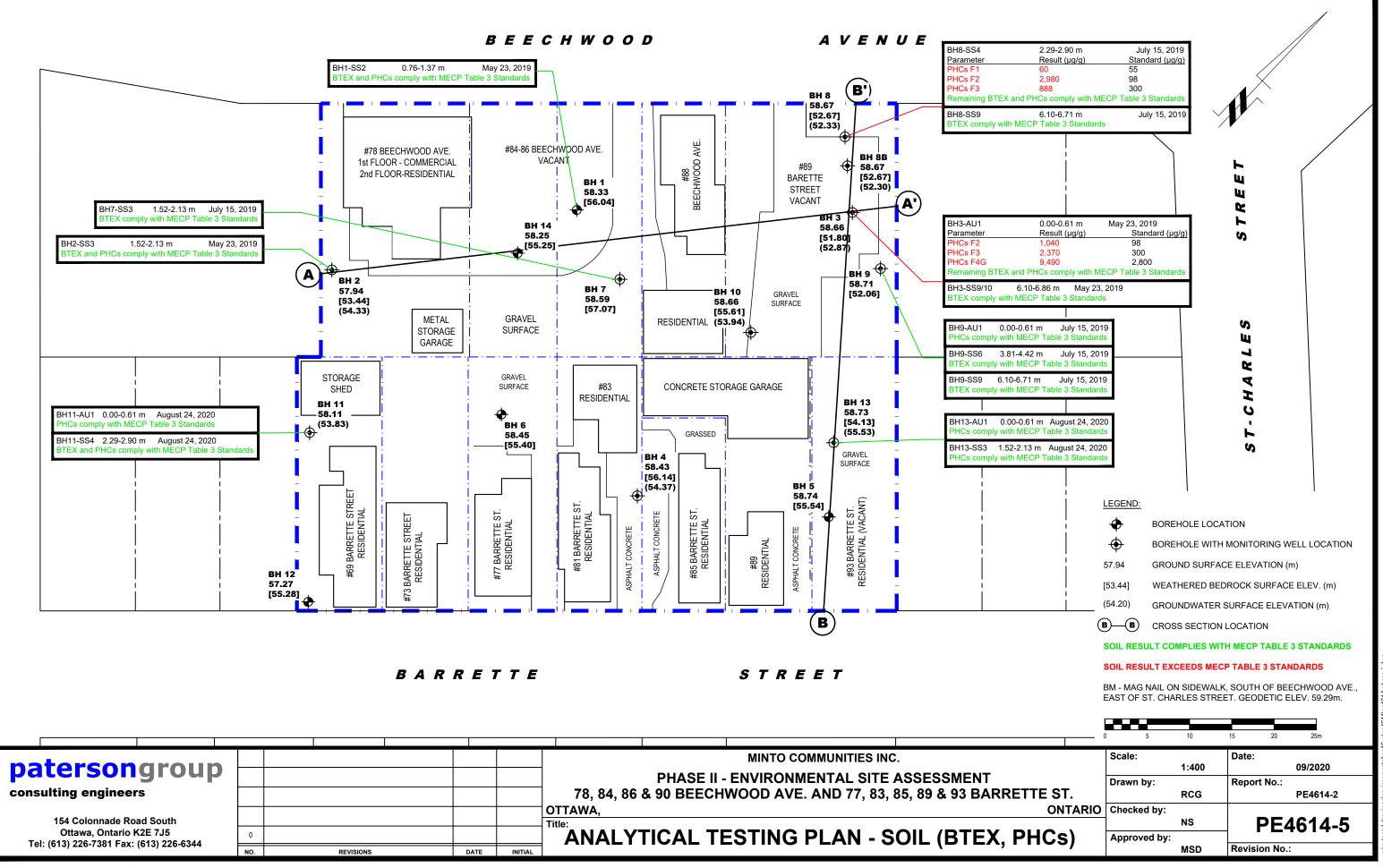
## patersongroup

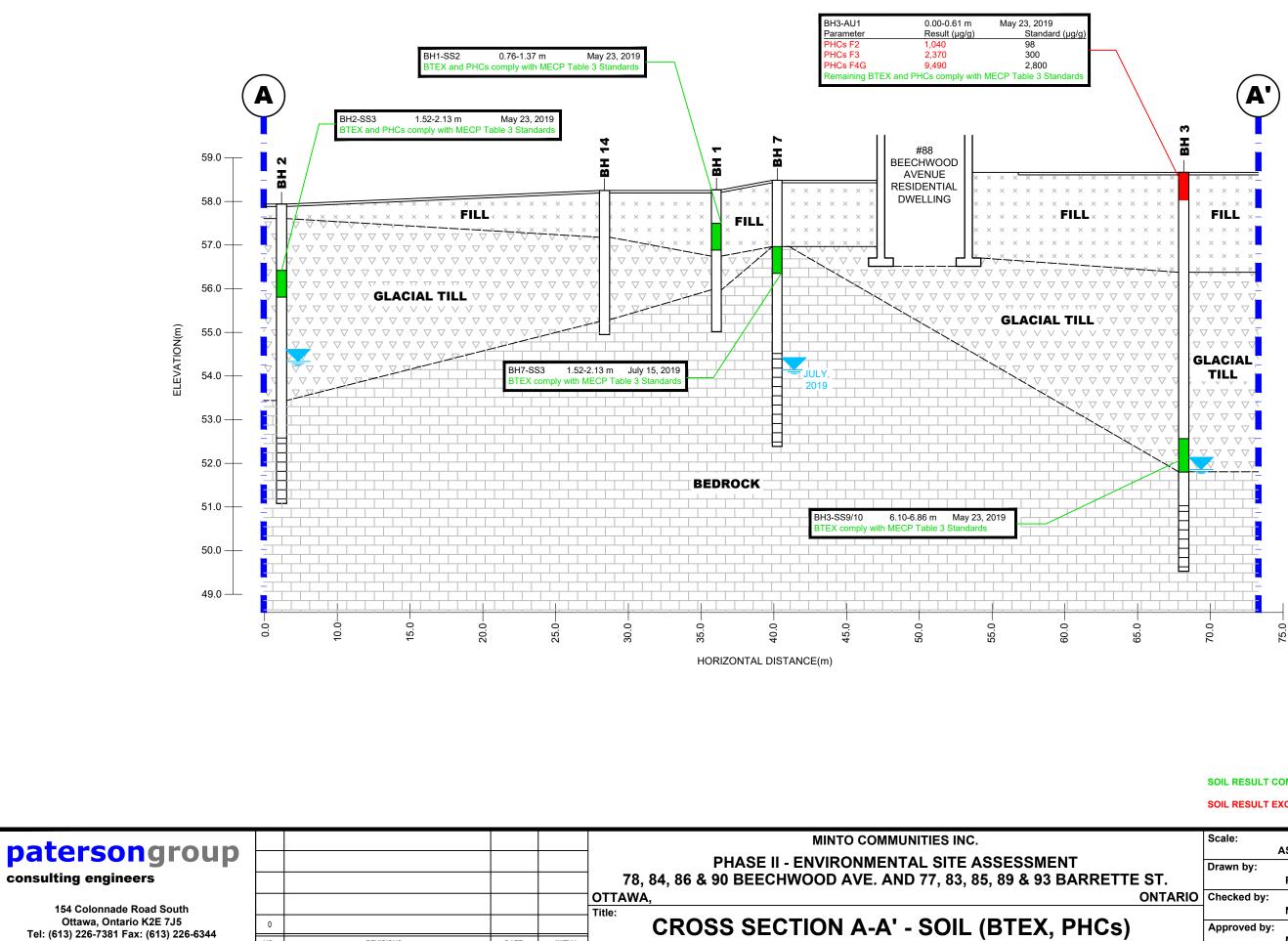
## <u>figure 1</u> KEY PLAN



BEECHWOOD







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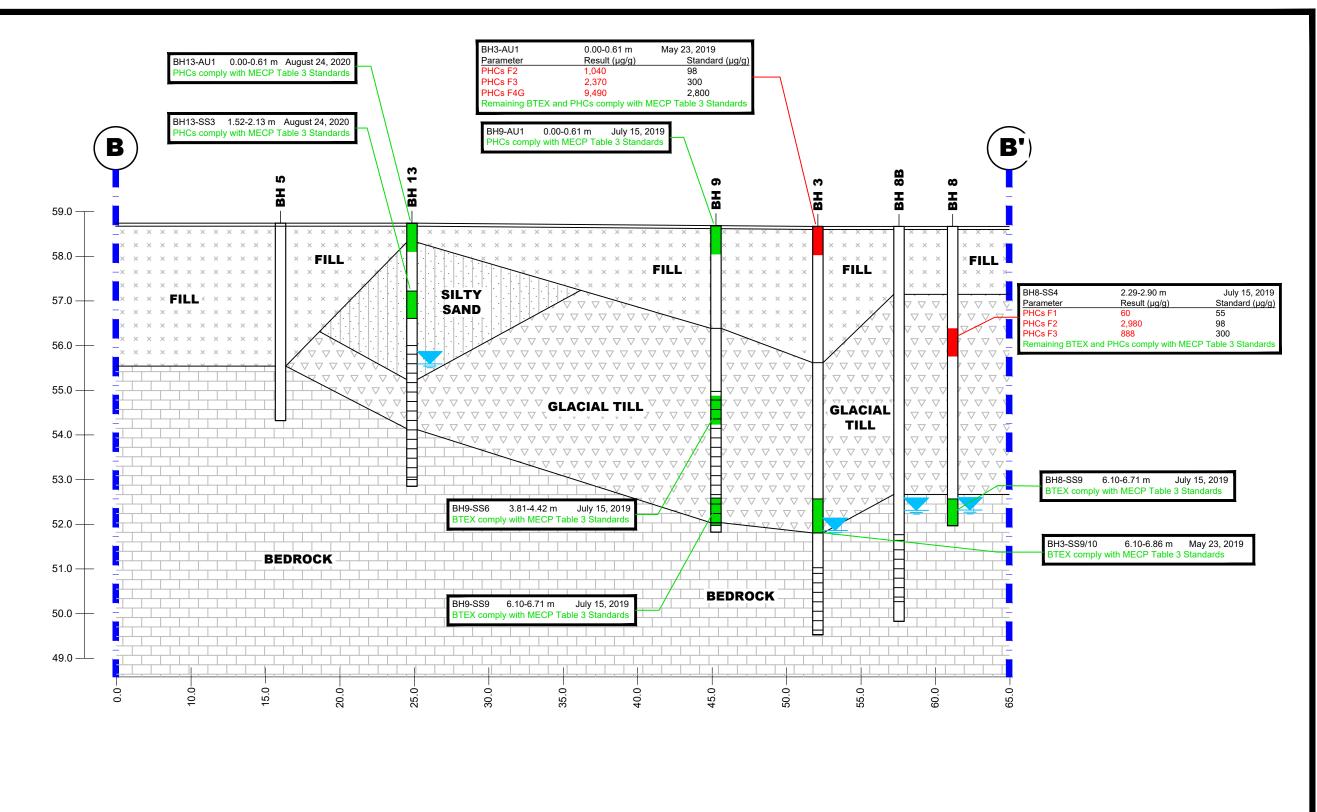
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natorcongroup					MINTO COMMUNITIES INC.
patersongroup					PHASE II - ENVIRONMENTAL SITE ASSESSMENT
consulting engineers					78, 84, 86 & 90 BEECHWOOD AVE. AND 77, 83, 85, 89 & 93 BARRET
					OTTAWA,
154 Colonnade Road South					Title:
Ottawa, Ontario K2E 7J5 Tel: (613) 226-7381 Fax: (613) 226-6344	0				CROSS SECTION B-B' - SOIL (BTEX, PHCs
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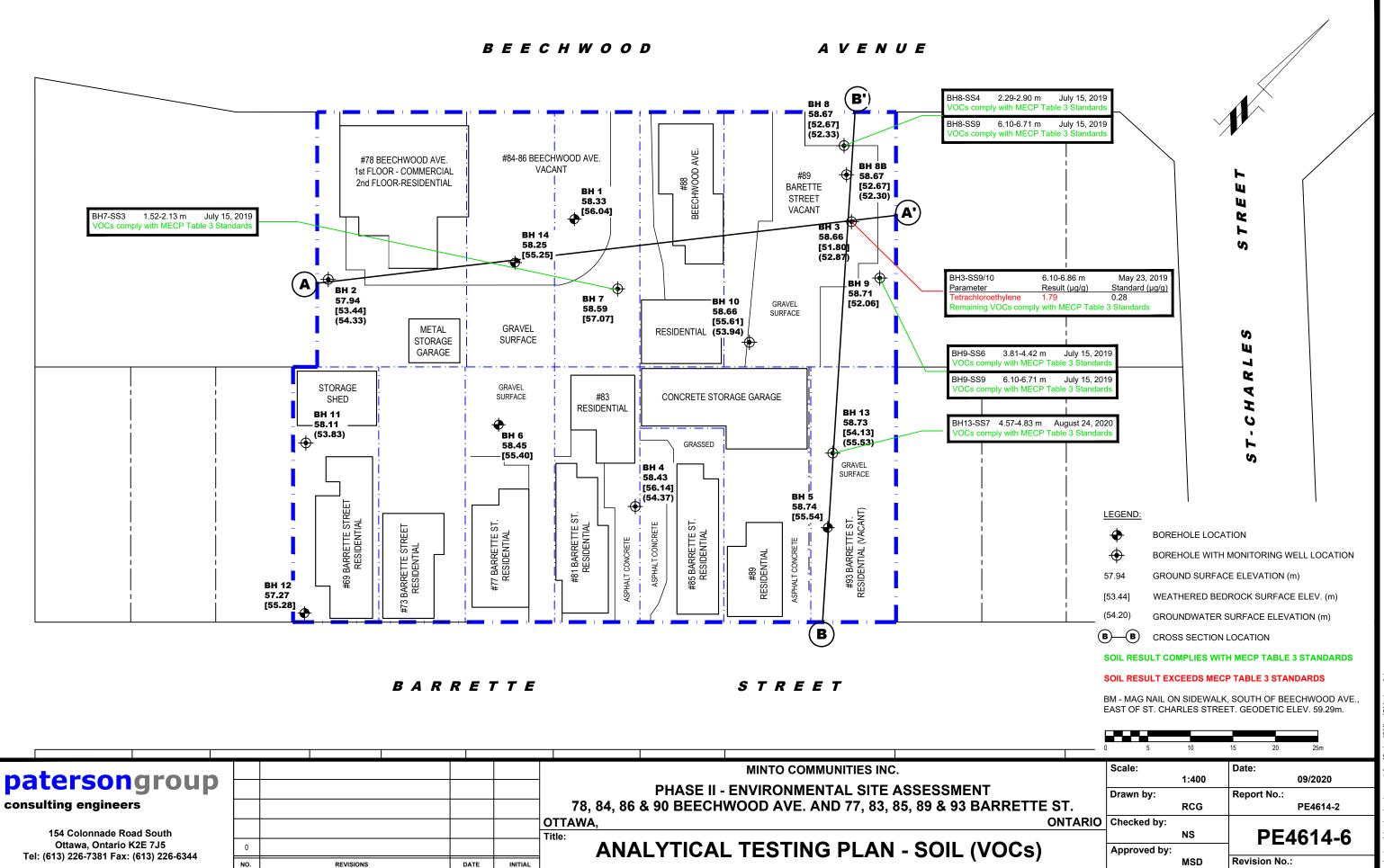
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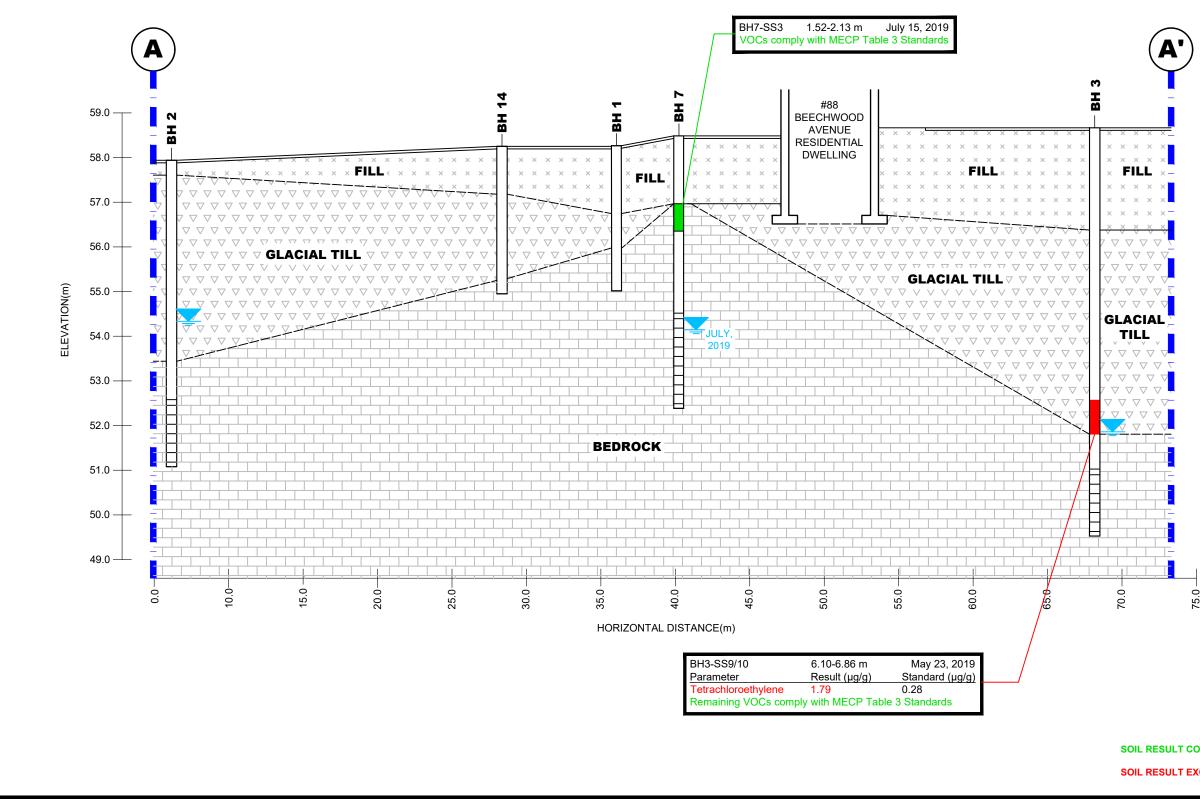
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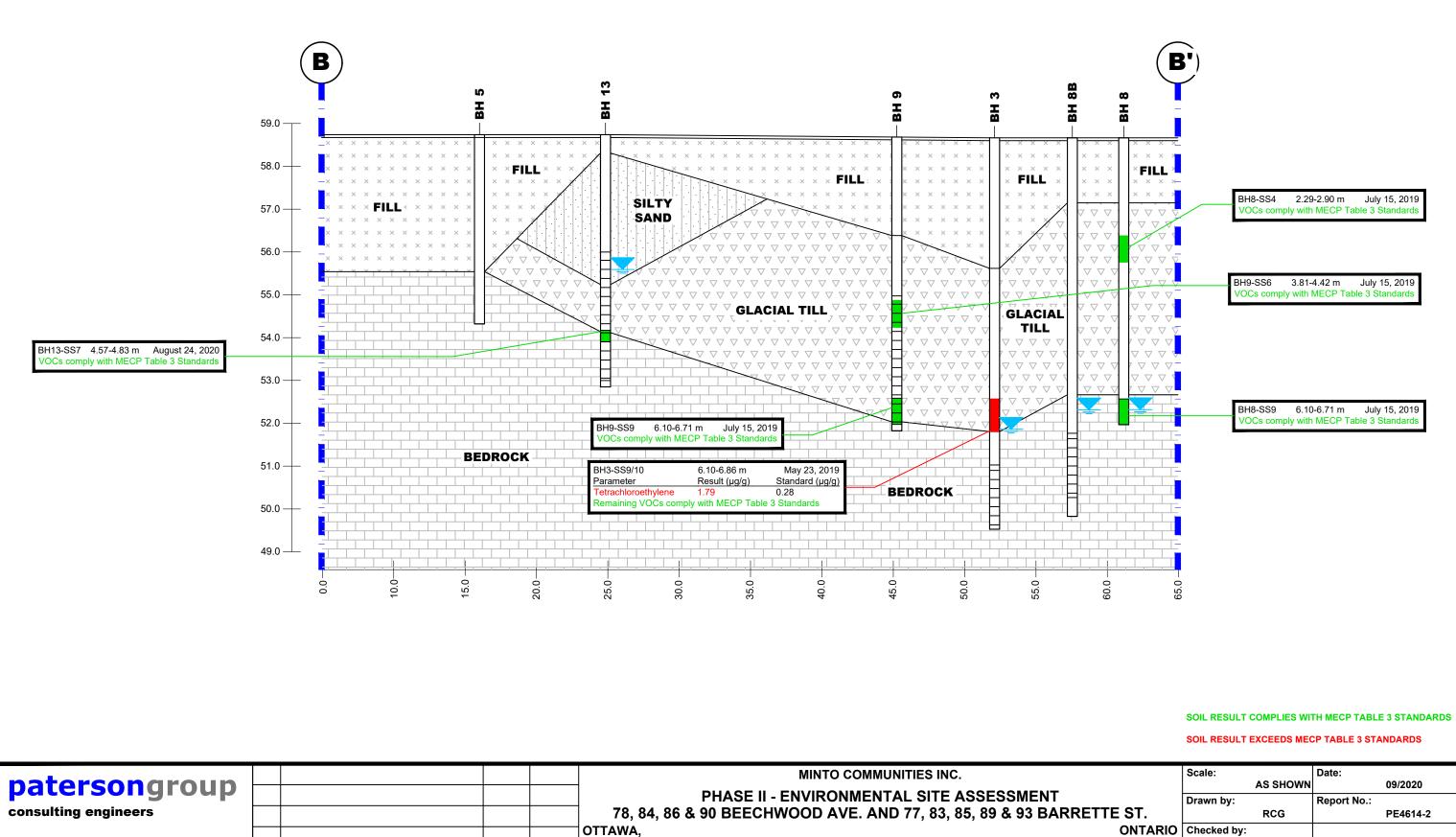




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Ottawa, Ontario K2E 7J5 Tel: (613) 226-7381 Fax: (613) 226-6344	0				CROSS SECTION A-A' - SOIL (VOCs)
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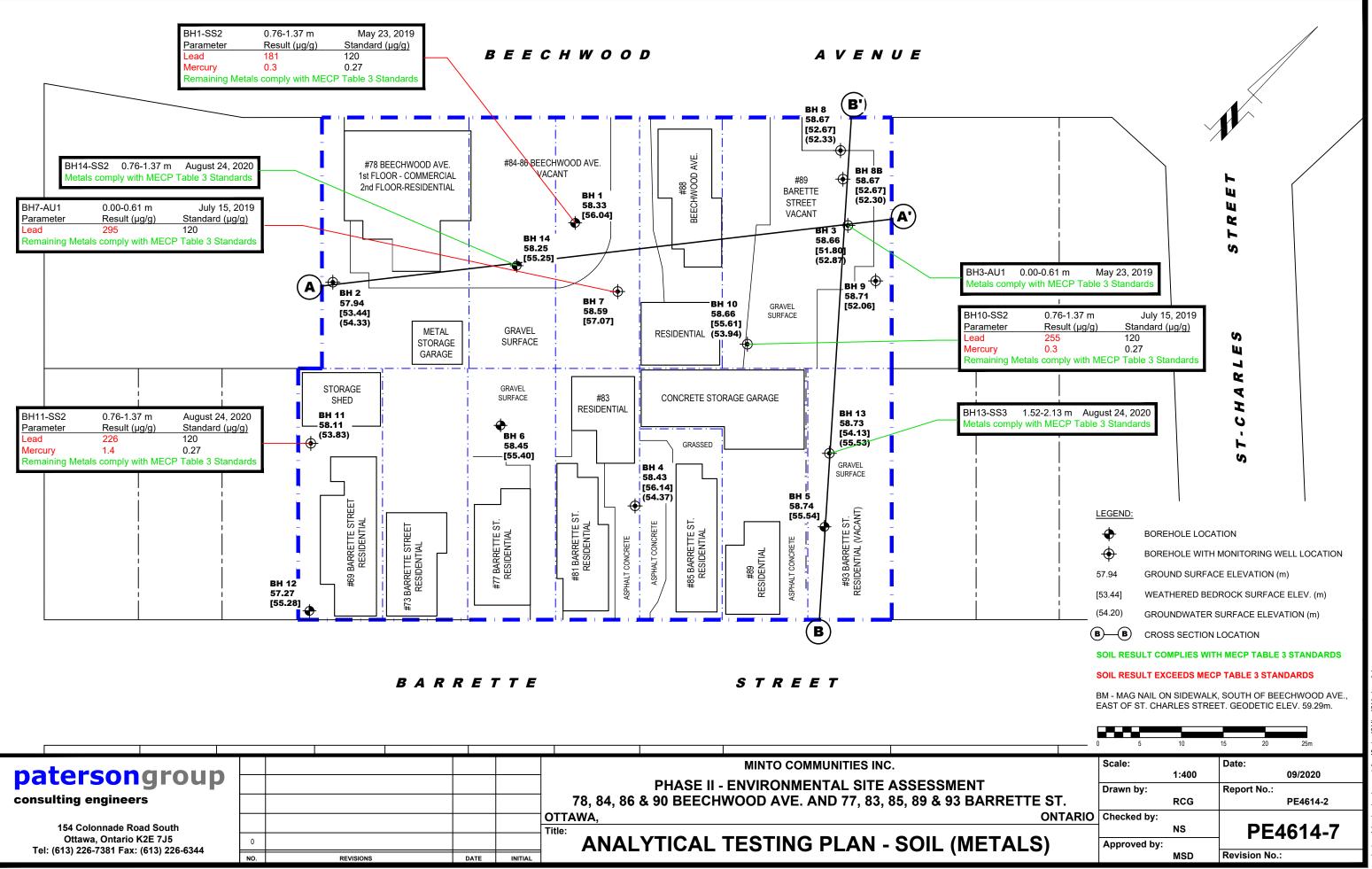
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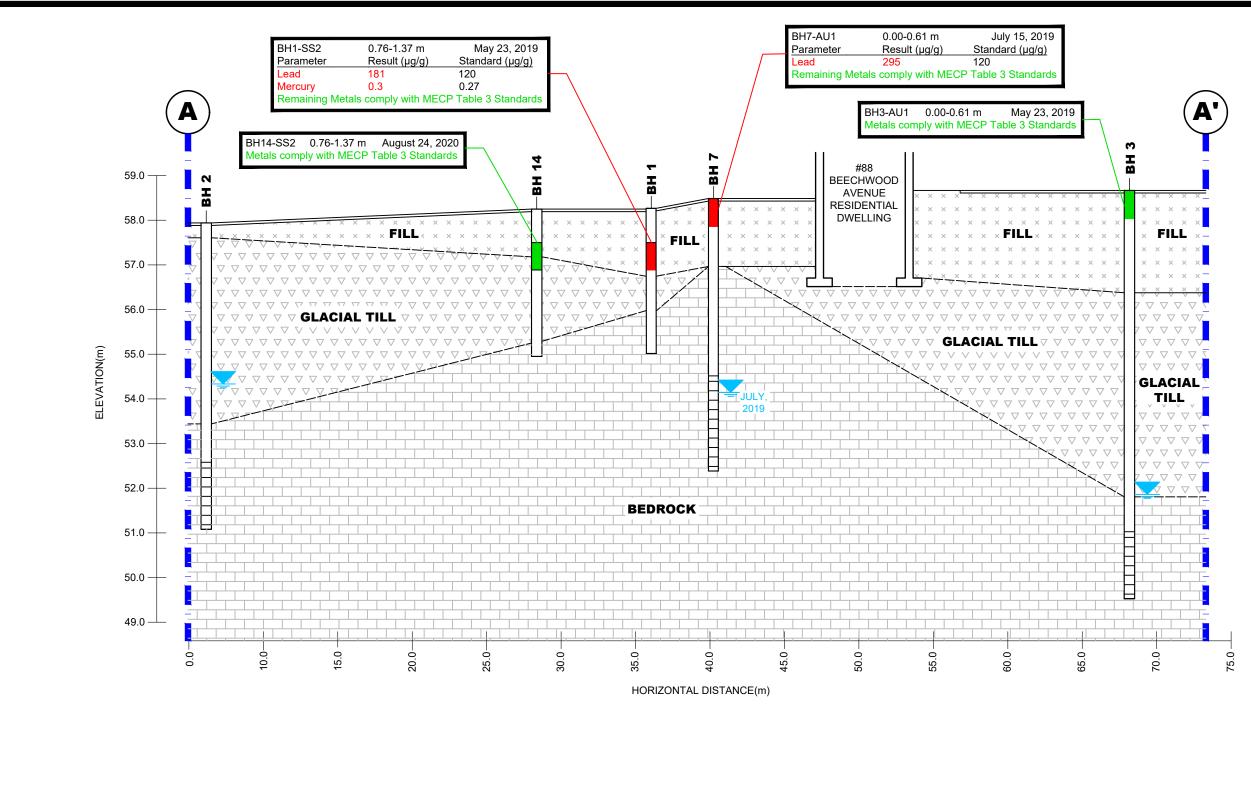
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**CROSS SECTION B-B' - SOIL (VOCs)** 

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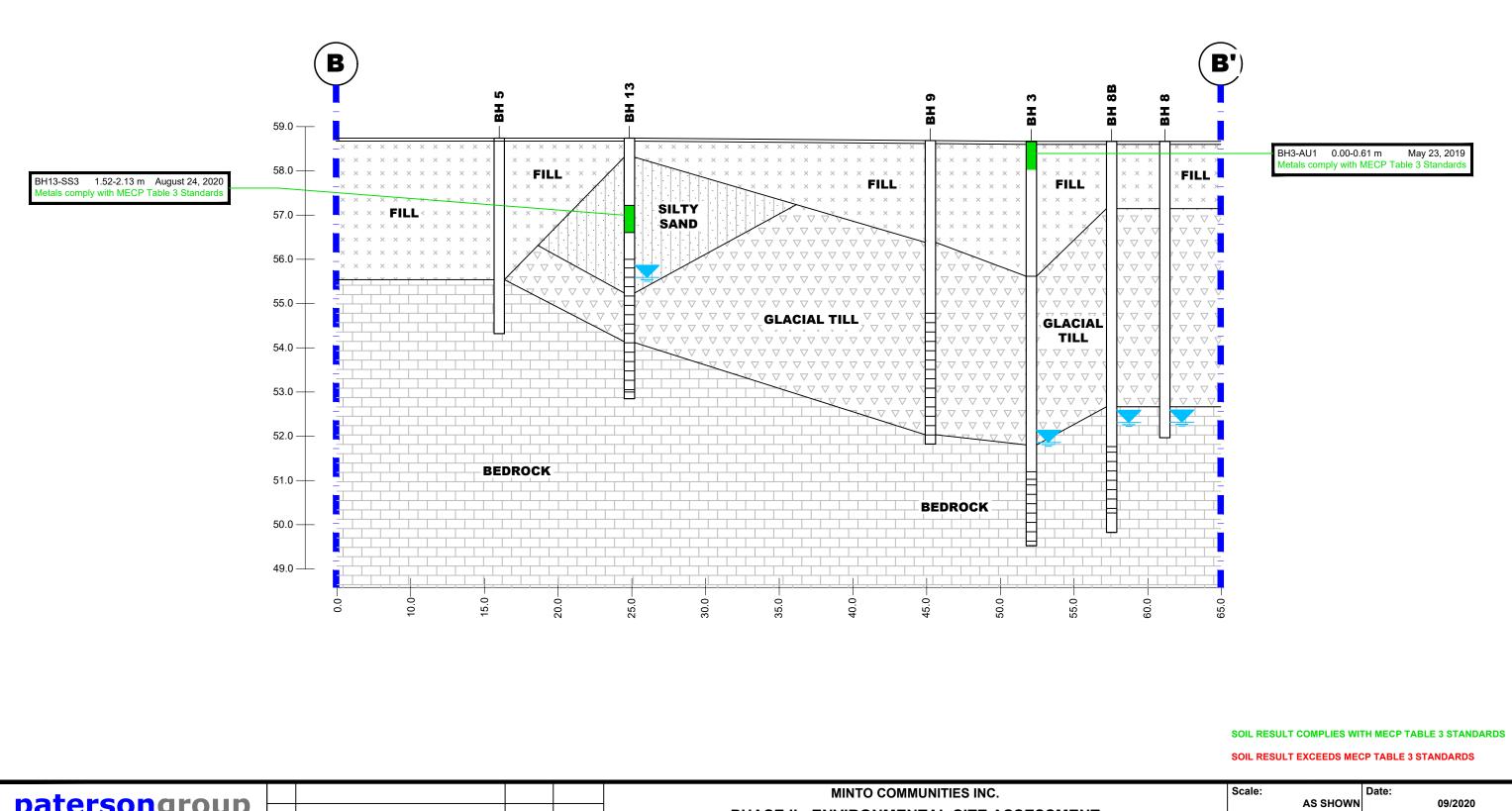
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					OTTAWA,
154 Colonnade Road South					Title:
Ottawa, Ontario K2E 7J5 Tel: (613) 226-7381 Fax: (613) 226-6344	0				CROSS SECTION A-A' - SOIL (METALS)
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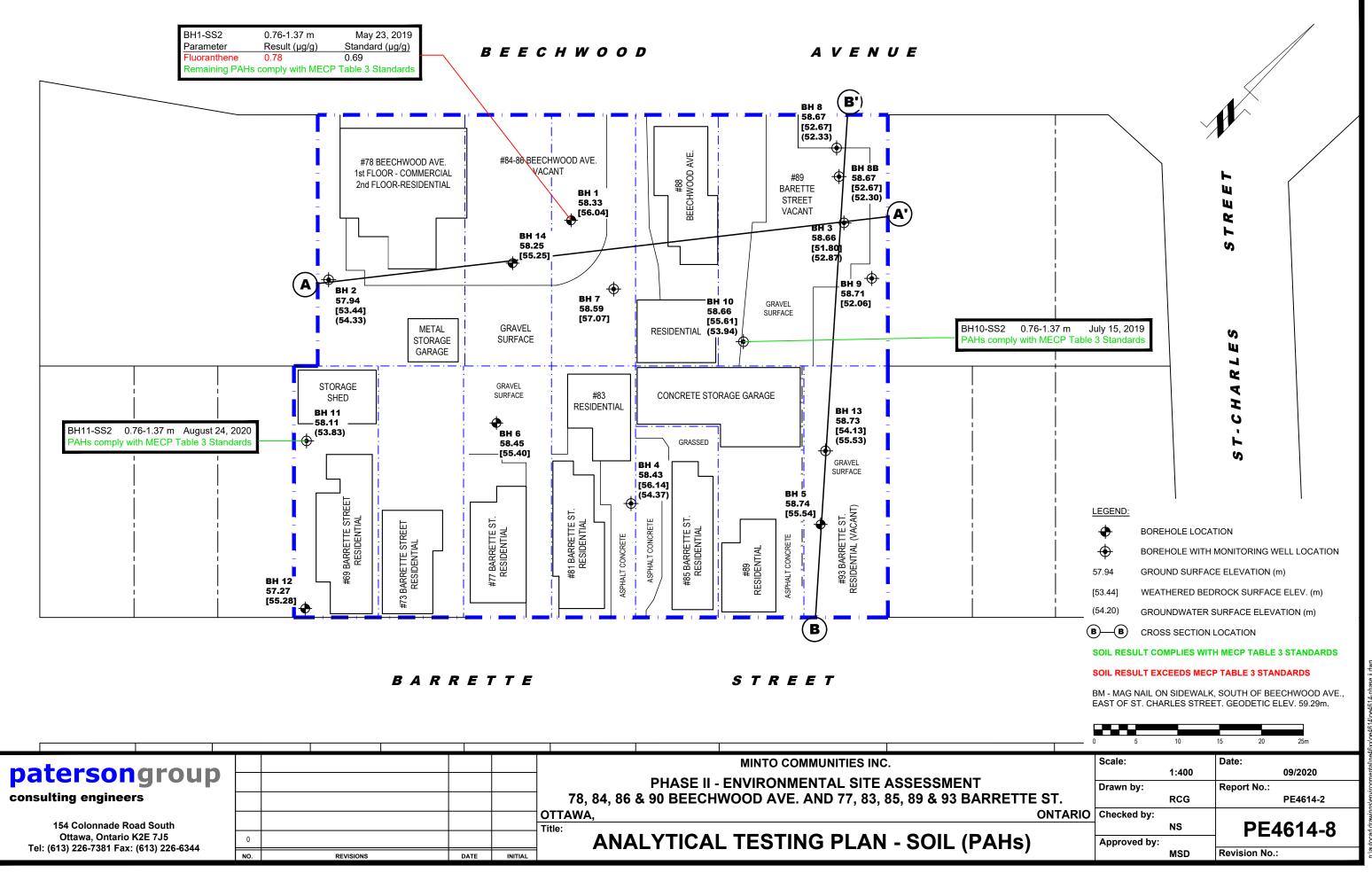
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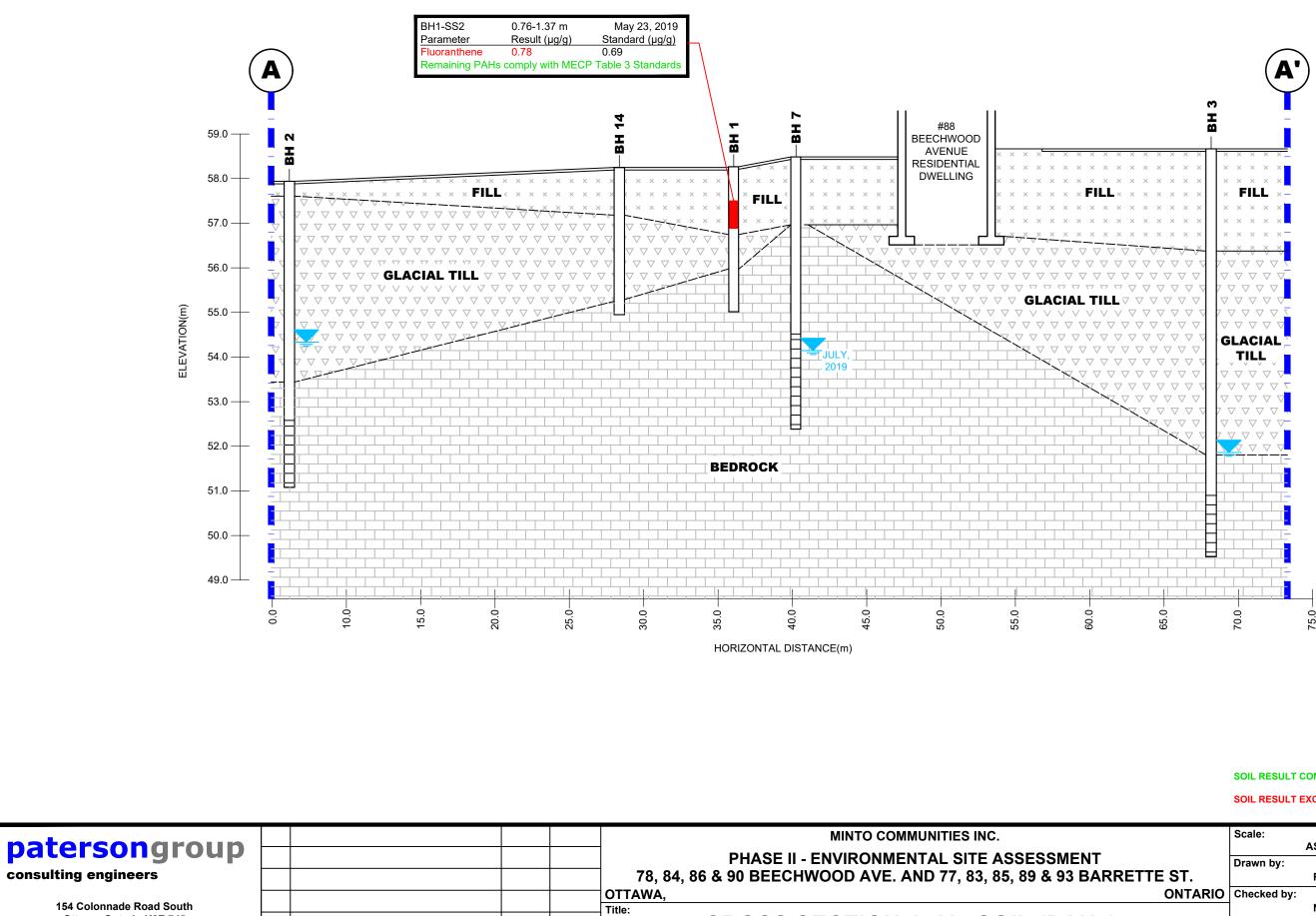
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patersongroup					PHASE II - ENVIRONMENTAL SITE ASSESSMENT
consulting engineers					78, 84, 86 & 90 BEECHWOOD AVE. AND 77, 83, 85, 89 & 93 BARRET
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154 Colonnade Road South					Title:
Ottawa, Ontario K2E 7J5 Tel: (613) 226-7381 Fax: (613) 226-6344	0				CROSS SECTION B-B' - SOIL (METALS)
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## **CROSS SECTION A-A' - SOIL (PAHs)**

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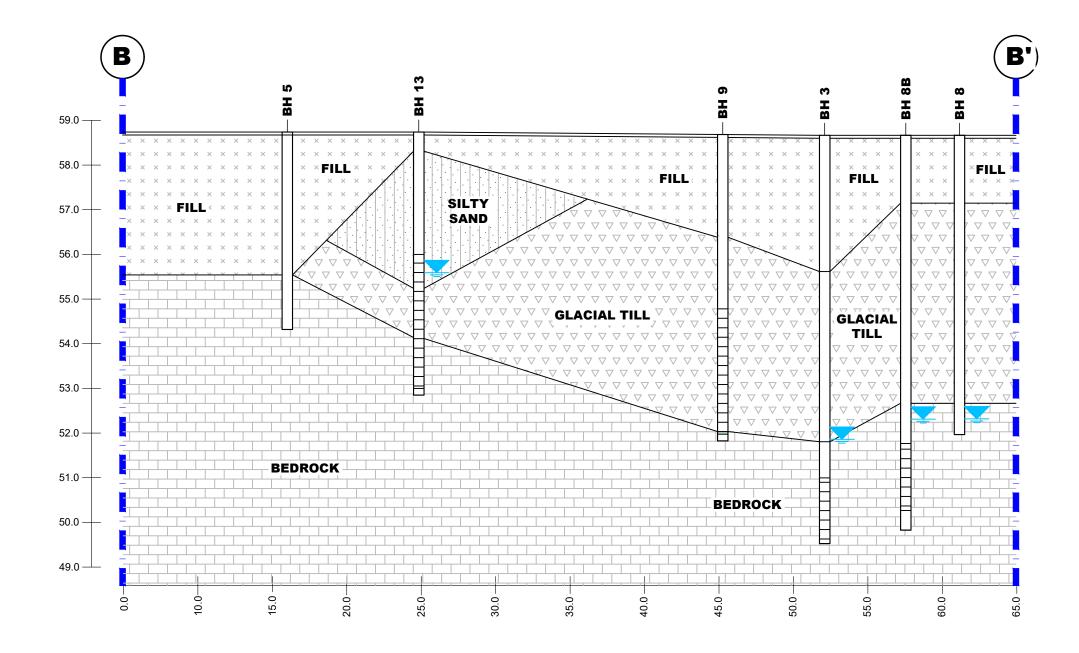
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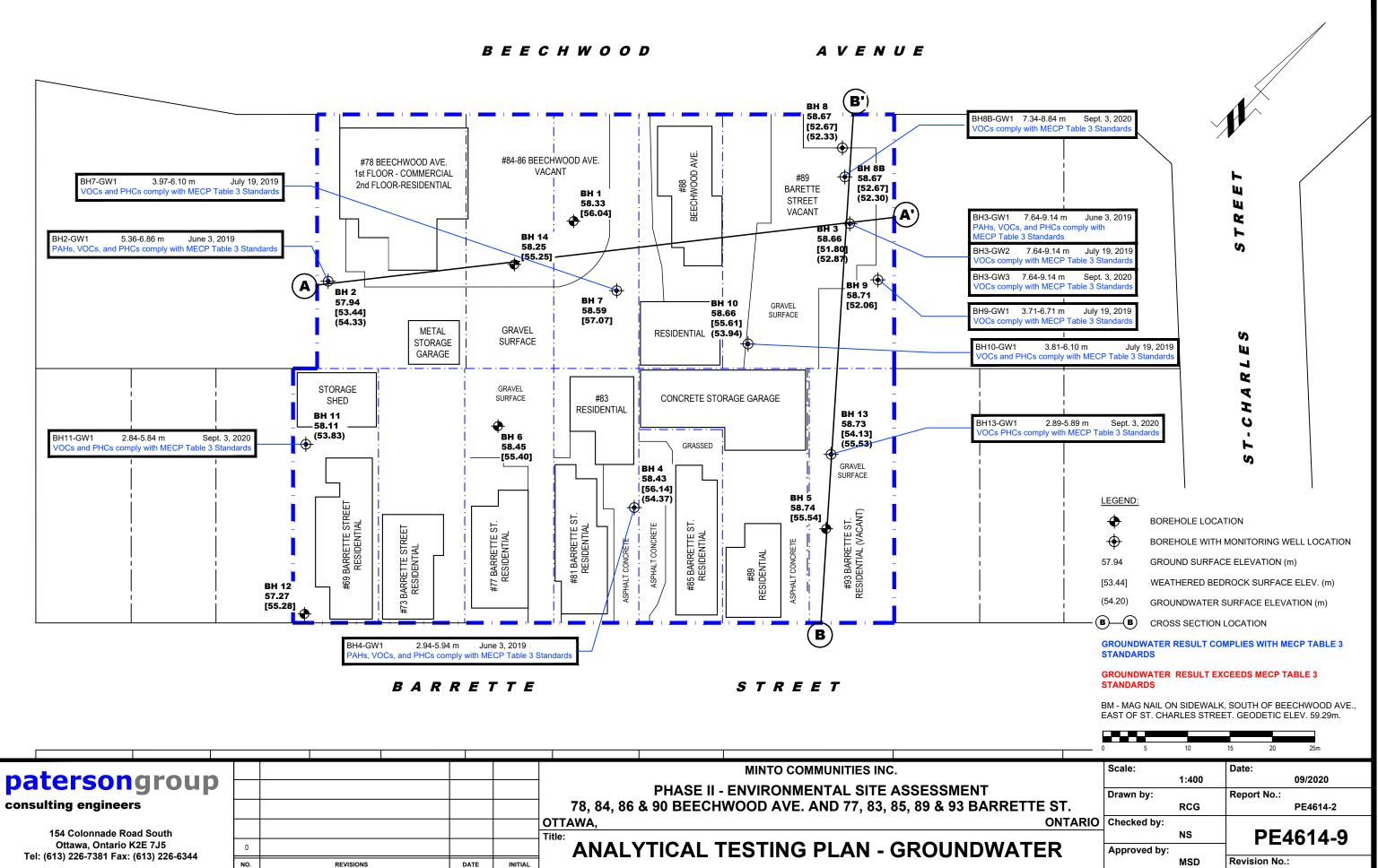
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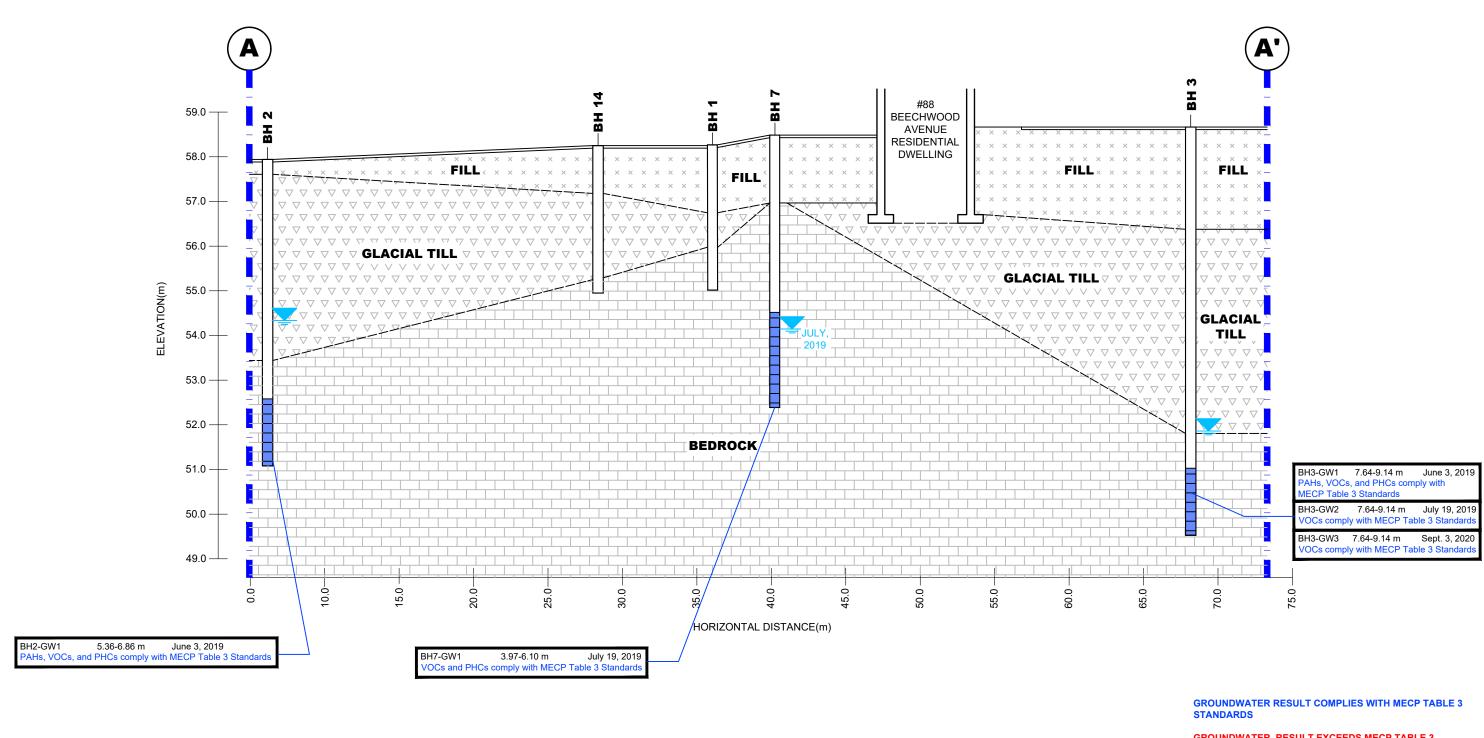


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consulting engineers					PHASE II - ENVIRONMENTAL SITE ASSESSMENT 78, 84, 86 & 90 BEECHWOOD AVE. AND 77, 83, 85, 89 & 93 BARRETTE ST.	Drawn by:	RCG	Report No.: PE4614-2
154 Colonnade Road South					Title:	Checked by:	NS	PE4614-8B
Ottawa, Ontario K2E 7J5 Tel: (613) 226-7381 Fax: (613) 226-6344	0 NO.	REVISIONS	DATE	INITIAL	CROSS SECTION B-B' - SOIL (PAHs)	Approved by		Revision No.:

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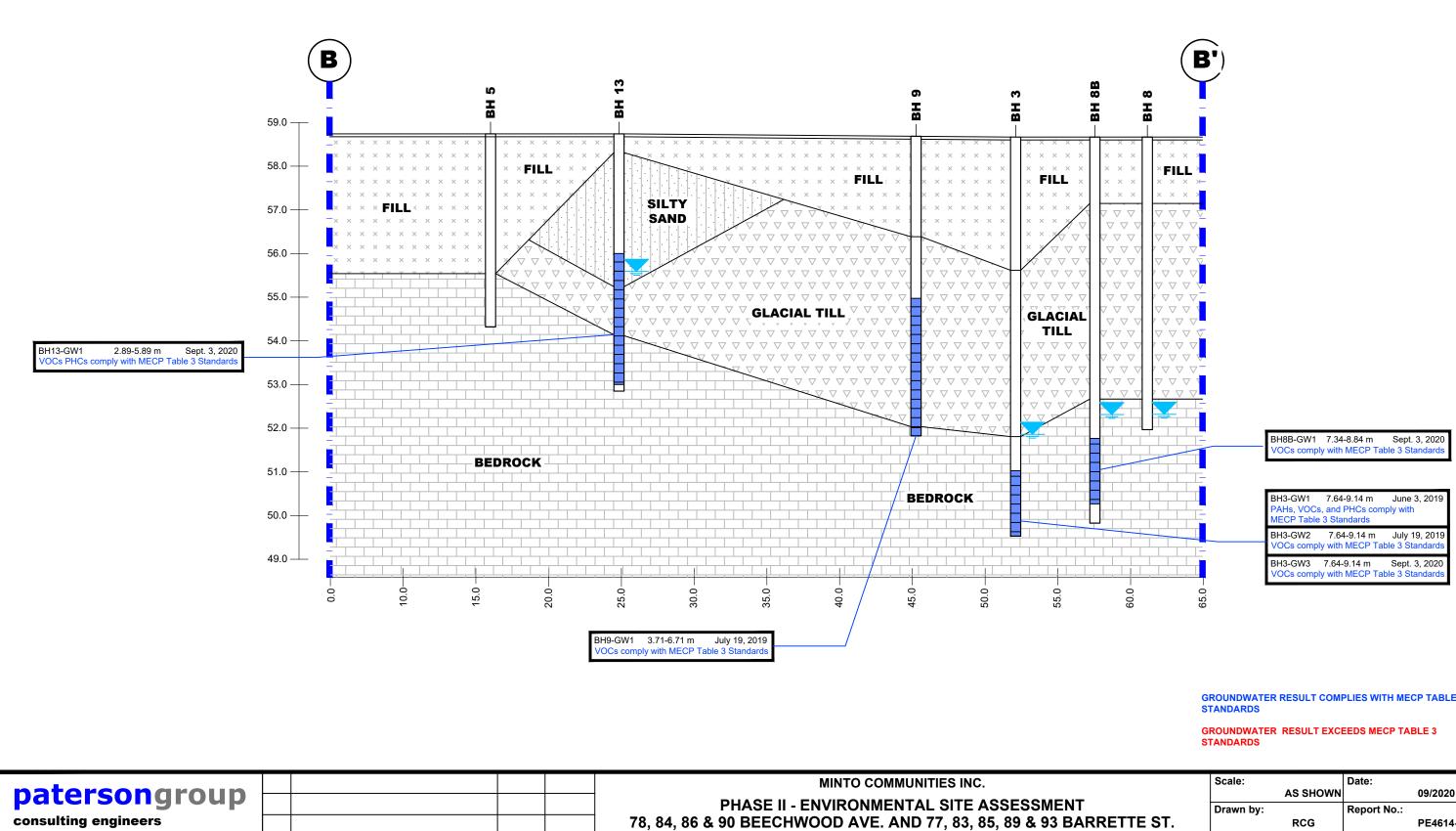




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patersongroup					PHASE II - ENVIRONMENTAL SITE ASSESSMENT
consulting engineers					78, 84, 86 & 90 BEECHWOOD AVE. AND 77, 83, 85, 89 & 93 BARRET
					OTTAWA,
154 Colonnade Road South					Title:
Ottawa, Ontario K2E 7J5 Tel: (613) 226-7381 Fax: (613) 226-6344	0				CROSS SECTION A-A' - GROUNDWATER
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GROUNDWATER RESULT EXCEEDS MECP TABLE 3 **STANDARDS** 

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OTTAWA,

Title:

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Tel: (613) 226-7381 Fax: (613) 226-6344

**CROSS SECTION B-B' - GROUNDWATER** 

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## **APPENDIX 1**

SAMPLING AND ANALYSIS PLAN

SOIL PROFILE AND TEST DATA SHEETS

SYMBOLS AND TERMS

LABORATORY CERTIFICATES OF ANALYSIS

# patersongroup

Geotechnical Engineering

Environmental Engineering

Hydrogeology

Geological Engineering

**Materials Testing** 

**Building Science** 

Archaeological Services

## Sampling & Analysis Plan

Phase II – Environmental Site Assessment 78-90 Beechwood Avenue & 69-93 Barrette Street Ottawa, Ontario

### **Prepared For**

**Minto Communities** 

#### Paterson Group Inc.

Consulting Engineers 154 Colonnade Road South Ottawa (Nepean), Ontario Canada K2E 7J5

Tel: (613) 226-7381 Fax: (613) 226-6344 www.patersongroup.ca May 1, 2019

Report: PE4614-SAP

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	ANALYTICAL TESTING PROGRAM STANDARD OPERATING PROCEEDURES 3.1 Environmental Drilling Procedure 3.2 Monitoring Well Installation Procedure 3.3 Monitoring Well Sampling Procedure QUALITY ASSURANCE/QUALITY CONTROL (QA/QC) DATA QUALITY OBJECTIVES

## 1.0 SAMPLING PROGRAM

Paterson Group Inc. (Paterson) was commissioned by Minto Communities, to conduct a Phase II – Environmental Site Assessment (Phase II ESA) for the properties addressed 78-90 Beechwood Avenue & 69-93 Barrette Street, in the City of Ottawa, Ontario.

Based on the findings of the Phase I ESA, the following subsurface investigation program was developed.

Borehole	Location & Rationale	Proposed Depth & Rationale
BH1	Central portion of subject site; to assess for potential impacts resulting from a former landfill site.	3-6 m; for general coverage purposes.
BH2	Western portion of subject site; to assess for potential impacts resulting from a former off-site retail fuel outlet.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH3	Northeastern portion of subject site: to assess for potential impacts resulting from an off-site dry-cleaners.	6-10 m; to intercept the groundwater table for the purpose of installing a deeper screen in the shale bedrock.
BH4	Southern portion of subject site; to assess for potential impacts resulting from a former on-site autobody shop.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH5	Southeastern portion of subject site; to assess for potential impacts resulting from a former on-site autobody shop.	3-6 m; for general coverage purposes.
BH6	Central portion of subject site; to assess for potential impacts resulting from a former landfill site.	3-6 m; for general coverage purposes.
BH7	Central portion of subject site; to assess for potential impacts resulting from a former landfill site.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH8	Northeastern portion of subject site: to assess for potential impacts resulting from an off-site dry-cleaners.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH8B	Northeastern portion of subject site: to assess for potential impacts resulting from an off-site dry-cleaners.	6-10 m; to intercept the groundwater table for the purpose of installing a deeper screen in the shale bedrock.
BH9	Northeastern portion of subject site: to assess for potential impacts resulting from an off-site dry-cleaners.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH10	Eastern portion of subject site; to assess for potential impacts resulting from a former on-site autobody shop.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH11	Western portion of subject site; to assess for potential impacts resulting from a former off-site retail fuel outlet.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH12	Southwestern portion of subject site; to assess for potential impacts resulting from a former landfill site.	3-6 m; for general coverage purposes.
BH13	Eastern portion of subject site; to assess for potential impacts resulting from a former on-site autobody shop.	6-10 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH14	Central portion of subject site; to assess for potential impacts resulting from a former landfill site.	3-6 m; for general coverage purposes.

Borehole locations are shown on Drawing PE4614-4 – Test Hole Location Plan, appended to the main report.

At each borehole, split-spoon samples of the overburden soils will be obtained at 0.76 m (2'6") intervals until practical refusal to augering. All soil samples will be retained, and samples will be selected for submission following a preliminary screening analysis.

Following the borehole drilling, groundwater monitoring wells will be installed in BH2, BH3, BH4, BH7, BH8, BH8B, BH9, BH10, BH11, and BH13 for the collection of groundwater samples.

## 2.0 ANALYTICAL TESTING PROGRAM

The analytical testing program for soil at the subject site is based on the following general considerations:

- □ At least one sample from each borehole should be submitted, in order to delineate the horizontal extent of contamination across the site.
- □ At least one sample from each stratigraphic unit should be submitted, in order to delineate the vertical extent of contamination at the site.
- □ In boreholes where there is visual or olfactory evidence of contamination, or where organic vapour meter or photoionization detector readings indicate the presence of contamination, the 'worst-case' sample from each borehole should be submitted for comparison with MECP site condition standards.
- In boreholes with evidence of contamination as described above, a sample should be submitted from the stratigraphic unit below the 'worst-case' sample to determine whether the contaminant(s) have migrated downward.
- Parameters analyzed should be consistent with the Contaminants of Potential Concern identified in the Phase I ESA.

The analytical testing program for soil at the subject site is based on the following general considerations:

Groundwater monitoring wells should be installed in all boreholes with visual or olfactory evidence of soil contamination, in stratigraphic units where soil contamination was encountered, where those stratigraphic units are at or below the water table (i.e. a water sample can be obtained).

- Groundwater monitoring well screens should straddle the water table at sites where the contaminants of concern are suspected to be LNAPLs.
- ☐ At least one groundwater monitoring well should be installed in a stratigraphic unit below the suspected contamination, where said stratigraphic unit is water-bearing.
- Parameters analyzed should be consistent with the Contaminants of Concern identified in the Phase I ESA and with the contaminants identified in the soil samples.

## 3.0 STANDARD OPERATING PROCEEDURES

### 3.1 Environmental Drilling Procedure

### Purpose

The purpose of environmental boreholes is to identify and/or delineate contamination within the soil and/or to install groundwater monitoring wells in order to identify contamination within the groundwater.

### Equipment

The following is a list of equipment that is in addition to regular drilling equipment stated in the geotechnical drilling SOP:

- Glass soil sample jars
- two buckets
- □ cleaning brush (toilet brush works well)
- dish detergent
- methyl hydrate
- d water (if not available on site water jugs available in trailer)
- □ latex or nitrile gloves (depending on suspected contaminant)
- RKI Eagle organic vapour meter or MiniRae photoionization detector (depending on contamination suspected)

### **Determining Borehole Locations**

If conditions on site are not as suspected, and planned borehole locations cannot be drilled, **call the office to discuss**. Alternative borehole locations will be determined in conversation with the field technician and supervising engineer. After drilling is completed a plan with the borehole locations must be provided. Distances and orientations of boreholes with respect to site features (buildings, roadways, etc.) must be provided. Distances should be measured using a measuring tape or wheel rather than paced off. Ground surface elevations at each borehole should be surveyed relative to a geodetic benchmark, if one is available, or a temporary site benchmark which can be tied in at a later date if necessary.

### **Drilling Procedure**

The actual drilling procedure for environmental boreholes is the same as geotechnical boreholes (see SOP for drilling and sampling) with a few exceptions as follows:

- Continuous split spoon samples (every 0.6 m or 2') or semi-continuous (every 0.76 m or 2'6") are required.
- □ Make sure samples are well sealed in plastic bags with no holes prior to screening and are kept cool but unfrozen.
- □ If sampling for VOCs, BTEX, or PHCs F<sub>1</sub>, a soil core from each soil sample, which may be analyzed, must be taken and placed in the laboratory-provided methanol vial.
- □ Note all and any odours or discolouration of samples.
- □ Split spoon samplers must be washed between samples.
- If obvious contamination is encountered, continue sampling until vertical extent of contamination is delineated.
- As a general rule, environmental boreholes should be deep enough to intercept the groundwater table (unless this is impossible/impractical - call project manager to discuss).
- If at all possible, soil samples should be submitted to a preliminary screening procedure on site, either using a RKI Eagle, PID, etc. depending on type of suspected contamination.

### Spoon Washing Procedure

All sampling equipment (spilt spoons, etc.) must be washed between samples in order to prevent cross contamination of soil samples.

- □ Obtain two buckets of water (preferably hot if available)
- Add a small amount of dish soap to one bucket
- □ Scrub spoons with brush in soapy water, inside and out, including tip
- **D** Rinse in clean water

- □ Apply a small amount of methyl hydrate to the inside of the spoon. (A spray bottle or water bottle with a small hole in the cap works well)
- □ Allow to dry (takes seconds)
- □ Rinse with distilled water, a spray bottle works well.

The methyl hydrate eliminates any soap residue that may be on the spoon and is especially important when dealing with suspected VOCs.

### Screening Procedure

The RKI Eagle is used to screen most soil samples, particularly where petroleum hydrocarbon contamination is suspected. The MiniRae is used when VOCs are suspected, however it also can be useful for detecting petroleum. These tools are for screening purposes only and cannot be used in place of laboratory testing. Vapour results obtained from the RKI Eagle and the PID are relative and must be interpreted.

Screening equipment should be calibrated on an approximately monthly basis, more frequently if heavily used.

- Samples should be brought to room temperature; this is specifically important in colder weather. Soil must not be frozen.
- □ Turn instrument on and allow to come to zero calibrate if necessary
- If using RKI Eagle, ensure instrument is in methane elimination mode unless otherwise directed.
- Ensure measurement units are ppm (parts per million) initially. RKI Eagle will automatically switch to %LEL (lower explosive limit) if higher concentrations are encountered.
- Break up large lumps of soil in the sample bag, taking care not to puncture bag.
- □ Insert probe into soil bag, creating a seal with your hand around the opening.
- Gently manipulate soil in bag while observing instrument readings.
- Record the highest value obtained in the first 15 to 25 seconds
- Make sure to indicate scale (ppm or LEL); also note which instrument was used (RKI Eagle 1 or 2, or MiniRae).
- □ Jar samples and refrigerate as per Sampling and Analysis Plan.

### 3.2 Monitoring Well Installation Procedure

### Equipment

- ☐ 5' x 2" threaded sections of Schedule 40 PVC slotted well screen (5' x 1 ¼" if installing in cored hole in bedrock)
- □ 5' x 2" threaded sections of Schedule 40 PVC riser pipe (5' x 1 ¼" if installing in cored hole in bedrock)
- □ Threaded end-cap
- □ Slip-cap or J-plug
- □ Asphalt cold patch or concrete
- □ Silica Sand
- Bentonite chips (Holeplug)
- □ Steel flushmount casing

### Procedure

- Drill borehole to required depth, using drilling and sampling procedures described above.
- If borehole is deeper than required monitoring well, backfill with bentonite chips to required depth. This should only be done on wells where contamination is not suspected, in order to prevent downward migration of contamination.
- □ Only one monitoring well should be installed per borehole.
- Monitoring wells should not be screened across more than one stratigraphic unit to prevent potential migration of contaminants between units.
- Where LNAPLs are the suspected contaminants of concern, monitoring wells should be screened straddling the water table in order to capture any free product floating on top of the water table.
- Thread the end cap onto a section of screen. Thread second section of screen if required. Thread risers onto screen. Lower into borehole to required depth. Ensure slip-cap or J-plug is inserted to prevent backfill materials entering well.
- □ As drillers remove augers, backfill borehole annulus with silica sand until the level of sand is approximately 0.3 m above the top of the screen.
- Backfill with holeplug until at least 0.3 m of holeplug is present above the top of the silica sand.
- Backfill remainder of borehole with holeplug or with auger cuttings (if contamination is not suspected).
- Install flushmount casing. Seal space between flushmount and borehole annulus with concrete, cold patch, or holeplug to match surrounding ground surface.

### 3.3 Monitoring Well Sampling Procedure

### Equipment

- □ Water level metre or interface probe on hydrocarbon/LNAPL sites
- Spray bottles containing water and methanol to clean water level tape or interface probe
- Peristaltic pump
- D Polyethylene tubing for peristaltic pump
- □ Flexible tubing for peristaltic pump
- □ Latex or nitrile gloves (depending on suspected contaminant)
- □ Allen keys and/or 9/16" socket wrench to remove well caps
- Graduated bucket with volume measurements
- D pH/Temperature/Conductivity combo pen
- □ Laboratory-supplied sample bottles

### Sampling Procedure

- □ Locate well and use socket wrench or Allan key to open metal flush mount protector cap. Remove plastic well cap.
- Measure water level, with respect to existing ground surface, using water level meter or interface probe. If using interface probe on suspected NAPL site, measure the thickness of free product.
- □ Measure total depth of well.
- Clean water level tape or interface probe using methanol and water. Change gloves between wells.
- □ Calculate volume of standing water within well and record.
- Insert polyethylene tubing into well and attach to peristaltic pump. Turn on peristaltic pump and purge into graduated bucket. Purge at least three well volumes of water from the well. Measure and record field chemistry. Continue to purge, measuring field chemistry after every well volume purged, until appearance or field chemistry stabilizes.
- Note appearance of purge water, including colour, opacity (clear, cloudy, silty), sheen, presence of LNAPL, and odour. Note any other unusual features (particulate matter, effervescence (bubbling) of dissolved gas, etc.).
- Fill required sample bottles. If sampling for metals, attach 75-micron filter to discharge tube and filter metals sample. If sampling for VOCs, use low flow rate to ensure continuous stream of non-turbulent flow into sample bottles. Ensure no headspace is present in VOC vials.
- □ Replace well cap and flushmount casing cap.

## 4.0 QUALITY ASSURANCE/QUALITY CONTROL (QA/QC)

The QA/QC program for this Phase II ESA is as follows:

- All non-dedicated sampling equipment (split spoons) will be decontaminated according to the SOPs listed above.
- □ All groundwater sampling equipment is dedicated (polyethylene and flexible peristaltic tubing is replaced for each well).
- Where groundwater samples are to be analyzed for VOCs, one laboratoryprovided trip blank will be submitted for analysis with every laboratory submission.
- Approximately one (1) field duplicate will be submitted for every ten (10) samples submitted for laboratory analysis. A minimum of one (1) field duplicate per project will be submitted. Field duplicates will be submitted for soil and groundwater samples
- Where combo pens are used to measure field chemistry, they will be calibrated on an approximately monthly basis, according to frequency of use.

## 5.0 DATA QUALITY OBJECTIVES

The purpose of setting data quality objectives (DQOs) is to ensure that the level of uncertainty in data collected during the Phase II ESA is low enough that decision-making is not affected, and that the overall objectives of the investigation are met.

The quality of data is assessed by comparing field duplicates with original samples. If the relative percent difference (RPD) between the duplicate and the sample is within 20%, the data are considered to be of sufficient quality so as not to affect decision-making. The RPD is calculated as follows:

$$RPD = \left| \frac{x_1 - x_2}{(x_1 + x_2)/2} \right| \times 100\%$$

Where  $x_1$  is the concentration of a given parameter in an original sample and  $x_2$  is the concentration of that same parameter in the field duplicate sample.

For the purpose of calculating the RPD, it is desirable to select field duplicates from samples for which parameters are present in concentrations above laboratory detection limits, i.e. samples which are expected to be contaminated. If parameters are below laboratory detection limits for selected samples or duplicates, the RPD may be calculated using a concentration equal to one half the laboratory detection limit.

It is also important to consider data quality in the overall context of the project. For example, if the DQOs are not met for a given sample, yet the concentrations of contaminants in both the sample and the duplicate exceed the MOE site remediation standards by a large margin, the decision-making usefulness of the sample may not be considered to be impaired. The proximity of other samples which meet the DQOs must also be considered in developing the Phase II Conceptual Site Model; often there are enough data available to produce a reliable Phase II Conceptual Site Model even if DQOs are not met for certain individual samples.

These considerations are discussed in the body of the report.

Sampling & Analysis Plan 78-90 Beechwood Avenue & 69-93 Barrette Street Ottawa, Ontario

#### 6.0 PHYSICAL IMPEDIMENTS

Ōttawa

Kingston

Physical impediments to the Sampling and Analysis plan may include:

- The location of underground utilities
- Poor recovery of split-spoon soil samples
- Insufficient groundwater volume for groundwater samples
- □ Breakage of sampling containers following sampling or while in transit to the laboratory
- □ Elevated detection limits due to matrix interference (generally related to soil colour or presence of organic material)
- Elevated detection limits due to high concentrations of certain parameters, necessitating dilution of samples in laboratory
- Drill rig breakdowns
- Winter conditions
- Other site-specific impediments

Site-specific impediments to the Sampling and Analysis plan are discussed in the body of the Phase II ESA report.

## patersongroup Consulting Engineers

## SOIL PROFILE AND TEST DATA

**Phase II - Environmental Site Assessment** 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

DATUM
-------

BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. Geodetic elevation = 59.29m. REMARKS BORINGS BY CME-55 Low Clearance Drill DATE May 23, 2019

HOLE NO. **BH 1** 

**PE4614** 

FILE NO.

SOIL DESCRIPTION			SAMPLE			DEPTH	ELEV.	Photo Ionization Detector     Volatile Organic Rdg. (ppm)			
GROUND SURFACE	STRATA PLOT	ТҮРЕ	NUMBER	°∂ RECOVERY	N VALUE or RQD	(m)	(m)	Photo Ionization Detector         ● Volatile Organic Rdg. (ppm)         ○ Lower Explosive Limit %         20       40       60       80			
Ell I · Crushed stone		AU	1			0-	-58.33	•			
FILL: Brown silty sand, some gravel		ss	2	50	19	1-	-57.33	•			
GLACIAL TILL: Very dense, prown clayey silty sand with gravel and shale	9	ss	3	62	52	2-	-56.33	•			
→→→→→→→		ss	4	100	50+			•			
3.2	5	∑ ss	5	100	50+	3-	-55.33				
Practical refusal to augering at 3.25m depth											
BH dry upon completion)											
								100 200 300 400 500 <b>RKI Eagle Rdg. (ppm)</b> ▲ Full Gas Resp. △ Methane Elim.			

6.86

## SOIL PROFILE AND TEST DATA

**PE4614** 

**BH 2** 

Phase II - Environmental Site Assessment 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

DATUM

End of Borehole

6.86m depth.

Practical refusal to augering at

(GWL @ 3.61m - Sept. 3, 2020)

BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. FILE NO. Geodetic elevation = 59.29m. REMARKS HOLE NO. BORINGS BY CME-55 Low Clearance Drill DATE May 23, 2019 SAMPLE **Photo Ionization Detector** PLOT DEPTH ELEV. SOIL DESCRIPTION Volatile Organic Rdg. (ppm) (m) (m) RECOVERY STRATA VALUE r RQD NUMBER TYPE o/0  $\bigcirc$ N OF **GROUND SURFACE** 20 40 0+57.94FILL: Crushed stone, some sand 0.33 AU 1 1+56.94

Monitoring Well Construction Lower Explosive Limit % 80 60 <u>ներերեր</u> 2 SS 79 14 SS 3 100 35 2+55.94 **GLACIAL TILL:** Brown clayey silt with sand and shaley gravel SS 4 100 50 +3+54.94SS 5 50 +100 4+53.94 SS 6 100 82 4.50 7 SS 100 50 +5+52.94 ≤ SS 8 100 50 +Weathered black shale **BEDROCK** 

6+51.94

100

200

RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

300

400

500

## SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 78, 84, 86 and 90 Beechwood Avenue

d South Ottawa Ontario K2F 7.15 154 ( do Do

154 Colonnade Road South, Ottawa, O	mano		15		83	, 85, 89, 9	3 Barret	te Street, (	Ottawa, Ont	ario	
DATUM BM - Mag nail on sidewa Geodetic elevation = 59. REMARKS	lk, sou 29m.	ith of E	Beech	wood	Ave.,	east of S	t. Charle	s Street.	FILE NO.	PE4614	ļ
BORINGS BY CME-55 Low Clearance	e Drill			D	ATE	May 23, 2	2019		HOLE NO.		
SOIL DESCRIPTION	PLOT		SAN	IPLE		DEPTH	ELEV.		onization De		g Well ction
	STRATA	ТҮРЕ	IUMBER	NUMBER % RECOVERY	N VALUE or RQD	(m)	(m)	○ Lowe	r Explosive	Limit %	Monitoring Well Construction
GROUND SURFACE	01		2	RE	N	0-	-58.66	20	40 60	80	Σ
FILL: Crushed stone, some sand 0.2	23 🔀	AU	1				-00.00	•			
FILL: Brown silty sand, trace gravel, brick and concrete		ss	2	92	7	1-	-57.66	•			
		ss	3	92	14	2-	-56.66				
FILL: Brown silty sand, some gravel, trace clay	<u>29 ~~ ~</u>	ss	4	91	88						
<u>3.0</u>	05	ss	5	80	50+	3-	-55.66	•	• • • • • • • • • • • • • • • • • • •		
		ss	6	92	70	4-	-54.66	•			
<b>GLACIAL TILL:</b> Brown sand silty/silty sand with shale		ss	7	100	50+	5-	-53.66	•			
		∦ ss	8	100	50+	6-	-52.66	•			¥
			9	100	50+						չյուներիներին։ Տերերերիներին
6.8	36	x SS	10	100	50+	7-	-51.66				<u>իրիրին։</u>
Weathered black shale <b>BEDROCK</b>						8-	-50.66				
9.14	4					9-	-49.66				
End of Borehole											
(GWL @ 5.79m - Sept. 3, 2020)								100	200 300	400 50	00
								RKI E	Eagle Rdg. (  as Resp. △ Me	ppm)	

# patersongroup Consulting Engineers

## SOIL PROFILE AND TEST DATA

**Phase II - Environmental Site Assessment** 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

DATUM

BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. FILE NO. Geodetic elevation = 59.29m. REMARKS HOLE NO. BOBINGS BY CME-55 Low Clearance Drill DATE May 23 2019

**PE4614** BH 4

BORINGS BY CME-55 Low Clearance	Drill			D	DATE	May 23, 2	2019	DI14	BH 4			
SOIL DESCRIPTION			SAN	IPLE	1	DEPTH (m)	ELEV. (m)	<ul> <li>Photo Ionization Detector</li> <li>Volatile Organic Rdg. (ppm)</li> </ul>	g Well			
	STRATA PLOT	ТҮРЕ	NUMBER	° ≈ © © ©	N VALUE or RQD		(,	Lower Explosive Limit %	Monitoring Well			
GROUND SURFACE				<u>щ</u>	<b></b>	0-	-58.43	20 40 60 80				
Asphaltic concrete0.00	\$XXX						00.40		. 클 .			
FILL: Crushed stone, some sand 0.28	3	B AU	1									
		ss	2	79	13	1-	-57.43					
GLACIAL TILL: Brown silty sand with shale		1 33	2	19	13							
with shale												
		∬ss	3	100	54							
		₩ 33	5		54	2+56.43	-56.43					
2.29												
		ss	4	100	50+							
		$\Delta$										
						3-	-55.43		-   E			
		ss	5	91	50+			•	E			
		x ss	6	80	50+				ΙE			
Veathered black shale <b>BEDROCK</b>		× 33	0	00	50+	4-	-54.43		- 星			
Vealliered black shale <b>BLDHOCK</b>												
		× SS	7	100	50+				E			
		A 33		100	50+							
						5-	-53.43		- E			
									ΙE			
5.94	1											
End of Borehole		1										
GWL @ 4.06m - Sept. 3, 2020)												
								100 200 300 400	 500			
								RKI Eagle Rdg. (ppm)				
								▲ Full Gas Resp. △ Methane Elim	ı.			
								, -				

## SOIL PROFILE AND TEST DATA

**Phase II - Environmental Site Assessment** 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

#### DATUM

BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. Geodetic elevation = 59.29m. REMARKS

**PE4614** HOLE NO

FILE NO.

BORINGS BY CME-55 Low Clearance			D	DATE	BH 5						
SOIL DESCRIPTION	PLOT	SAMPLE				DEPTH			Photo Ionization Detector     Volatile Organic Rdg. (ppm)		
GROUND SURFACE	STRATA P	ТҮРЕ	NUMBER	°8 ©	N VALUE or RQD	(m)	(m)		r Explosive Li	mit %	Construction
Asphaltic concrete 0.05 FILL: Brown silty sand with 0.25 crushed stone		AU	1			0-	-58.74	•			
		ss	2	88	10	1-	-57.74	•			
<b>FILL:</b> Brown silty sand, trace gravel		ss	3	92	15	2-	-56.74	•			
3.20	ж Ж	∬ ss ∑ ss	4 5	88 100	23 50+	3-	-55.74	•		· · · · · · · · · · · · · · · · · · ·	
Weathered black shale <b>BEDROCK</b>		x ss	6	67	50+	4-	-54.74	•			
4.42	2	-								<u> </u>	
Practical refusal to augering at 4.42m depth (BH dry upon completion)											
									200 300 a Eagle Rdg. (pp as Resp. △ Metha		

## SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

DATUM
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BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. Geodetic elevation = 59.29m. REMARKS

**PE4614** HOLE NO.

FILE NO.

#### **BH 6** BORINGS BY CME-55 Low Clearance Drill DATE May 24, 2019 Monitoring Well Construction SAMPLE **Photo Ionization Detector** PLOT DEPTH ELEV. SOIL DESCRIPTION Volatile Organic Rdg. (ppm) (m) (m) RECOVERY STRATA VALUE r RQD NUMBER TYPE o/0 Lower Explosive Limit % $\bigcirc$ N OF **GROUND SURFACE** 80 20 40 60 0+58.45Asphaltic concrete 0.06 AU 1 0.30 **FILL:** Brown silty sand with crushed stone 1+57.452 SS 79 16 GLACIAL TILL: Brown clayey silt with shale, trace sand and gravel SS 3 83 38 2+56.45SS 4 91 89 <u>3</u>.05 3+55.45SS 5 67 50 +Weathered black shale BEDROCK 3.66 End of Borehole Practical refusal to augering at 3.66m depth (BH dry upon completion) 100 200 300 400 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

## SOIL PROFILE AND TEST DATA

100

200

RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

300

400

500

Monitoring Well Construction

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**Phase II - Environmental Site Assessment** 78, 84, 86 and 90 Beechwood Avenue

154 Colonnade Road South, Ottawa, Ontario K2E 7J5 83, 85, 89, 93 Barrette Street, Ottawa, Ontario BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. FILE NO. Geodetic elevation = 59.29m. **PE4614** REMARKS HOLE NO. **BH 7** BORINGS BY CME 55 Power Auger DATE July 15, 2019 SAMPLE **Photo Ionization Detector** STRATA PLOT DEPTH ELEV. SOIL DESCRIPTION • Volatile Organic Rdg. (ppm) (m) (m) N VALUE or RQD RECOVERY NUMBER TYPE o/0 Lower Explosive Limit %  $\cap$ **GROUND SURFACE** 80 20 40 60 0+58.59FILL: Brown silty sand with 0.25 AU 1 crushed stone FILL: Brown/black silty sand, some gravel 1 + 57.592 8 SS 25

1.52	ss	3	83	67	2-	56.59	
	ss	4	42	70			
Weathered black shale <b>BEDROCK</b>	ss	5	75	38	3+	55.59	
	∛ss	6	88	50+	4-	54.59	
	s ss	7	80	50+	5-	53.59	
	≤ SS	8		50+			
End of Borehole					6-	52.59	

# pa

## SOIL PROFILE AND TEST DATA

100

200

RKI Eagle Rdg. (ppm) • Full Gas Resp.  $\triangle$  Methane Elim.

300

400

500

Monitoring Well Construction

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naters	nnr			Con	sulting	g								
154 Colonnade Road Sou			-		ineers	Phase II - Environmental Site Assessment 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario								
DATUM BM - Mag na Geodetic ele REMARKS	ail on sidewalk evation = 59.29	, sou m.	th of B	Beech	wood	Ave.,	east of S	t. Charle	s Street.	FILE		PE	461	4
BORINGS BY CME 55 P	ower Auger				D	ATE .	July 15, 2	019		HOLE	: NO.	BH	8	
SOIL DESCRIPTION				SAN	IPLE		DEPTH	ELEV.		onization Detector tile Organic Rdg. (ppm)				
			ТҮРЕ	NUMBER	% RECOVERY	N VALUE of RQD	(m)	(m)		er Expl			,	Monitoring Woll
GROUND SURFACE		STRATA		ĨŇ	RE	zö	0	-58.67	20	40	60	80		
<b>FILL:</b> Brown slty sand w	vith gravel, <sub>0.30</sub>		AU	1				-30.07	•					
FILL: Brown silty sand v and construction debris	vith gravel		ss	2	29	9	1-	-57.67	•					ներերե
<u>1.52</u>		ss	3	38	49	2-	-56.67						իրիկիրինի	
		ss	4	33	34		FF 07			•			որդորդու	
GLACIAL TILL: Dense t dense, brown silty sand	o very with gravel,		ss	5	58	73	3-	-55.67	•				· · · · · · · · · · · · · · · · · · ·	կուրու
cobbles and shale			ss	6	75	80	4 -	-54.67	•					
			ss	7	83	89	5-	-53.67	•				· · · · · · · · · · · · · · · · · · ·	
	6.00		ss	8	71	61	6-	-52.67	•			·····		
Weathered black shale	BEDROCK 6.70		ss	9		56			•				· · · · · · · · · · · · · · · · · · ·	
End of Borehole														
(GWL @ 6.34m - Sept.	3, 2020)													

## SOIL PROFILE AND TEST DATA

▲ Full Gas Resp.  $\triangle$  Methane Elim.

Phase II - Environmental Site Assessment 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontari

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

83, 85, 89, 93 Barrette Street, Ottawa, Ontario											
<b>DATUM</b> BM - Mag nail on sidewalk Geodetic elevation = 59.29	, sout )m.	h of E	Beech	wood	Ave.,	east of S	t. Charle	s Street.	FILE NO.	PE4614	L
REMARKS									HOLE NO.	BH 8B	
BORINGS BY CME 55 Power Auger	DATE July 15, 2019										, 
SOIL DESCRIPTION	PLOT		SAN			DEPTH (m)	ELEV. (m)	Photo I • Vola	dg. (ppm)	Monitoring Well Construction	
	STRATA	TYPE NUMBER % RECOVERY N VALUE	VALUE r RQD			• Lowe	r Explosive	nitorir			
GROUND SURFACE	ะง		NI	REC	N OL			20	40 60	80	ΣO
FILL: Brown slty sand with gravel, 0.30		-				0-	-58.67				
<b>FILL:</b> Brown silty sand with gravel and construction debris						1-	-57.67				<u>իկկկկկի</u>
1.52		-								· · · · · · · · · · · · · · · · · · ·	<u>իրիկիիի</u>
<b>GLACIAL TILL:</b> Dense to very dense, brown silty sand with gravel, cobbles and shale						2-	-56.67		· · · · · · · · · · · · · · · · · · ·		<u>լիդվորիկու</u>
						3-	-55.67				<u>ներերերի</u>
						4-	-54.67				անալուներները երկերությունը երկերիները երկերությունը երկերությունը երկերությունը երկերությունը երկերությունը ե ԱՄՆ ընդերը երկերությունը երկերությունը երկերությունը երկերությունը երկերությունը երկերությունը երկերությունը երկ
						5-	5-53.67		· · · · · · · · · · · · · · · · · · ·	· · · · · · · · · · · · · · · · · · ·	րիկորիրին սրկություն
											<u>իսիրսիրի</u> սորերեր
6. <u>00</u>		-				6-	-52.67				
Weathered black shale <b>BEDROCK</b>						7-	-51.67				
						8-	-50.67				
8.84 End of Borehole		-									
(GWL @ 6.37m - Sept. 3, 2020)											
								100 RKI E	200 300 Eagle Rdg. (	400 50 ( <b>ppm</b> )	bo

## SOIL PROFILE AND TEST DATA

**Phase II - Environmental Site Assessment** 78, 84, 86 and 90 Beechwood Avenue

154 Colonnade Road South, Ottawa, Ontario K2E 7J5 83, 85, 89, 93 Barrette Street, Ottawa, Ontario BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. FILE NO. **PE4614** Geodetic elevation = 59.29m. REMARKS HOLE NO. **BH 9** BORINGS BY CME 55 Power Auger DATE July 15, 2019 PLOT SAMPLE **Photo Ionization Detector** DEPTH ELEV. SOIL DESCRIPTION • Volatile Organic Rdg. (ppm) (m) (m) OVERY RATA ALUE ROD **IBER** ζPE O Lower Explosive Limit %

SOIL DESCRIPTION			SAN	IPLE		DEPTH	ELEV.	Photo Ionization Detector
	STRATA PLOT	ЭДХТ	NUMBER	°% RECOVERY	N VALUE or RQD	(m)	(m)	Photo Ionization Detector       ■ unit         ● Volatile Organic Rdg. (ppm)       ■ Nition Distribution of the provided of the pr
GROUND SURFACE	S		N	RE	z °			20 40 60 80 2
<b>FILL:</b> Brown silty sand with 0.33		XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX	1			0-	-58.71	
FILL: Brown silty sand		ss	2	8	9	1-	-57.71	
2.30		ss	3	42	18	2-	-56.71	
		ss	4	53	50+	2.	-55.71	
		x ss	5	80	50+	5	-55.71	
<b>GLACIAL TILL:</b> Dense to very dense, grey silty sand with gravel and shale fragments		ss	6	75	38	4-	-54.71	
and shale fragments		ss	7	71	51	5-	-53.71	
		ss	8	71	52	6-	-52.71	
6.65 Weathered black shale <b>BEDROCK</b> 6.86 End of Borehole		X SS	8	64	66			
								100 200 300 400 500 <b>RKI Eagle Rdg. (ppm)</b> ▲ Full Gas Resp. △ Methane Elim.

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## SOIL PROFILE AND TEST DATA

**PE4614** 

**BH10** 

**Phase II - Environmental Site Assessment** 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. FILE NO. Geodetic elevation = 59.29m. REMARKS HOLE NO. BORINGS BY CME 55 Power Auger **DATE** July 15, 2019

SOIL DESCRIPTION	E O						ELEV.	Photo Ionization Detector ● Volatile Organic Rdg. (ppm)
GROUND SURFACE	STRATA P	TYPE	NUMBER	°% RECOVERY	N VALUE or RQD	(m)	(m)	Photo Ionization Detector       III of unitarian providence of the sector <ul> <li>Volatile Organic Rdg. (ppm)</li> <li>C Lower Explosive Limit %</li> <li>20</li> <li>40</li> <li>60</li> <li>80</li> </ul> <li>III of unitarian providence of the sector</li> <li>IIII of unitarian providence of the sector</li>
FILL: Brown silty sand with         0.2           Crushed stone	5	au Bartes	1			0-	-58.66	
FILL: Brown silty sand, some gravel, trace organics		ss	2	62	9	1-	-57.66	
GLACIAL TILL: Dense to very		ss	3	100	34	2-	-56.66	
dense, brown silty sand, some gravel and shale		ss	4	100	50+			
3.0	5\^^^^	∑ss	5	70	50+	3-	-55.66	
		X SS	6	100	50+	4-	-54.66	
Weathered black shale <b>BEDROCK</b>		∑ss	7	100	50+	5-	-53.66	•
		x ss	8	100	50+			
6.7		X SS	9	50	50+	6-	-52.66	
6.7 End of Borehole (GWL @ 4.72m - Sept. 3, 2020)								
								100 200 300 400 500 <b>RKI Eagle Rdg. (ppm)</b> ▲ Full Gas Resp. △ Methane Elim.

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## SOIL PROFILE AND TEST DATA

100

200

300

**RKI Eagle Rdg. (ppm)** ▲ Full Gas Resp. △ Methane Elim.

400

500

Monitoring Well Construction

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Dalerson	][(			Con	suitin	g	001			- / \\						
154 Colonnade Road South, Ottaw			-		ineers	78	ase II - E , 84, 86 a , 85, 89, 9	nd 90 Be	echw	vood	A٧	/enue				
DATUM BM - Mag nail on side Geodetic elevation =	ewalk, 59.29	sout m.	h of B	eech	wood	Ave.,	east of S	t. Charle	s Stre	eet.	F	FILE NO.		PE <sup>,</sup>	461 <sup>,</sup>	4
REMARKS											ŀ	HOLE NC	).	рμ	14 4	
BORINGS BY CME-55 Low Cleara	ance D	Drill			D	ATE /	August 24	4, 2020	1					BH		_
SOIL DESCRIPTION		PLOT		SAN			DEPTH (m)	ELEV. (m)				<b>ization</b> Organic				
GROUND SURFACE		STRATA	ТҮРЕ	NUMBER	% RECOVERY	N VALUE or RQD			0	Lowe		Explosi		_imit 80	: %	
Asphaltic concrete FILL: Brown silty sand with	0.10 0.60		AU	1			- 0-	-58.11	•							1
			ss	2	62	6	1-	-57.11	•							
<b>GLACIAL TILL:</b> Dense to very dense, brown silty sand with shale	e		ss	3	54	40	2-	-56.11	•							
			ss	4	79	44	3-	-55.11								
	<u>3.80</u>		ss	5	76	50+									) - ( ) - (- ) - (	
			⊻ ss ∝ ss	6 7	78 100	50+ 50+	4-	-54.11								
Possible weathered <b>BEDROCK</b>			∆ 00	7	100	50+	5-	-53.11								
End of Borehole	<u>5.84</u>		-							· · · · · · · · · · · · · · · · · · ·				<u></u>	<u>.</u>	
(GWL @ 4.28m - Sept. 3, 2020)																

## SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

#### DATUM

BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. Geodetic elevation = 59.29m.

**PE4614** 

FILE NO.

#### REMARKS HOLE NO. **BH12** BORINGS BY CME-55 Low Clearance Drill DATE August 24, 2020 Monitoring Well Construction SAMPLE **Photo Ionization Detector** STRATA PLOT DEPTH ELEV. SOIL DESCRIPTION Volatile Organic Rdg. (ppm) (m) (m) RECOVERY N VALUE or RQD NUMBER TYPE o/0 Lower Explosive Limit % $\bigcirc$ **GROUND SURFACE** 80 20 40 60 0+57.57Asphaltic concrete 0.10 FILL: Brown silty sand with AU 1 crushed stone 0.63 1+56.57SS 2 46 10 GLACIAL TILL: Compact to very dense, brown silty sand with shale SS 3 58 52 2 + 55.572.30 SS 4 71 50 +Weathered black shale **BEDROCK** 3+54.57 50+ 3.28 SS 5 11 End of Borehole

RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

300

400

500

200

100

## SOIL PROFILE AND TEST DATA

Monitoring Well Construction

**Phase II - Environmental Site Assessment** 78 84 86 and 90 Reechwood Avenue

154 Colonnade Road South, Ottawa, Ont	ario ł	<2E 7J	5			, ,		te Street, (	Ottawa, On	tario	
DATUM BM - Mag nail on sidewalk Geodetic elevation = 59.29		th of B	eech	wood	Ave.,	east of S	t. Charle	s Street.	FILE NO.	PE4614	ł
REMARKS BORINGS BY CME-55 Low Clearance I	Drill			D	ATE /	August 24	4, 2020		HOLE NO.	BH13	_
SOIL DESCRIPTION			SAM	IPLE		DEPTH	ELEV.		onization D tile Organic Re		
GROUND SURFACE	STRATA PL	ТҮРЕ	NUMBER	% RECOVERY	N VALUE or RQD		(m)		r Explosive		1
FILL: Brown silty sand with crushed stone0.43		AU	1			- 0-	-58.73				T FULLING LINE L
		ss	2	83	11	1-	-57.73				
Compact, brown <b>SILTY SAND,</b> some gravel		ss	3	67	24	2-	-56.73				

75

62

90

17

20

50+

SS

SS

SS

3.50

GLACIAL TILL: Very dense, brown

Weathered black shale **BEDROCK** 

(GWL @ 3.20m - Sept. 3, 2020)

silty sand with shale

End of Borehole

4

5

6

4.60 SS 7 50+ 90 5+53.73 5.89

100

200

**RKI Eagle Rdg. (ppm)** ▲ Full Gas Resp. △ Methane Elim.

300

400

500

3+55.73

4+54.73

# patersongroup Consulting Engineers

## SOIL PROFILE AND TEST DATA

**Phase II - Environmental Site Assessment** 78, 84, 86 and 90 Beechwood Avenue 83, 85, 89, 93 Barrette Street, Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

#### DATUM

BM - Mag nail on sidewalk, south of Beechwood Ave., east of St. Charles Street. Geodetic elevation = 59.29m. REMARKS

**PE4614** 

FILE NO.

BORINGS BY CME-55 Low Clearance	Drill			D	ATE	August 24	4, 2020		HOLE	B	H14	
SOIL DESCRIPTION	PLOT		SAN	IPLE	1	DEPTH	ELEV.			<b>ion Detec</b> anic Rdg. (p		y Well
	STRATA	ТҮРЕ	NUMBER	°⊗ RECOVERY	N VALUE or RQD	(m)	(m)			osive Lin		Monitoring Well Construction
GROUND SURFACE	ß		N	RE	z °	0	-58.25	20	40	60 B	30	ž
FILL: Brown silty sand with crushed stone		AU	1									
<u>1.07</u>		ss	2	67	38	1-	-57.25	•				
<b>GLACIAL TILL:</b> Very dense, brown/black silty sand with shale		ss	3	71	60	2-	-56.25	•				
		ss	4	73	50+							
3.00		A 33	4	13	50+	3-	-55.25					
Weathered black shale <b>BEDROCK</b> <u>3.30</u> End of Borehole		ss	5	50	50+	5	55.25	•····				
										300 44 Rdg. (ppn ). △ Methar		00

## SYMBOLS AND TERMS

#### SOIL DESCRIPTION

Behavioural properties, such as structure and strength, take precedence over particle gradation in describing soils. Terminology describing soil structure are as follows:

Desiccated	-	having visible signs of weathering by oxidation of clay minerals, shrinkage cracks, etc.
Fissured	-	having cracks, and hence a blocky structure.
Varved	-	composed of regular alternating layers of silt and clay.
Stratified	-	composed of alternating layers of different soil types, e.g. silt and sand or silt and clay.
Well-Graded	-	Having wide range in grain sizes and substantial amounts of all intermediate particle sizes (see Grain Size Distribution).
Uniformly-Graded	-	Predominantly of one grain size (see Grain Size Distribution).

The standard terminology to describe the relative strength of cohesionless soils is the compactness condition, usually inferred from the results of the Standard Penetration Test (SPT) 'N' value. The SPT N value is the number of blows of a 63.5 kg hammer, falling 760 mm, required to drive a 51 mm O.D. split spoon sampler 300 mm into the soil after an initial penetration of 150 mm. An SPT N value of "P" denotes that the split-spoon sampler was pushed 300 mm into the soil without the use of a falling hammer.

Compactness Condition	'N' Value	Relative Density %
Very Loose	<4	<15
Loose	4-10	15-35
Compact	10-30	35-65
Dense	30-50	65-85
Very Dense	>50	>85

The standard terminology to describe the strength of cohesive soils is the consistency, which is based on the undisturbed undrained shear strength as measured by the in situ or laboratory shear vane tests, unconfined compression tests, or occasionally by the Standard Penetration Test (SPT). Note that the typical correlations of undrained shear strength to SPT N value (tabulated below) tend to underestimate the consistency for sensitive silty clays, so Paterson reviews the applicable split spoon samples in the laboratory to provide a more representative consistency value based on tactile examination.

Consistency	Undrained Shear Strength (kPa)	'N' Value
Very Soft	<12	<2
Soft	12-25	2-4
Firm	25-50	4-8
Stiff	50-100	8-15
Very Stiff	100-200	15-30
Hard	>200	>30

### SYMBOLS AND TERMS (continued)

#### **SOIL DESCRIPTION (continued)**

Cohesive soils can also be classified according to their "sensitivity". The sensitivity, St, is the ratio between the undisturbed undrained shear strength and the remoulded undrained shear strength of the soil. The classes of sensitivity may be defined as follows:

Low Sensitivity:	St < 2
Medium Sensitivity:	$2 < S_t < 4$
Sensitive:	$4 < S_t < 8$
Extra Sensitive:	8 < St < 16
Quick Clay:	St > 16

#### **ROCK DESCRIPTION**

The structural description of the bedrock mass is based on the Rock Quality Designation (RQD).

The RQD classification is based on a modified core recovery percentage in which all pieces of sound core over 100 mm long are counted as recovery. The smaller pieces are considered to be a result of closely-spaced discontinuities (resulting from shearing, jointing, faulting, or weathering) in the rock mass and are not counted. RQD is ideally determined from NQ or larger size core. However, it can be used on smaller core sizes, such as BQ, if the bulk of the fractures caused by drilling stresses (called "mechanical breaks") are easily distinguishable from the normal in situ fractures.

#### RQD % ROCK QUALITY

90-100	Excellent, intact, very sound
75-90	Good, massive, moderately jointed or sound
50-75	Fair, blocky and seamy, fractured
25-50	Poor, shattered and very seamy or blocky, severely fractured
0-25	Very poor, crushed, very severely fractured

#### SAMPLE TYPES

SS	-	Split spoon sample (obtained in conjunction with the performing of the Standard Penetration Test (SPT))					
TW	-	Thin wall tube or Shelby tube, generally recovered using a piston sampler					
G	-	"Grab" sample from test pit or surface materials					
AU	-	Auger sample or bulk sample					
WS	-	Wash sample					
RC	-	Rock core sample (Core bit size BQ, NQ, HQ, etc.). Rock core samples are obtained with the use of standard diamond drilling bits.					

### SYMBOLS AND TERMS (continued)

### PLASTICITY LIMITS AND GRAIN SIZE DISTRIBUTION

WC%	-	Natural water content or water content of sample, %					
LL	-	Liquid Limit, % (water content above which soil behaves as a liquid)					
PL	-	Plastic Limit, % (water content above which soil behaves plastically)					
PI	-	Plasticity Index, % (difference between LL and PL)					
Dxx	-	Grain size at which xx% of the soil, by weight, is of finer grain sizes These grain size descriptions are not used below 0.075 mm grain size					
D10	-	Grain size at which 10% of the soil is finer (effective grain size)					
D60	-	Grain size at which 60% of the soil is finer					
Сс	-	Concavity coefficient = $(D30)^2 / (D10 \times D60)$					
Cu	-	Uniformity coefficient = D60 / D10					
	0	we also access the supplicer of several and supplices					

Cc and Cu are used to assess the grading of sands and gravels: Well-graded gravels have: 1 < Cc < 3 and Cu > 4Well-graded sands have: 1 < Cc < 3 and Cu > 6Sands and gravels not meeting the above requirements are poorly-graded or uniformly-graded. Cc and Cu are not applicable for the description of soils with more than 10% silt and clay (more than 10% finer than 0.075 mm or the #200 sieve)

### **CONSOLIDATION TEST**

p'o	-	Present effective overburden pressure at sample depth
p'c	-	Preconsolidation pressure of (maximum past pressure on) sample
Ccr	-	Recompression index (in effect at pressures below p'c)
Cc	-	Compression index (in effect at pressures above p'c)
OC Ratio	)	Overconsolidaton ratio = p'c / p'o
Void Rati	io	Initial sample void ratio = volume of voids / volume of solids
Wo	-	Initial water content (at start of consolidation test)

### PERMEABILITY TEST

k - Coefficient of permeability or hydraulic conductivity is a measure of the ability of water to flow through the sample. The value of k is measured at a specified unit weight for (remoulded) cohesionless soil samples, because its value will vary with the unit weight or density of the sample during the test.

### SYMBOLS AND TERMS (continued) STRATA PLOT Topsoil Asphalt Peat Sand Silty Sand Fill $\nabla$ Sandy Silt Clay Silty Clay Clayey Silty Sand Glacial Till Shale Bedrock

### MONITORING WELL AND PIEZOMETER CONSTRUCTION



PIEZOMETER CONSTRUCTION





RELIABLE.

## Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Mark D'Arcy

Client PO: 26494 Project: PE4614 Custody: 122147

Report Date: 30-May-2019 Order Date: 24-May-2019

Order #: 1921505

This Certificate of Analysis contains analytical data applicable to the following samples as submitted:

Paracel ID	Client ID
1921505-01	BH1-SS2
1921505-02	BH2-SS3
1921505-03	BH3-AU1
1921505-04	BH3-SS9/10

Approved By:

Mark Foto

Mark Foto, M.Sc. Lab Supervisor

Any use of these results implies your agreement that our total liability in connection with this work, however arising, shall be limited to the amount paid by you for this work, and that our employees or agents shall not under any circumstances be liable to you in connection with this work.



### **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	29-May-19	30-May-19
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	24-May-19	28-May-19
Mercury by CVAA	EPA 7471B - CVAA, digestion	30-May-19	30-May-19
pH, soil	EPA 150.1 - pH probe @ 25 °C, CaCl buffered ext.	28-May-19	28-May-19
PHC F1	CWS Tier 1 - P&T GC-FID	29-May-19	30-May-19
PHC F4G (gravimetric)	CWS Tier 1 - Extraction Gravimetric	29-May-19	30-May-19
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	24-May-19	27-May-19
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	29-May-19	29-May-19
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	24-May-19	27-May-19
REG 153: VOCs by P&T GC/MS	EPA 8260 - P&T GC-MS	29-May-19	30-May-19
Solids, %	Gravimetric, calculation	29-May-19	29-May-19

Report Date: 30-May-2019 Order Date: 24-May-2019



Order #: 1921505

Report Date: 30-May-2019 Order Date: 24-May-2019

	Client ID: Sample Date:	BH1-SS2 23-May-19 09:00	BH2-SS3 23-May-19 10:30	BH3-AU1 23-May-19 12:00	BH3-SS9/10 23-May-19 09:00
	Sample ID:	1921505-01	1921505-02	1921505-03	1921505-04
	MDL/Units	Soil	Soil	Soil	Soil
Physical Characteristics					
% Solids	0.1 % by Wt.	96.7	90.5	98.6	89.8
General Inorganics					
рН	0.05 pH Units	7.94	7.39	-	-
Metals			I		
Antimony	1.0 ug/g dry	<1.0	-	<1.0	-
Arsenic	1.0 ug/g dry	7.7	-	4.0	-
Barium	1.0 ug/g dry	256	-	54.9	-
Beryllium	0.5 ug/g dry	0.6	-	<0.5	-
Boron	5.0 ug/g dry	15.8	-	6.8	-
Cadmium	0.5 ug/g dry	<0.5	-	<0.5	-
Chromium	5.0 ug/g dry	30.1	-	10.1	-
Chromium (VI)	0.2 ug/g dry	<0.2	-	<0.2	-
Cobalt	1.0 ug/g dry	11.7	-	4.7	-
Copper	5.0 ug/g dry	37.5	-	13.5	-
Lead	1.0 ug/g dry	181	-	44.7	-
Mercury	0.1 ug/g dry	0.3	-	<0.1	-
Molybdenum	1.0 ug/g dry	3.3	-	2.2	-
Nickel	5.0 ug/g dry	31.6	-	16.0	-
Selenium	1.0 ug/g dry	<1.0	-	<1.0	-
Silver	0.3 ug/g dry	0.5	-	<0.3	-
Thallium	1.0 ug/g dry	<1.0	-	<1.0	-
Uranium	1.0 ug/g dry	1.4	-	<1.0	-
Vanadium	10.0 ug/g dry	27.5	-	24.7	-
Zinc	20.0 ug/g dry	143	-	41.6	-
Volatiles					
Acetone	0.50 ug/g dry	-	-	-	<0.50
Benzene	0.02 ug/g dry	-	-	-	<0.02
Bromodichloromethane	0.05 ug/g dry	-	-	-	<0.05
Bromoform	0.05 ug/g dry	-	-	-	<0.05
Bromomethane	0.05 ug/g dry	-	-	-	<0.05
Carbon Tetrachloride	0.05 ug/g dry	-	-	-	<0.05
Chlorobenzene	0.05 ug/g dry	-	-	-	<0.05
Chloroform	0.05 ug/g dry	-	-	-	<0.05
Dibromochloromethane	0.05 ug/g dry	-	-	-	<0.05
Dichlorodifluoromethane	0.05 ug/g dry	-	-	-	<0.05



Report Date: 30-May-2019 Order Date: 24-May-2019

-	Client ID: Sample Date: Sample ID:	BH1-SS2 23-May-19 09:00 1921505-01	BH2-SS3 23-May-19 10:30 1921505-02	BH3-AU1 23-May-19 12:00 1921505-03	BH3-SS9/10 23-May-19 09:00 1921505-04
	MDL/Units	Soil	Soil	Soil	Soil
1,2-Dichlorobenzene	0.05 ug/g dry	-	-	-	<0.05
1,3-Dichlorobenzene	0.05 ug/g dry	-	-	-	<0.05
1,4-Dichlorobenzene	0.05 ug/g dry	-	-	-	<0.05
1,1-Dichloroethane	0.05 ug/g dry	-	-	-	<0.05
1,2-Dichloroethane	0.05 ug/g dry	-	-	-	<0.05
1,1-Dichloroethylene	0.05 ug/g dry	-	-	-	<0.05
cis-1,2-Dichloroethylene	0.05 ug/g dry	-	-	-	<0.05
trans-1,2-Dichloroethylene	0.05 ug/g dry	-	-	-	<0.05
1,2-Dichloropropane	0.05 ug/g dry	-	-	-	<0.05
cis-1,3-Dichloropropylene	0.05 ug/g dry	-	-	-	<0.05
trans-1,3-Dichloropropylene	0.05 ug/g dry	-	-	-	<0.05
1,3-Dichloropropene, total	0.05 ug/g dry	-	-	-	<0.05
Ethylbenzene	0.05 ug/g dry	-	-	-	<0.05
Ethylene dibromide (dibromoethan	0.05 ug/g dry	-	-	-	<0.05
Hexane	0.05 ug/g dry	-	-	-	<0.05
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	-	-	-	<0.50
Methyl Isobutyl Ketone	0.50 ug/g dry	-	-	-	<0.50
Methyl tert-butyl ether	0.05 ug/g dry	-	-	-	<0.05
Methylene Chloride	0.05 ug/g dry	-	-	-	<0.05
Styrene	0.05 ug/g dry	-	-	-	<0.05
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	-	-	-	<0.05
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	-	-	-	<0.05
Tetrachloroethylene	0.05 ug/g dry	-	-	-	1.79
Toluene	0.05 ug/g dry	-	-	-	<0.05
1,1,1-Trichloroethane	0.05 ug/g dry	-	-	-	<0.05
1,1,2-Trichloroethane	0.05 ug/g dry	-	-	-	<0.05
Trichloroethylene	0.05 ug/g dry	-	-	-	<0.05
Trichlorofluoromethane	0.05 ug/g dry	-	-	-	<0.05
Vinyl chloride	0.02 ug/g dry	-	-	-	<0.02
m,p-Xylenes	0.05 ug/g dry	-	-	-	0.21
o-Xylene	0.05 ug/g dry	-	-	-	0.09
Xylenes, total	0.05 ug/g dry	-	-	-	0.29
4-Bromofluorobenzene	Surrogate	-	-	-	89.7%
Dibromofluoromethane	Surrogate	-	-	-	81.5%
Toluene-d8	Surrogate	-	-	-	79.5%
Benzene	0.02 ug/g dry	<0.02	<0.02	<0.02	-



Order #: 1921505

Report Date: 30-May-2019 Order Date: 24-May-2019

	Client ID: Sample Date: Sample ID: MDL/Units	BH1-SS2 23-May-19 09:00 1921505-01 Soil	BH2-SS3 23-May-19 10:30 1921505-02 Soil	BH3-AU1 23-May-19 12:00 1921505-03 Soil	BH3-SS9/10 23-May-19 09:00 1921505-04 Soil
Ethylbenzene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Toluene	0.05 ug/g dry	0.05	<0.05	<0.05	-
m,p-Xylenes	0.05 ug/g dry	0.67	<0.05	<0.05	-
o-Xylene	0.05 ug/g dry	0.06	<0.05	<0.05	-
Xylenes, total	0.05 ug/g dry	0.73	<0.05	<0.05	-
Toluene-d8	Surrogate	75.3%	76.7%	88.2%	-
Hydrocarbons					
F1 PHCs (C6-C10)	7 ug/g dry	44	<7	15	-
F2 PHCs (C10-C16)	4 ug/g dry	22	80	1040	-
F3 PHCs (C16-C34)	8 ug/g dry	42	29	2370	-
F4 PHCs (C34-C50)	6 ug/g dry	12	<6	2590 [1]	-
F4G PHCs (gravimetric)	50 ug/g dry	-	-	9490	-
Semi-Volatiles			-	-	
Acenaphthene	0.02 ug/g dry	0.06	-	-	-
Acenaphthylene	0.02 ug/g dry	0.05	-	-	-
Anthracene	0.02 ug/g dry	0.13	-	-	-
Benzo [a] anthracene	0.02 ug/g dry	0.37	-	-	-
Benzo [a] pyrene	0.02 ug/g dry	0.29	-	-	-
Benzo [b] fluoranthene	0.02 ug/g dry	0.45	-	-	-
Benzo [g,h,i] perylene	0.02 ug/g dry	0.20	-	-	-
Benzo [k] fluoranthene	0.02 ug/g dry	0.31	-	-	-
Chrysene	0.02 ug/g dry	0.38	-	-	-
Dibenzo [a,h] anthracene	0.02 ug/g dry	0.07	-	-	-
Fluoranthene	0.02 ug/g dry	0.78	-	-	-
Fluorene	0.02 ug/g dry	0.06	-	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	0.16	-	-	-
1-Methylnaphthalene	0.02 ug/g dry	<0.02	-	-	-
2-Methylnaphthalene	0.02 ug/g dry	0.03	-	-	-
Methylnaphthalene (1&2)	0.04 ug/g dry	0.04	-	-	-
Naphthalene	0.01 ug/g dry	0.04	-	-	-
Phenanthrene	0.02 ug/g dry	0.58	-	-	-
Pyrene	0.02 ug/g dry	0.62	-	-	-
2-Fluorobiphenyl	Surrogate	84.3%	-	-	-
Terphenyl-d14	Surrogate	117%	-	-	-



Order #: 1921505

Report Date: 30-May-2019

Order Date: 24-May-2019

Project Description: PE4614

## Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g						
F2 PHCs (C10-C16)	ND	4	ug/g						
F3 PHCs (C16-C34)	ND	8	ug/g						
F4 PHCs (C34-C50)	ND	6	ug/g						
F4G PHCs (gravimetric)	ND	50	ug/g						
Metals									
Antimony	ND	1.0	ug/g						
Arsenic	ND	1.0	ug/g						
Barium	ND	1.0	ug/g						
Beryllium	ND	0.5	ug/g						
Boron	ND	5.0	ug/g						
Cadmium	ND	0.5	ug/g						
Chromium (VI)	ND	0.2	ug/g						
Chromium	ND	5.0	ug/g						
Cobalt	ND	1.0	ug/g						
Copper	ND	5.0	ug/g						
Lead	ND	1.0	ug/g						
Mercury	ND	0.1	ug/g						
Molybdenum	ND	1.0	ug/g						
Nickel	ND	5.0	ug/g						
Selenium	ND	1.0	ug/g						
Silver	ND	0.3	ug/g						
Thallium	ND	1.0	ug/g						
Uranium	ND	1.0	ug/g						
Vanadium	ND ND	10.0	ug/g						
Zinc Somi Volotiloo	ND	20.0	ug/g						
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g						
Acenaphthylene	ND	0.02	ug/g						
Anthracene	ND	0.02	ug/g						
Benzo [a] anthracene	ND	0.02	ug/g						
Benzo [a] pyrene	ND ND	0.02 0.02	ug/g						
Benzo [b] fluoranthene Benzo [g,h,i] perylene	ND	0.02	ug/g						
Benzo [k] fluoranthene	ND	0.02	ug/g ug/g						
Chrysene	ND	0.02	ug/g ug/g						
Dibenzo [a,h] anthracene	ND	0.02	ug/g ug/g						
Fluoranthene	ND	0.02	ug/g ug/g						
Fluorene	ND	0.02	ug/g						
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g						
1-Methylnaphthalene	ND	0.02	ug/g						
2-Methylnaphthalene	ND	0.02	ug/g						
Methylnaphthalene (1&2)	ND	0.04	ug/g						
Naphthalene	ND	0.01	ug/g						
Phenanthrene	ND	0.02	ug/g						
Pyrene	ND	0.02	ug/g						
Surrogate: 2-Fluorobiphenyl	1.38		ug/g		104	50-140			
Surrogate: Terphenyl-d14	1.66		ug/g		125	50-140			
Volatiles									
Acetone	ND	0.50	ug/g						
Benzene	ND	0.02	ug/g						
Bromodichloromethane	ND	0.05	ug/g						
Bromoform	ND	0.05	ug/g						
Bromomethane	ND	0.05	ug/g						
Carbon Tetrachloride	ND	0.05	ug/g						
Chlorobenzene	ND	0.05	ug/g						
Chloroform	ND	0.05	ug/g						
Dibromochloromethane	ND	0.05	ug/g						



#### Order #: 1921505

Report Date: 30-May-2019

Order Date: 24-May-2019

Project Description: PE4614

### Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Dichlorodifluoromethane	ND	0.05	ug/g						
1.2-Dichlorobenzene	ND	0.05	ug/g						
1,3-Dichlorobenzene	ND	0.05	ug/g						
1,4-Dichlorobenzene	ND	0.05	ug/g						
1,1-Dichloroethane	ND	0.05	ug/g						
1,2-Dichloroethane	ND	0.05	ug/g						
1,1-Dichloroethylene	ND	0.05	ug/g						
cis-1,2-Dichloroethylene	ND	0.05	ug/g						
trans-1,2-Dichloroethylene	ND	0.05	ug/g						
1,2-Dichloropropane	ND	0.05	ug/g						
cis-1,3-Dichloropropylene	ND	0.05	ug/g						
trans-1,3-Dichloropropylene	ND	0.05	ug/g						
1,3-Dichloropropene, total	ND	0.05	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Ethylene dibromide (dibromoethane	ND	0.05	ug/g						
Hexane	ND	0.05	ug/g						
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g						
Methyl Isobutyl Ketone	ND	0.50	ug/g						
Methyl tert-butyl ether	ND	0.05	ug/g						
Methylene Chloride	ND	0.05	ug/g						
Styrene	ND	0.05	ug/g						
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g						
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g						
Tetrachloroethylene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
1,1,1-Trichloroethane	ND	0.05	ug/g						
1,1,2-Trichloroethane	ND	0.05	ug/g						
Trichloroethylene	ND	0.05	ug/g						
Trichlorofluoromethane	ND	0.05	ug/g						
Vinyl chloride	ND	0.02	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g		100	50 1 10			
Surrogate: 4-Bromofluorobenzene	3.47		ug/g		108	50-140			
Surrogate: Dibromofluoromethane	3.76		ug/g		117	50-140			
Surrogate: Toluene-d8	3.03		ug/g		94.6	50-140			
Benzene	ND	0.02	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g						
Surrogate: Toluene-d8	3.03		ug/g		94.6	50-140			



Order #: 1921505

Report Date: 30-May-2019

Order Date: 24-May-2019

Project Description: PE4614

### Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
General Inorganics		0.07		0.70					
pH	6.82	0.05	pH Units	6.79			0.4	10	
Hydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g dry	ND				40	
F2 PHCs (C10-C16)	ND	4	ug/g dry	ND				30	
F3 PHCs (C16-C34)	ND	8	ug/g dry	ND				30	
F4 PHCs (C34-C50)	ND	6	ug/g dry	ND				30	
Metals									
Antimony	ND	1.0	ug/g dry	ND			0.0	30	
Arsenic	7.1	1.0	ug/g dry	6.4			9.0	30	
Barium Beryllium	60.9 0.6	1.0 0.5	ug/g dry	54.7 0.5			10.8 10.3	30 30	
Boron	9.7	5.0	ug/g dry ug/g dry	0.5 8.5			13.2	30	
Cadmium	ND	0.5	ug/g dry	ND			0.0	30	
Chromium (VI)	ND	0.2	ug/g dry	ND			0.0	35	
Chromium	18.0	5.0	ug/g dry	16.5			8.4	30	
Cobalt	7.0	1.0	ug/g dry	6.0			14.5	30	
Copper	45.5	5.0	ug/g dry	40.3			12.0	30	
Lead	26.1	1.0	ug/g dry	29.0			10.5	30	
Mercury	0.245	0.1	ug/g dry	0.259			5.6	30	
Molybdenum Nickel	ND 15.9	1.0 5.0	ug/g dry	ND 14.8			0.0 6.9	30 30	
Selenium	ND	1.0	ug/g dry ug/g dry	ND			0.9	30	
Silver	ND	0.3	ug/g dry	ND			0.0	30	
Thallium	ND	1.0	ug/g dry	ND			0.0	30	
Uranium	ND	1.0	ug/g dry	ND			0.0	30	
Vanadium	26.4	10.0	ug/g dry	22.9			14.4	30	
Zinc	124	20.0	ug/g dry	125			0.9	30	
Physical Characteristics									
% Solids	90.5	0.1	% by Wt.	90.2			0.3	25	
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g dry	ND				40	
Acenaphthylene	ND	0.02	ug/g dry	ND			0.0	40	
Anthracene	ND	0.02	ug/g dry	ND			0.0	40	
Benzo [a] anthracene	0.056	0.02	ug/g dry	0.059			5.4	40	
Benzo [a] pyrene	0.079	0.02	ug/g dry	0.098			22.4	40	
Benzo [b] fluoranthene Benzo [g,h,i] perylene	0.130 0.067	0.02 0.02	ug/g dry	0.132 0.089			1.6 27.9	40 40	
Benzo [k] fluoranthene	0.051	0.02	ug/g dry ug/g dry	0.009			30.1	40	
Chrysene	0.087	0.02	ug/g dry	0.086			0.6	40	
Dibenzo [a,h] anthracene	0.022	0.02	ug/g dry	ND			0.0	40	
Fluoranthene	0.118	0.02	ug/g dry	0.116			1.8	40	
Fluorene	ND	0.02	ug/g dry	ND				40	
Indeno [1,2,3-cd] pyrene	0.059	0.02	ug/g dry	0.077			26.2	40	
1-Methylnaphthalene	ND	0.02	ug/g dry	ND			0.0	40	
2-Methylnaphthalene Naphthalene	ND ND	0.02 0.01	ug/g dry	ND ND			0.0 0.0	40 40	
Phenanthrene	0.041	0.01	ug/g dry	0.042			0.0 3.4	40 40	
Pyrene	0.103	0.02	ug/g dry ug/g dry	0.042			3.4	40	
Surrogate: 2-Fluorobiphenyl	1.41	0.02	ug/g dry	0.100	80.6	50-140	0.2		
Surrogate: Terphenyl-d14	1.57		ug/g dry		90.1	50-140			
Volatiles									
Acetone	ND	0.50	ug/g dry	ND				50	
Benzene	ND	0.02	ug/g dry ug/g dry	ND				50	
Bromodichloromethane	ND	0.05	ug/g dry	ND				50	
Bromoform	ND	0.05	ug/g dry	ND				50	
Bromomethane	ND	0.05	ug/g dry	ND				50	



Analyte

Carbon Tetrachloride

Surrogate: 4-Bromofluorobenzene

Surrogate: Dibromofluoromethane

Surrogate: Toluene-d8

Surrogate: Toluene-d8

Benzene

Toluene

o-Xylene

Ethylbenzene

m,p-Xylenes

#### Certificate of Analysis **Client: Paterson Group Consulting Engineers** Client PO: 26494

### Method Quality Control: Duplicate

RPD

Limit

50

50

50

50

50

50

RPD

%REC

Limit

Report Date: 30-May-2019 Order Date: 24-May-2019

Project Description: PE4614

Notes

Chlorobenzene ND 0.05 ug/g dry ND	50 50
	50
Chloroform ND 0.05 ug/g dry ND	50
Dibromochloromethane ND 0.05 ug/g dry ND	50
Dichlorodifluoromethane ND 0.05 ug/g dry ND	50
1,2-Dichlorobenzene ND 0.05 ug/g dry ND	50
1,3-Dichlorobenzene ND 0.05 ug/g dry ND	50
1,4-Dichlorobenzene ND 0.05 ug/g dry ND	50
1,1-Dichloroethane ND 0.05 ug/g dry ND	50
1,2-Dichloroethane ND 0.05 ug/g dry ND	50
1,1-Dichloroethylene ND 0.05 ug/g dry ND	50
cis-1,2-Dichloroethylene ND 0.05 ug/g dry ND	50
trans-1,2-Dichloroethylene ND 0.05 ug/g dry ND	50
1,2-Dichloropropane ND 0.05 ug/g dry ND	50
cis-1,3-Dichloropropylene ND 0.05 ug/g dry ND	50
trans-1,3-Dichloropropylene ND 0.05 ug/g dry ND	50
Ethylbenzene ND 0.05 ug/g dry ND	50
Ethylene dibromide (dibromoethane ND 0.05 ug/g dry ND	50
Hexane ND 0.05 ug/g dry ND	50
Methyl Ethyl Ketone (2-Butanone) ND 0.50 ug/g dry ND	50
Methyl Isobutyl Ketone ND 0.50 ug/g dry ND	50
Methyl tert-butyl ether ND 0.05 ug/g dry ND	50
Methylene Chloride ND 0.05 ug/g dry ND	50
StyreneND 0.05 ug/g dry ND	50
1,1,1,2-Tetrachloroethane ND 0.05 ug/g dry ND	50
1,1,2,2-Tetrachloroethane ND 0.05 ug/g dry ND	50
Tetrachloroethylene ND 0.05 ug/g dry ND	50
Toluene ND 0.05 ug/g dry ND	50
1,1,1-Trichloroethane ND 0.05 ug/g dry ND	50
1,1,2-Trichloroethane ND 0.05 ug/g dry ND	50
Trichloroethylene ND 0.05 ug/g dry ND	50
Trichlorofluoromethane ND 0.05 ug/g dry ND	50
Vinyl chloride ND 0.02 ug/g dry ND	50
m,p-Xylenes ND 0.05 ug/g dry ND	50
o-Xylene ND 0.05 ug/g dry ND	50

Source

Result

ND

%REC

98.6

113

98.6

98.6

ND

ND

ND

ND

ND

50-140

50-140

50-140

50-140

Units

ug/g dry

Reporting

Limit

0.05

Result

ND

3.40

3.88

3.40

ND

ND

ND

ND

ND

3.40

0.02

0.05

0.05

0.05

0.05



## Method Quality Control: Spike

Report Date: 30-May-2019

Order Date: 24-May-2019

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	201	7	ug/g		100	80-120			
F2 PHCs (C10-C16)	95	4	ug/g	ND	78.8	60-140			
F3 PHCs (C16-C34)	259	8	ug/g	ND	87.3	60-140			
F4 PHCs (C34-C50)	153	6	ug/g	ND	81.4	60-140			
F4G PHCs (gravimetric)	810	50	ug/g		81.0	80-120			
Metals									
Antimony	44.9		ug/L	ND	89.6	70-130			
Arsenic	57.1		ug/L	2.6	109	70-130			
Barium	71.2		ug/L	21.9	98.7	70-130			
Beryllium	53.3		ug/L	ND	106	70-130			
Boron	50.1		ug/L	ND	93.4	70-130			
Cadmium	47.2		ug/L	ND	94.2	70-130			
Chromium (VI)	4.3	0.2	ug/g	ND	74.0	70-130			
Chromium	61.5		ug/L	6.6	110	70-130			
Cobalt	55.1		ug/L	2.4	105	70-130			
Copper	67.1		ug/L	16.1	102	70-130			
Lead	61.6		ug/L	11.6	100	70-130			
Mercury	1.97	0.1	ug/g	0.259	114	70-130			
Molybdenum	52.5		ug/L	ND	105	70-130			
Nickel	57.7		ug/L	5.9	103	70-130			
Selenium	54.5		ug/L	ND	109	70-130			
Silver	49.6		ug/L	ND	99.2	70-130			
Thallium	50.5		ug/L	ND	101	70-130			
Uranium	54.2		ug/L	ND	108	70-130			
Vanadium	62.8		ug/L	ND	107	70-130			
Zinc	102		ug/L	50.1	105	70-130			
Semi-Volatiles			-						
Acenaphthene	0.180	0.02	ug/g	ND	82.4	50-140			
Acenaphthylene	0.179	0.02	ug/g	ND	82.2	50-140			
Anthracene	0.182	0.02	ug/g	ND	83.6	50-140			
Benzo [a] anthracene	0.244	0.02	ug/g	0.059	84.8	50-140			
Benzo [a] pyrene	0.233	0.02	ug/g	0.098	61.7	50-140			
Benzo [b] fluoranthene	0.430	0.02	ug/g	0.132	137	50-140			
Benzo [g,h,i] perylene	0.335	0.02	ug/g	0.089	113	50-140			
Benzo [k] fluoranthene	0.347	0.02	ug/g	0.070	127	50-140			
Chrysene	0.288	0.02	ug/g	0.086	92.5	50-140			
Dibenzo [a,h] anthracene	0.299	0.02	ug/g	ND	137	50-140			
Fluoranthene	0.306	0.02	ug/g	0.116	87.3	50-140			
Fluorene	0.170	0.02	ug/g	ND	77.9	50-140			
Indeno [1,2,3-cd] pyrene	0.362	0.02	ug/g	0.077	130	50-140			
1-Methylnaphthalene	0.152	0.02	ug/g	ND	69.7	50-140			
2-Methylnaphthalene	0.171	0.02	ug/g	ND	78.6	50-140			
Naphthalene	0.181	0.01	ug/g	ND	83.0	50-140			
Phenanthrene	0.232	0.02	ug/g	0.042	87.2	50-140			
Pyrene	0.278	0.02	ug/g	0.100	81.6	50-140			
Surrogate: 2-Fluorobiphenyl	1.34		ug/g		76.8	50-140			
Volatiles									
Acetone	9.83	0.50	ug/g		98.3	50-140			
Benzene	3.44	0.02	ug/g		85.9	60-130			
Bromodichloromethane	3.55	0.05	ug/g		88.8	60-130			



## Method Quality Control: Spike

Report Date: 30-May-2019 Order Date: 24-May-2019

Analyte	Result	Reporting Limit	Units	Source Result			RPD	RPD Limit	Notes
Bromoform	3.95	0.05	ug/g		98.9	60-130			
Bromomethane	2.47	0.05	ug/g		61.7				
Carbon Tetrachloride	3.22	0.05	ug/g		80.4				
Chlorobenzene	2.98	0.05	ug/g		74.5	60-130			
Chloroform	3.56	0.05	ug/g		89.0	60-130			
Dibromochloromethane	4.40	0.05	ug/g		110	60-130			
Dichlorodifluoromethane	2.44	0.05	ug/g		61.0	50-140			
1,2-Dichlorobenzene	3.54	0.05	ug/g		88.5	60-130			
1,3-Dichlorobenzene	3.48	0.05	ug/g		87.0	60-130			
1,4-Dichlorobenzene	2.82	0.05	ug/g		70.5	60-130			
1,1-Dichloroethane	3.39	0.05	ug/g		84.7	60-130			
1,2-Dichloroethane	3.17	0.05	ug/g		79.2	60-130			
1,1-Dichloroethylene	3.30	0.05	ug/g		82.4	60-130			
cis-1,2-Dichloroethylene	2.64	0.05	ug/g		66.0	60-130			
trans-1,2-Dichloroethylene	3.01	0.05	ug/g		75.2	60-130			
1,2-Dichloropropane	3.61	0.05	ug/g		90.1	60-130			
cis-1,3-Dichloropropylene	3.12	0.05	ug/g		78.1	60-130			
trans-1,3-Dichloropropylene	3.49	0.05	ug/g		87.3	60-130			
Ethylbenzene	2.83	0.05	ug/g		70.8	60-130			
Ethylene dibromide (dibromoethane	3.85	0.05	ug/g		96.2	60-130			
Hexane	3.68	0.05	ug/g		92.0	60-130			
Methyl Ethyl Ketone (2-Butanone)	9.05	0.50	ug/g		90.5	50-140			
Methyl Isobutyl Ketone	9.74	0.50	ug/g		97.4	50-140			
Methyl tert-butyl ether	8.00	0.05	ug/g		80.0	50-140			
Methylene Chloride	3.41	0.05	ug/g		85.3	60-130			
Styrene	2.80	0.05	ug/g		70.0	60-130			
1,1,1,2-Tetrachloroethane	3.86	0.05	ug/g		96.5	60-130			
1,1,2,2-Tetrachloroethane	4.61	0.05	ug/g		115	60-130			
Tetrachloroethylene	3.40	0.05	ug/g		85.1	60-130			
Toluene	3.61	0.05	ug/g		90.3	60-130			
1,1,1-Trichloroethane	3.39	0.05	ug/g		84.8	60-130			
1,1,2-Trichloroethane	4.29	0.05	ug/g		107	60-130			
Trichloroethylene	4.30	0.05	ug/g		108	60-130			
Trichlorofluoromethane	2.75	0.05	ug/g		68.8	50-140			
Vinyl chloride	2.94	0.02	ug/g		73.6	50-140			
m,p-Xylenes	8.49	0.05	ug/g		106	60-130			
o-Xylene	3.75	0.05	ug/g		93.7	60-130			
Benzene	3.44	0.02	ug/g		85.9	60-130			
Ethylbenzene	2.83	0.05	ug/g		70.8	60-130			
Toluene	3.61	0.05	ug/g		90.3	60-130			
m,p-Xylenes	8.49	0.05	ug/g		106	60-130			
o-Xylene	3.75	0.05	ug/g		93.7	60-130			



Page 12 of 12

#### **Qualifier Notes:**

#### Sample Qualifiers :

1: GC-FID signal did not return to baseline by C50

#### Sample Data Revisions

None

#### Work Order Revisions / Comments:

None

#### **Other Report Notes:**

n/a: not applicable ND: Not Detected MDL: Method Detection Limit Source Result: Data used as source for matrix and duplicate samples %REC: Percent recovery. RPD: Relative percent difference.

Soil results are reported on a dry weight basis when the units are denoted with 'dry'. Where %Solids is reported, moisture loss includes the loss of volatile hydrocarbons.

#### CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.

- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.
- When reported, data for F4G has been processed using a silica gel cleanup.

Paracel ID: 1921505 Head Office 300-2319 St. Laurent Blvd. Ottawa, Ontario K1G 4J8 p: 1-800-749-1947 e: paracel@paracellabs.com				ij8	Chain of Custody (Lab Use Only) . No. 122147													
LABORATORIES LTD.															Page	· / of	1	
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RELIABLE.

## Certificate of Analysis

### **Paterson Group Consulting Engineers**

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Mark D'Arcy

Client PO: 26976 Project: PE4614 Custody: 122828

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

Order #: 1929255

This Certificate of Analysis contains analytical data applicable to the following samples as submitted:

Paracel ID	Client ID
1929255-01	BH7-AU1
1929255-02	BH7-SS3
1929255-03	BH8-SS4
1929255-04	BH8-SS9
1929255-05	BH9-AU1
1929255-06	BH9-SS6
1929255-07	BH9-SS9
1929255-08	BH10-SS2

Approved By:

Dale Robertson, BSc Laboratory Director

Any use of these results implies your agreement that our total liability in connection with this work, however arising, shall be limited to the amount paid by you for this work, and that our employees or agents shall not under any circumstances be liable to you in connection with this work.



#### **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	17-Jul-19	19-Jul-19
Mercury by CVAA	EPA 7471B - CVAA, digestion	18-Jul-19	19-Jul-19
PHC F1	CWS Tier 1 - P&T GC-FID	18-Jul-19	20-Jul-19
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	17-Jul-19	19-Jul-19
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	18-Jul-19	19-Jul-19
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	17-Jul-19	22-Jul-19
REG 153: VOCs by P&T GC/MS	EPA 8260 - P&T GC-MS	18-Jul-19	19-Jul-19
Solids, %	Gravimetric, calculation	20-Jul-19	20-Jul-19

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019



Order #: 1929255

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

	Client ID: Sample Date: Sample ID: MDL/Units	BH7-AU1 15-Jul-19 10:00 1929255-01 Soil	BH7-SS3 15-Jul-19 10:00 1929255-02 Soil	BH8-SS4 15-Jul-19 11:00 1929255-03 Soil	BH8-SS9 15-Jul-19 11:00 1929255-04 Soil
Physical Characteristics	WDL/Onits		001	001	
% Solids	0.1 % by Wt.	85.0	90.8	89.5	93.1
Metals		00.0	00.0		00.1
Antimony	1.0 ug/g dry	1.5	-	-	-
Arsenic	1.0 ug/g dry	10.8	-	-	-
Barium	1.0 ug/g dry	293	-	-	-
Beryllium	0.5 ug/g dry	0.7	-	-	-
Boron	5.0 ug/g dry	7.8	-	-	-
Cadmium	0.5 ug/g dry	0.7	-	-	-
Chromium	5.0 ug/g dry	26.8	-	-	-
Chromium (VI)	0.2 ug/g dry	<0.2	-	-	-
Cobalt	1.0 ug/g dry	9.1	-	-	-
Copper	5.0 ug/g dry	77.6	-	-	-
Lead	1.0 ug/g dry	295	-	-	-
Mercury	0.1 ug/g dry	0.2	-	-	-
Molybdenum	1.0 ug/g dry	2.7	-	-	-
Nickel	5.0 ug/g dry	27.8	-	-	-
Selenium	1.0 ug/g dry	1.0	-	-	-
Silver	0.3 ug/g dry	0.3	-	-	-
Thallium	1.0 ug/g dry	<1.0	-	-	-
Uranium	1.0 ug/g dry	<1.0	-	-	-
Vanadium	10.0 ug/g dry	28.0	-	-	-
Zinc	20.0 ug/g dry	279	-	-	-
/olatiles					
Acetone	0.50 ug/g dry	-	<0.50	<0.50	<0.50
Benzene	0.02 ug/g dry	-	<0.02	<0.02	<0.02
Bromodichloromethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Bromoform	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Bromomethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Carbon Tetrachloride	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Chlorobenzene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Chloroform	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Dibromochloromethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Dichlorodifluoromethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,2-Dichlorobenzene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,3-Dichlorobenzene	0.05 ug/g dry	-	<0.05	<0.05	<0.05



Order #: 1929255

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

-	Client ID: Sample Date: Sample ID:	BH7-AU1 15-Jul-19 10:00 1929255-01	BH7-SS3 15-Jul-19 10:00 1929255-02	BH8-SS4 15-Jul-19 11:00 1929255-03	BH8-SS9 15-Jul-19 11:00 1929255-04
r	MDL/Units	Soil	Soil	Soil	Soil
1,4-Dichlorobenzene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,1-Dichloroethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,2-Dichloroethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,1-Dichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
cis-1,2-Dichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
trans-1,2-Dichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,2-Dichloropropane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
cis-1,3-Dichloropropylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
trans-1,3-Dichloropropylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,3-Dichloropropene, total	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Ethylbenzene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Ethylene dibromide (dibromoethan	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Hexane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	-	<0.50	<0.50	<0.50
Methyl Isobutyl Ketone	0.50 ug/g dry	-	<0.50	<0.50	<0.50
Methyl tert-butyl ether	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Methylene Chloride	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Styrene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Tetrachloroethylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Toluene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,1,1-Trichloroethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
1,1,2-Trichloroethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Trichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Trichlorofluoromethane	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Vinyl chloride	0.02 ug/g dry	-	<0.02	<0.02	<0.02
m,p-Xylenes	0.05 ug/g dry	-	<0.05	<0.05	<0.05
o-Xylene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Xylenes, total	0.05 ug/g dry	-	<0.05	<0.05	<0.05
4-Bromofluorobenzene	Surrogate	-	119%	106%	132%
Dibromofluoromethane	Surrogate	-	95.4%	74.2%	99.2%
Toluene-d8	Surrogate	-	83.9%	101%	89.9%
Hydrocarbons	7.49/8.4.		1	a-	
F1 PHCs (C6-C10)	7 ug/g dry	-	-	60	-
F2 PHCs (C10-C16)	4 ug/g dry	-	-	2980	-



Order #: 1929255

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

	Client ID: Sample Date: Sample ID:	15-Jul-19 10:00	BH7-SS3 15-Jul-19 10:00 1929255-02	BH8-SS4 15-Jul-19 11:00 1929255-03	BH8-SS9 15-Jul-19 11:00 1929255-04
	MDL/Units	Soil	Soil	Soil	Soil
F3 PHCs (C16-C34)	8 ug/g dry	-	-	888	-
F4 PHCs (C34-C50)	6 ug/g dry	-	-	<6	-

# PARACEL LABORATORIES LTD.

### Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 26976

Order #: 1929255

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

	Client ID: Sample Date: Sample ID: MDL/Units	BH9-AU1 15-Jul-19 12:30 1929255-05 Soil	BH9-SS6 15-Jul-19 12:30 1929255-06 Soil	BH9-SS9 15-Jul-19 12:30 1929255-07 Soil	BH10-SS2 15-Jul-19 13:30 1929255-08 Soil
Physical Characteristics					
% Solids	0.1 % by Wt.	94.2	87.8	92.5	82.3
Metals	1.0 ug/g dry		1		
Antimony		-	-	-	<1.0
Arsenic	1.0 ug/g dry	-	-	-	5.8
Barium	1.0 ug/g dry	-	-	-	171
Beryllium	0.5 ug/g dry	-	-	-	0.6
Boron	5.0 ug/g dry	-	-	-	6.7
Cadmium	0.5 ug/g dry	-	-	-	<0.5
Chromium	5.0 ug/g dry	-	-	-	19.7
Chromium (VI)	0.2 ug/g dry	-	-	-	<0.2
Cobalt	1.0 ug/g dry	-	-	-	8.1
Copper	5.0 ug/g dry	-	-	-	43.8
Lead	1.0 ug/g dry	-	-	-	255
Mercury	0.1 ug/g dry	-	-	-	0.3
Molybdenum	1.0 ug/g dry	-	-	-	1.5
Nickel	5.0 ug/g dry	-	-	-	21.0
Selenium	1.0 ug/g dry	-	-	-	<1.0
Silver	0.3 ug/g dry	-	-	-	<0.3
Thallium	1.0 ug/g dry	-	-	-	<1.0
Uranium	1.0 ug/g dry	-	-	-	<1.0
Vanadium	10.0 ug/g dry	-	-	-	23.0
Zinc	20.0 ug/g dry	-	-	-	186
Volatiles					
Acetone	0.50 ug/g dry	-	<0.50	<0.50	-
Benzene	0.02 ug/g dry	-	<0.02	<0.02	-
Bromodichloromethane	0.05 ug/g dry	-	<0.05	<0.05	-
Bromoform	0.05 ug/g dry	-	<0.05	<0.05	-
Bromomethane	0.05 ug/g dry	-	<0.05	<0.05	-
Carbon Tetrachloride	0.05 ug/g dry	-	<0.05	<0.05	-
Chlorobenzene	0.05 ug/g dry	-	<0.05	<0.05	-
Chloroform	0.05 ug/g dry	-	<0.05	<0.05	-
Dibromochloromethane	0.05 ug/g dry	-	<0.05	<0.05	-
Dichlorodifluoromethane	0.05 ug/g dry	-	<0.05	<0.05	-
1,2-Dichlorobenzene	0.05 ug/g dry	-	<0.05	<0.05	_
1,3-Dichlorobenzene	0.05 ug/g dry	-	< 0.05	<0.05	



Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

	Client ID: Sample Date: Sample ID: MDL/Units	BH9-AU1 15-Jul-19 12:30 1929255-05 Soil	BH9-SS6 15-Jul-19 12:30 1929255-06 Soil	BH9-SS9 15-Jul-19 12:30 1929255-07 Soil	BH10-SS2 15-Jul-19 13:30 1929255-08 Soil
1,4-Dichlorobenzene	0.05 ug/g dry	-	<0.05	<0.05	-
1,1-Dichloroethane	0.05 ug/g dry	-	<0.05	<0.05	-
1,2-Dichloroethane	0.05 ug/g dry	-	<0.05	<0.05	-
1,1-Dichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	-
1,2-Dichloropropane	0.05 ug/g dry	-	<0.05	<0.05	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	-	<0.05	<0.05	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	-	<0.05	<0.05	-
1,3-Dichloropropene, total	0.05 ug/g dry	-	<0.05	<0.05	-
Ethylbenzene	0.05 ug/g dry	-	<0.05	<0.05	-
Ethylene dibromide (dibromoethar	0.05 ug/g dry	-	<0.05	<0.05	-
Hexane	0.05 ug/g dry	-	<0.05	<0.05	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	-	<0.50	<0.50	-
Methyl Isobutyl Ketone	0.50 ug/g dry	-	<0.50	<0.50	-
Methyl tert-butyl ether	0.05 ug/g dry	-	<0.05	<0.05	-
Methylene Chloride	0.05 ug/g dry	-	<0.05	<0.05	-
Styrene	0.05 ug/g dry	-	<0.05	<0.05	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	-	<0.05	<0.05	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	-	<0.05	<0.05	-
Tetrachloroethylene	0.05 ug/g dry	-	<0.05	<0.05	-
Toluene	0.05 ug/g dry	-	<0.05	<0.05	-
1,1,1-Trichloroethane	0.05 ug/g dry	-	<0.05	<0.05	-
1,1,2-Trichloroethane	0.05 ug/g dry	-	<0.05	<0.05	-
Trichloroethylene	0.05 ug/g dry	-	<0.05	<0.05	-
Trichlorofluoromethane	0.05 ug/g dry	-	<0.05	<0.05	-
Vinyl chloride	0.02 ug/g dry	-	<0.02	<0.02	-
m,p-Xylenes	0.05 ug/g dry	-	0.61	<0.05	-
o-Xylene	0.05 ug/g dry	-	0.26	<0.05	-
Xylenes, total	0.05 ug/g dry	-	0.87	<0.05	-
4-Bromofluorobenzene	Surrogate	-	115%	135%	-
Dibromofluoromethane	Surrogate	-	66.7%	96.8%	-
Toluene-d8	Surrogate	-	114%	96.1%	-
Hydrocarbons					
F2 PHCs (C10-C16)	4 ug/g dry	<4	-	-	-



Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

	Client ID: Sample Date: Sample ID:	BH9-AU1 15-Jul-19 12:30 1929255-05	BH9-SS6 15-Jul-19 12:30 1929255-06	BH9-SS9 15-Jul-19 12:30 1929255-07	BH10-SS2 15-Jul-19 13:30 1929255-08
	MDL/Units	Soil	Soil	Soil	Soil
F3 PHCs (C16-C34)	8 ug/g dry	98	-	-	-
F4 PHCs (C34-C50)	6 ug/g dry	171 [1]	-	-	-
Semi-Volatiles			-	-	•
Acenaphthene	0.02 ug/g dry	-	-	-	<0.02
Acenaphthylene	0.02 ug/g dry	-	-	-	<0.02
Anthracene	0.02 ug/g dry	-	-	-	<0.02
Benzo [a] anthracene	0.02 ug/g dry	-	-	-	0.04
Benzo [a] pyrene	0.02 ug/g dry	-	-	-	0.03
Benzo [b] fluoranthene	0.02 ug/g dry	-	-	-	0.04
Benzo [g,h,i] perylene	0.02 ug/g dry	-	-	-	0.03
Benzo [k] fluoranthene	0.02 ug/g dry	-	-	-	0.03
Chrysene	0.02 ug/g dry	-	-	-	0.06
Dibenzo [a,h] anthracene	0.02 ug/g dry	-	-	-	<0.02
Fluoranthene	0.02 ug/g dry	-	-	-	0.09
Fluorene	0.02 ug/g dry	-	-	-	<0.02
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	-	-	-	<0.02
1-Methylnaphthalene	0.02 ug/g dry	-	-	-	<0.02
2-Methylnaphthalene	0.02 ug/g dry	-	-	-	<0.02
Methylnaphthalene (1&2)	0.04 ug/g dry	-	-	-	<0.04
Naphthalene	0.01 ug/g dry	-	-	-	<0.01
Phenanthrene	0.02 ug/g dry	-	-	-	0.07
Pyrene	0.02 ug/g dry	-	-	-	0.09
2-Fluorobiphenyl	Surrogate	-	-	-	69.1%
Terphenyl-d14	Surrogate	-	-	-	66.1%



Order #: 1929255

Report Date: 22-Jul-2019

Order Date: 16-Jul-2019

Project Description: PE4614

## Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g						
F2 PHCs (C10-C16)	ND	4	ug/g ug/g						
F3 PHCs (C16-C34)	ND	8	ug/g ug/g						
F4 PHCs (C34-C50)	ND	о 6	ug/g ug/g						
Metals		v	~ <del>~</del> 9						
Antimony	ND	1.0	ug/g						
Arsenic	ND	1.0	ug/g ug/g						
Barium	ND	1.0							
	ND	0.5	ug/g						
Beryllium Boron	ND ND	0.5 5.0	ug/g ug/g						
	ND ND	5.0 0.5	ug/g						
Cadmium			ug/g						
Chromium (VI)		0.2	ug/g						
Cobalt	ND	5.0	ug/g						
Cobalt	ND	1.0	ug/g						
Copper	ND	5.0	ug/g						
Lead	ND	1.0	ug/g						
Mercury	ND	0.1	ug/g						
Molybdenum	ND	1.0	ug/g						
Nickel	ND	5.0	ug/g						
Selenium	ND	1.0	ug/g						
Silver	ND	0.3	ug/g						
Thallium	ND	1.0	ug/g						
Uranium	ND	1.0	ug/g						
Vanadium	ND	10.0	ug/g						
Zinc	ND	20.0	ug/g						
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g						
Acenaphthylene	ND	0.02	ug/g						
Anthracene	ND	0.02	ug/g						
Benzo [a] anthracene	ND	0.02	ug/g						
Benzo [a] pyrene	ND	0.02	ug/g						
Benzo [b] fluoranthene	ND	0.02	ug/g						
Benzo [g,h,i] perylene	ND	0.02	ug/g						
Benzo [k] fluoranthene	ND	0.02	ug/g						
Chrysene	ND	0.02	ug/g						
Dibenzo [a,h] anthracene	ND	0.02	ug/g						
Fluoranthene	ND	0.02	ug/g						
Fluorene	ND	0.02	ug/g						
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g						
1-Methylnaphthalene	ND	0.02	ug/g						
2-Methylnaphthalene	ND	0.02	ug/g						
Methylnaphthalene (1&2)	ND	0.02	ug/g						
Naphthalene	ND	0.04	ug/g ug/g						
Phenanthrene	ND	0.02	ug/g ug/g						
Pyrene	ND	0.02	ug/g ug/g						
Surrogate: 2-Fluorobiphenyl	0.936		ug/g ug/g		70.2	50-140			
Surrogate: Terphenyl-d14	0.930		ug/g ug/g		70.2	50-140 50-140			
Volatiles	0.011		3'3						
		0.50							
Acetone	ND	0.50	ug/g						
Benzene	ND	0.02	ug/g						
Bromodichloromethane	ND	0.05	ug/g						
Bromoform	ND	0.05	ug/g						
Bromomethane	ND	0.05	ug/g						
Carbon Tetrachloride	ND	0.05	ug/g						
Chlorobenzene	ND	0.05	ug/g						
Chloroform	ND	0.05	ug/g						
Dibromochloromethane	ND	0.05	ug/g						
Dichlorodifluoromethane	ND	0.05	ug/g						



# Order #: 1929255

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

Project Description: PE4614

# Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
1,2-Dichlorobenzene	ND	0.05	ug/g						
1,3-Dichlorobenzene	ND	0.05	ug/g						
1,4-Dichlorobenzene	ND	0.05	ug/g						
1,1-Dichloroethane	ND	0.05	ug/g						
1,2-Dichloroethane	ND	0.05	ug/g						
1,1-Dichloroethylene	ND	0.05	ug/g						
cis-1,2-Dichloroethylene	ND	0.05	ug/g						
trans-1,2-Dichloroethylene	ND	0.05	ug/g						
1,2-Dichloropropane	ND	0.05	ug/g						
cis-1,3-Dichloropropylene	ND	0.05	ug/g						
trans-1,3-Dichloropropylene	ND	0.05	ug/g						
1,3-Dichloropropene, total	ND	0.05	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Ethylene dibromide (dibromoethane	ND	0.05	ug/g						
Hexane	ND	0.05	ug/g						
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g						
Methyl Isobutyl Ketone	ND	0.50	ug/g						
Methyl tert-butyl ether	ND	0.05	ug/g						
Methylene Chloride	ND	0.05	ug/g						
Styrene	ND	0.05	ug/g						
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g						
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g						
Tetrachloroethylene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
1,1,1-Trichloroethane	ND	0.05	ug/g						
1,1,2-Trichloroethane	ND	0.05	ug/g						
Trichloroethylene	ND	0.05	ug/g						
Trichlorofluoromethane	ND	0.05	ug/g						
Vinyl chloride	ND	0.02	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g						
Surrogate: 4-Bromofluorobenzene	3.46		ug/g		108	50-140			
Surrogate: Dibromofluoromethane	3.26		ug/g		102	50-140			
Surrogate: Toluene-d8	2.47		ug/g		77.1	50-140			



Report Date: 22-Jul-2019

Order Date: 16-Jul-2019

**Project Description: PE4614** 

# Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g dry	ND				40	
F2 PHCs (C10-C16)	ND	4	ug/g dry	ND				30	
F3 PHCs (C16-C34)	ND	8	ug/g dry	ND				30	
F4 PHCs (C34-C50)	ND	6	ug/g dry	ND				30	
Metals									
Antimony	1.7	1.0	ug/g dry	1.5			14.2	30	
Arsenic	8.8	1.0	ug/g dry	8.4			5.6	30	
Barium	133	1.0	ug/g dry	137			3.5	30	
Beryllium	0.8	0.5	ug/g dry	0.8			2.1	30	
Boron	12.2	5.0	ug/g dry	11.2			9.0	30	
Cadmium	3.4	0.5	ug/g dry	3.3			1.1	30	
Chromium (VI)	ND	0.2	ug/g dry	ND				35	
Chromium	28.1	5.0	ug/g dry	26.8			4.6	30	
Cobalt	8.7	1.0	ug/g dry	8.6			1.9	30	
Copper	61.8	5.0	ug/g dry	60.9			1.4	30	
Lead	175	1.0	ug/g dry	103			51.9	30	
Mercury	ND	0.1	ug/g dry	0.113			0.0	30	
Molybdenum	2.0	1.0	ug/g dry	1.6			20.3	30	
Nickel	25.6	5.0	ug/g dry	24.0			6.4	30	
Selenium	ND	1.0	ug/g dry	ND			0.0	30	
Silver	ND	0.3	ug/g dry	ND			0.0	30	
Thallium	ND	1.0	ug/g dry	ND			0.0	30	
Uranium	ND	1.0	ug/g dry	ND			0.0	30	
Vanadium	33.8	10.0	ug/g dry	32.2			4.9	30	
Zinc	206	20.0	ug/g dry	193			6.5	30	
Physical Characteristics									
% Solids	83.6	0.1	% by Wt.	85.0			1.6	25	
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g dry	ND				40	
Acenaphthylene	ND	0.02	ug/g dry	ND			0.0	40	
Anthracene	ND	0.02	ug/g dry	ND			0.0	40	
Benzo [a] anthracene	0.040	0.02	ug/g dry	0.044			9.4	40	
Benzo [a] pyrene	0.032	0.02	ug/g dry	0.034			4.5	40	
Benzo [b] fluoranthene	0.039	0.02	ug/g dry	0.042			7.3	40	
Benzo [g,h,i] perylene	0.024	0.02	ug/g dry	0.025			5.4	40	
Benzo [k] fluoranthene	ND	0.02	ug/g dry	0.026			0.0	40	
Chrysene	0.048	0.02	ug/g dry	0.056			14.7	40	
Dibenzo [a,h] anthracene	ND	0.02	ug/g dry	ND				40	
Fluoranthene	0.082	0.02	ug/g dry	0.094			13.5	40	
Fluorene	ND	0.02	ug/g dry	ND			0.0	40	
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g dry	ND			0.0	40	
1-Methylnaphthalene	ND	0.02	ug/g dry	ND				40	
2-Methylnaphthalene	ND	0.02	ug/g dry	ND				40	
Naphthalene	ND	0.01	ug/g dry	ND				40	
Phenanthrene	0.047	0.02	ug/g dry	0.071			42.2		QR-01
Pyrene	0.074	0.02	ug/g dry	0.088			17.8	40	
Surrogate: 2-Fluorobiphenyl	0.868		ug/g dry		53.6	50-140			
Surrogate: Terphenyl-d14	0.840		ug/g dry		51.9	50-140			
Volatiles									
Acetone	ND	0.50	ug/g dry	ND				50	
Benzene	ND	0.02	ug/g dry	ND				50	
Bromodichloromethane	ND	0.05	ug/g dry	ND				50	
Bromoform	ND	0.05	ug/g dry	ND				50	
Bromomethane	ND	0.05	ug/g dry	ND				50	
Carbon Tetrachloride	ND	0.05	ug/g dry	ND				50	
Chlorobenzene	ND	0.05	ug/g dry	ND				50	



### Certificate of Analysis **Client: Paterson Group Consulting Engineers** Client PO: 26976

# Method Quality

		Reporting		Source		%REC		RPD	RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes	
Chloroform	ND	0.05	ug/g dry	ND				50		
Dibromochloromethane	ND	0.05	ug/g dry	ND				50		
Dichlorodifluoromethane	ND	0.05	ug/g dry	ND				50		
1,2-Dichlorobenzene	ND	0.05	ug/g dry	ND				50		
1,3-Dichlorobenzene	ND	0.05	ug/g dry	ND				50		
1,4-Dichlorobenzene	ND	0.05	ug/g dry	ND				50		
1,1-Dichloroethane	ND	0.05	ug/g dry	ND				50		
1,2-Dichloroethane	ND	0.05	ug/g dry	ND				50		
1,1-Dichloroethylene	ND	0.05	ug/g dry	ND				50		
cis-1,2-Dichloroethylene	ND	0.05	ug/g dry	ND				50		
trans-1,2-Dichloroethylene	ND	0.05	ug/g dry	ND				50		
1,2-Dichloropropane	ND	0.05	ug/g dry	ND				50		
cis-1,3-Dichloropropylene	ND	0.05	ug/g dry	ND				50		

1,2-Dichloroethane	ND	0.05	ug/g ury	ND			50	
1,1-Dichloroethylene	ND	0.05	ug/g dry	ND			50	
cis-1,2-Dichloroethylene	ND	0.05	ug/g dry	ND			50	
trans-1,2-Dichloroethylene	ND	0.05	ug/g dry	ND			50	
1,2-Dichloropropane	ND	0.05	ug/g dry	ND			50	
cis-1,3-Dichloropropylene	ND	0.05	ug/g dry	ND			50	
trans-1,3-Dichloropropylene	ND	0.05	ug/g dry	ND			50	
Ethylbenzene	ND	0.05	ug/g dry	ND			50	
Ethylene dibromide (dibromoethane	ND	0.05	ug/g dry	ND			50	
Hexane	ND	0.05	ug/g dry	ND			50	
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g dry	ND			50	
Methyl Isobutyl Ketone	ND	0.50	ug/g dry	ND			50	
Methyl tert-butyl ether	ND	0.05	ug/g dry	ND			50	
Methylene Chloride	ND	0.05	ug/g dry	ND			50	
Styrene	ND	0.05	ug/g dry	ND			50	
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g dry	ND			50	
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g dry	ND			50	
Tetrachloroethylene	ND	0.05	ug/g dry	ND			50	
Toluene	ND	0.05	ug/g dry	ND			50	
1,1,1-Trichloroethane	ND	0.05	ug/g dry	ND			50	
1,1,2-Trichloroethane	ND	0.05	ug/g dry	ND			50	
Trichloroethylene	ND	0.05	ug/g dry	ND			50	
Trichlorofluoromethane	ND	0.05	ug/g dry	ND			50	
Vinyl chloride	ND	0.02	ug/g dry	ND			50	
m,p-Xylenes	ND	0.05	ug/g dry	ND			50	
o-Xylene	ND	0.05	ug/g dry	ND			50	
Surrogate: 4-Bromofluorobenzene	4.34		ug/g dry		123	50-140		
Surrogate: Dibromofluoromethane	3.53		ug/g dry		100	50-140		
Surrogate: Toluene-d8	2.98		ug/g dry		84.6	50-140		

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019



# Method Quality Control: Spike

Report Date: 22-Jul-2019

Order Date: 16-Jul-2019

**Project Description: PE4614** 

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	196	7	ug/g		97.9	80-120			
F2 PHCs (C10-C16)	96	4	ug/g	ND	111	60-140			
F3 PHCs (C16-C34)	282	8	ug/g	ND	133	60-140			
F4 PHCs (C34-C50)	164	6	ug/g	ND	122	60-140			
Metals									
Antimony	39.3		ug/L	ND	77.4	70-130			
Arsenic	52.3		ug/L	3.3	98.0	70-130			
Barium	105		ug/L	54.9	101	70-130			
Beryllium	47.4		ug/L	ND	94.2	70-130			
Boron	47.3		ug/L	ND	85.7	70-130			
Cadmium	47.4		ug/L	1.3	92.1	70-130			
Chromium (VI)	3.7	0.2	ug/g		74.5	70-130			
Chromium	59.8		ug/L	10.7	98.1	70-130			
Cobalt	51.1		ug/L	3.4	95.3	70-130			
Copper	72.8		ug/L	24.4	97.0	70-130			
Lead	87.4		ug/L	41.3	92.2	70-130			
Mercury	1.51	0.1	ug/g	0.113	93.2	70-130			
Molybdenum	47.4		ug/L	ND	93.5	70-130			
Nickel	56.7		ug/L	9.6	94.2	70-130			
Selenium	48.4		ug/L	ND	96.2	70-130			
Silver	39.4		ug/L	ND	78.7	70-130			
Thallium	44.3		ug/L	ND	88.4	70-130			
Uranium	46.6		ug/L	ND	92.5	70-130			
Vanadium	61.9		ug/L	12.9	98.1	70-130			
Zinc	124		ug/L	77.2	94.3	70-130			
Semi-Volatiles									
Acenaphthene	0.134	0.02	ug/g	ND	66.3	50-140			
Acenaphthylene	0.128	0.02	ug/g	ND	63.2	50-140			
Anthracene	0.150	0.02	ug/g	ND	74.1	50-140			
Benzo [a] anthracene	0.187	0.02	ug/g	0.044	70.6	50-140			
Benzo [a] pyrene	0.146	0.02	ug/g	0.034	55.2	50-140			
Benzo [b] fluoranthene	0.236	0.02	ug/g	0.042	96.3	50-140			
Benzo [g,h,i] perylene	0.130	0.02	ug/g	0.025	51.7	50-140			
Benzo [k] fluoranthene	0.182	0.02	ug/g	0.026	77.2	50-140			
Chrysene	0.236	0.02	ug/g	0.056	89.0	50-140			
Dibenzo [a,h] anthracene	0.099	0.02	ug/g		59.2	50-140			
Fluoranthene	0.247	0.02	ug/g	0.094	75.5	50-140			
Fluorene	0.125	0.02	ug/g	ND	61.8	50-140			
Indeno [1,2,3-cd] pyrene	0.125	0.02	ug/g	ND	61.8	50-140			
1-Methylnaphthalene	0.140	0.02	ug/g	ND	69.0	50-140			
2-Methylnaphthalene	0.152	0.02	ug/g	ND	75.2	50-140			
Naphthalene	0.133	0.01	ug/g	ND	65.7	50-140			
Phenanthrene	0.219	0.02	ug/g	0.071	73.1	50-140			
Pyrene	0.245	0.02	ug/g	0.088	77.6	50-140			
Surrogate: 2-Fluorobiphenyl	0.998		ug/g		61.6	50-140			
Volatiles	10.0	0.50			400	50 110			
Acetone	13.0	0.50	ug/g		130	50-140			
Benzene	4.13	0.02	ug/g		103	60-130			
Bromodichloromethane	4.50	0.05	ug/g		112	60-130			
Bromoform	4.08	0.05	ug/g		102	60-130			



# Method Quality Control: Spike

Report Date: 22-Jul-2019 Order Date: 16-Jul-2019

**Project Description: PE4614** 

%REC Limit	RPD	RPD Limit	Notes

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Bromomethane	2.37	0.05	ug/g		59.2	50-140			
Carbon Tetrachloride	3.97	0.05	ug/g		99.3	60-130			
Chlorobenzene	3.78	0.05	ug/g		94.5	60-130			
Chloroform	4.73	0.05	ug/g		118	60-130			
Dibromochloromethane	4.06	0.05	ug/g		101	60-130			
Dichlorodifluoromethane	2.49	0.05	ug/g		62.2	50-140			
1,2-Dichlorobenzene	2.84	0.05	ug/g		71.0	60-130			
1,3-Dichlorobenzene	2.96	0.05	ug/g		74.1	60-130			
1,4-Dichlorobenzene	2.93	0.05	ug/g		73.2	60-130			
1,1-Dichloroethane	4.73	0.05	ug/g		118	60-130			
1,2-Dichloroethane	4.82	0.05	ug/g		120	60-130			
1,1-Dichloroethylene	3.91	0.05	ug/g		97.9	60-130			
cis-1,2-Dichloroethylene	4.16	0.05	ug/g		104	60-130			
trans-1,2-Dichloroethylene	3.70	0.05	ug/g		92.5	60-130			
1,2-Dichloropropane	4.67	0.05	ug/g		117	60-130			
cis-1,3-Dichloropropylene	2.74	0.05	ug/g		68.6	60-130			
trans-1,3-Dichloropropylene	3.17	0.05	ug/g		79.1	60-130			
Ethylbenzene	4.19	0.05	ug/g		105	60-130			
Ethylene dibromide (dibromoethane	4.01	0.05	ug/g		100	60-130			
Hexane	3.96	0.05	ug/g		99.1	60-130			
Methyl Ethyl Ketone (2-Butanone)	12.0	0.50	ug/g		120	50-140			
Methyl Isobutyl Ketone	11.4	0.50	ug/g		114	50-140			
Methyl tert-butyl ether	9.57	0.05	ug/g		95.7	50-140			
Methylene Chloride	5.12	0.05	ug/g		128	60-130			
Styrene	4.17	0.05	ug/g		104	60-130			
1,1,1,2-Tetrachloroethane	3.95	0.05	ug/g		98.9	60-130			
1,1,2,2-Tetrachloroethane	4.32	0.05	ug/g		108	60-130			
Tetrachloroethylene	3.27	0.05	ug/g		81.8	60-130			
Toluene	3.99	0.05	ug/g		99.7	60-130			
1,1,1-Trichloroethane	3.89	0.05	ug/g		97.4	60-130			
1,1,2-Trichloroethane	4.87	0.05	ug/g		122	60-130			
Trichloroethylene	3.61	0.05	ug/g		90.3	60-130			
Trichlorofluoromethane	3.61	0.05	ug/g		90.4	50-140			
Vinyl chloride	2.68	0.02	ug/g		67.1	50-140			
m,p-Xylenes	7.76	0.05	ug/g		97.1	60-130			
o-Xylene	4.48	0.05	ug/g		112	60-130			



### Sample Qualifiers :

1: GC-FID signal did not return to baseline by C50

### QC Qualifiers :

QR-01: Duplicate RPD is high, however, the sample result is less than 10x the MDL.

### Sample Data Revisions

None

### Work Order Revisions / Comments:

None

### **Other Report Notes:**

n/a: not applicable ND: Not Detected MDL: Method Detection Limit Source Result: Data used as source for matrix and duplicate samples %REC: Percent recovery. RPD: Relative percent difference.

Soil results are reported on a dry weight basis when the units are denoted with 'dry'. Where %Solids is reported, moisture loss includes the loss of volatile hydrocarbons.

### CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.

- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.
- When reported, data for F4G has been processed using a silica gel cleanup.

LABORATORIES LTD. Client Name: Poterson Contact Name: Mork D'Arsy Address: Telephone: 226 - 7381 Criteria: DO. Reg. 153/04 (As Amended) Table _ CRSC Fil		0. Reg		Project Reference Quote # PO #	e. j	PE	.4	61	4						10.0	: 0	(	
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Paracel Order Number: 929255 Sample ID/Location Name 1 T3 H 7-AU1 2 T3 H 7-SS3	S S Matrix	Air Volume	2 A # of Containers	Sample Date Jy 15 2019	Time Time To am	PHCs F1-F4+	VOCS	PAHs Metals by ICP		< GNI	B (HWS)	PH CA)			120 Misore		Ja(	
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Chain of Custody (Env) - Rev 0.7 Feb. 2016



RELIABLE.

300 - 2319 St. Laurent Blvd Ottawa, ON, K1G 4J8 1-800-749-1947 www.paracellabs.com

# Certificate of Analysis

### **Paterson Group Consulting Engineers**

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Mark D'Arcy

Client PO: 30555 Project: PE4614 Custody: 125769

Report Date: 2-Sep-2020 Order Date: 27-Aug-2020

Order #: 2035556

This Certificate of Analysis contains analytical data applicable to the following samples as submitted:

Paracel ID	Client ID
2035556-01	BH11-20-AU1
2035556-02	BH11-20-SS2
2035556-03	BH11-20-SS4
2035556-04	BH13-20-AU1
2035556-05	BH13-20-SS3
2035556-06	BH13-20-SS7
2035556-07	BH14-20-SS2

Approved By:

Dale Robertson, BSc Laboratory Director

Any use of these results implies your agreement that our total liability in connection with this work, however arising, shall be limited to the amount paid by you for this work, and that our employees or agents shall not under any circumstances be liable to you in connection with this work.



### **Analysis Summary Table**

Order #	<b>#: 2035556</b>
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Report Date: 02-Sep-2020 Order Date: 27-Aug-2020

Project Description: PE4614

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	28-Aug-20	28-Aug-20
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	28-Aug-20	29-Aug-20
Mercury by CVAA	EPA 7471B - CVAA, digestion	1-Sep-20	1-Sep-20
PHC F1	CWS Tier 1 - P&T GC-FID	28-Aug-20	28-Aug-20
PHC F4G (gravimetric)	CWS Tier 1 - Extraction Gravimetric	1-Sep-20	2-Sep-20
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	28-Aug-20	2-Sep-20
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	31-Aug-20	31-Aug-20
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	28-Aug-20	1-Sep-20
REG 153: VOCs by P&T GC/MS	EPA 8260 - P&T GC-MS	28-Aug-20	28-Aug-20
Solids, %	Gravimetric, calculation	28-Aug-20	31-Aug-20



### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 30555

Report Date: 02-Sep-2020

Order Date: 27-Aug-2020

Project Description: PE4614

	Client ID: Sample Date: Sample ID: MDL/Units	BH11-20-AU1 24-Aug-20 09:00 2035556-01 Soil	BH11-20-SS2 24-Aug-20 09:00 2035556-02 Soil	BH11-20-SS4 24-Aug-20 09:00 2035556-03 Soil	BH13-20-AU1 24-Aug-20 09:00 2035556-04 Soil
Physical Characteristics					
% Solids	0.1 % by Wt.	94.8	85.3	90.8	91.6
Metals	•		•		
Antimony	1.0 ug/g dry	-	<1.0	-	-
Arsenic	1.0 ug/g dry	-	8.2	-	-
Barium	1.0 ug/g dry	-	180	-	-
Beryllium	0.5 ug/g dry	-	0.8	-	-
Boron	5.0 ug/g dry	-	9.0	-	-
Cadmium	0.5 ug/g dry	-	<0.5	-	-
Chromium	5.0 ug/g dry	-	30.4	-	-
Chromium (VI)	0.2 ug/g dry	-	<0.2	-	-
Cobalt	1.0 ug/g dry	-	9.5	-	-
Copper	5.0 ug/g dry	-	63.1	-	-
Lead	1.0 ug/g dry	-	226	-	-
Mercury	0.1 ug/g dry	-	1.4	-	-
Molybdenum	1.0 ug/g dry	-	2.2	-	-
Nickel	5.0 ug/g dry	-	34.8	-	-
Selenium	1.0 ug/g dry	-	<1.0	-	-
Silver	0.3 ug/g dry	-	1.1	-	-
Thallium	1.0 ug/g dry	-	<1.0	-	-
Uranium	1.0 ug/g dry	-	2.4	-	-
Vanadium	10.0 ug/g dry	-	30.5	-	-
Zinc	20.0 ug/g dry	-	175	-	-
Volatiles	+		4	ł	• •
Benzene	0.02 ug/g dry	-	-	<0.02	-
Ethylbenzene	0.05 ug/g dry	-	-	<0.05	-
Toluene	0.05 ug/g dry	-	-	<0.05	-
m,p-Xylenes	0.05 ug/g dry	-	-	<0.05	-
o-Xylene	0.05 ug/g dry	-	-	<0.05	-
Xylenes, total	0.05 ug/g dry	-	-	<0.05	-
Toluene-d8	Surrogate	-	-	108%	-
Hydrocarbons					
F1 PHCs (C6-C10)	7 ug/g dry	-	-	<7	-
F2 PHCs (C10-C16)	4 ug/g dry	<40 [1]	-	35	5
F3 PHCs (C16-C34)	8 ug/g dry	131	-	83	54
F4 PHCs (C34-C50)	6 ug/g dry	877 [2]	-	36	57

# PARACEL LABORATORIES LTD.

# Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 30555

Report Date: 02-Sep-2020 Order Date: 27-Aug-2020

	Client ID: Sample Date: Sample ID:	BH11-20-AU1 24-Aug-20 09:00 2035556-01 Soil	BH11-20-SS2 24-Aug-20 09:00 2035556-02 Soil	BH11-20-SS4 24-Aug-20 09:00 2035556-03 Soil	BH13-20-AU1 24-Aug-20 09:00 2035556-04 Soil
F4G PHCs (gravimetric)	MDL/Units 50 ug/g dry	1730		-	-
Semi-Volatiles		1760	ļ -	ļ	_
Acenaphthene	0.02 ug/g dry	-	<0.02	-	-
Acenaphthylene	0.02 ug/g dry	-	0.05	-	-
Anthracene	0.02 ug/g dry	-	0.05	-	-
Benzo [a] anthracene	0.02 ug/g dry	-	0.13	-	-
Benzo [a] pyrene	0.02 ug/g dry	-	0.14	-	-
Benzo [b] fluoranthene	0.02 ug/g dry	-	0.19	-	-
Benzo [g,h,i] perylene	0.02 ug/g dry	-	0.10	-	-
Benzo [k] fluoranthene	0.02 ug/g dry	-	0.10	-	-
Chrysene	0.02 ug/g dry	-	0.13	-	-
Dibenzo [a,h] anthracene	0.02 ug/g dry	-	0.03	-	-
Fluoranthene	0.02 ug/g dry	-	0.27	-	-
Fluorene	0.02 ug/g dry	-	<0.02	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	-	0.09	-	-
1-Methylnaphthalene	0.02 ug/g dry	-	<0.02	-	-
2-Methylnaphthalene	0.02 ug/g dry	-	<0.02	-	-
Methylnaphthalene (1&2)	0.04 ug/g dry	-	<0.04	-	-
Naphthalene	0.01 ug/g dry	-	<0.01	-	-
Phenanthrene	0.02 ug/g dry	-	0.15	-	-
Pyrene	0.02 ug/g dry	-	0.24	-	-
2-Fluorobiphenyl	Surrogate	-	92.6%	-	-
Terphenyl-d14	Surrogate	-	110%	-	-

# PARACEL LABORATORIES LTD.

### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 30555

Order #: 2035556

Report Date: 02-Sep-2020

Order Date: 27-Aug-2020

Project Description: PE4614

	Client ID: Sample Date: Sample ID: MDL/Units	BH13-20-SS3 24-Aug-20 09:00 2035556-05 Soil	BH13-20-SS7 24-Aug-20 09:00 2035556-06 Soil	BH14-20-SS2 24-Aug-20 09:00 2035556-07 Soil	- - - -
Physical Characteristics					
% Solids	0.1 % by Wt.	96.7	89.8	94.0	-
Metals					·
Antimony	1.0 ug/g dry	<1.0	-	3.9	-
Arsenic	1.0 ug/g dry	2.0	-	2.3	-
Barium	1.0 ug/g dry	14.8	-	99.5	-
Beryllium	0.5 ug/g dry	<0.5	-	<0.5	-
Boron	5.0 ug/g dry	<5.0	-	7.6	-
Cadmium	0.5 ug/g dry	<0.5	-	<0.5	-
Chromium	5.0 ug/g dry	7.9	-	13.7	-
Chromium (VI)	0.2 ug/g dry	<0.2	-	0.3	-
Cobalt	1.0 ug/g dry	3.5	-	5.4	-
Copper	5.0 ug/g dry	8.0	-	18.4	-
Lead	1.0 ug/g dry	3.6	-	22.7	-
Mercury	0.1 ug/g dry	<0.1	-	0.2	-
Molybdenum	1.0 ug/g dry	<1.0	-	<1.0	-
Nickel	5.0 ug/g dry	10.1	-	11.6	-
Selenium	1.0 ug/g dry	<1.0	-	<1.0	-
Silver	0.3 ug/g dry	<0.3	-	<0.3	-
Thallium	1.0 ug/g dry	<1.0	-	<1.0	-
Uranium	1.0 ug/g dry	<1.0	-	<1.0	-
Vanadium	10.0 ug/g dry	14.6	-	16.5	-
Zinc	20.0 ug/g dry	<20.0	-	36.6	-
Volatiles	+ +				
Acetone	0.50 ug/g dry	-	<0.50	-	-
Benzene	0.02 ug/g dry	-	<0.02	-	-
Bromodichloromethane	0.05 ug/g dry	-	<0.05	-	-
Bromoform	0.05 ug/g dry	-	<0.05	-	-
Bromomethane	0.05 ug/g dry	-	<0.05	-	-
Carbon Tetrachloride	0.05 ug/g dry	-	<0.05	-	-
Chlorobenzene	0.05 ug/g dry	-	<0.05	-	-
Chloroform	0.05 ug/g dry	-	<0.05	-	-
Dibromochloromethane	0.05 ug/g dry	-	<0.05	-	-
Dichlorodifluoromethane	0.05 ug/g dry	-	<0.05	-	-
1,2-Dichlorobenzene	0.05 ug/g dry	-	<0.05	-	-
1,3-Dichlorobenzene	0.05 ug/g dry	-	<0.05	-	-



Order #: 2035556

Report Date: 02-Sep-2020 Order Date: 27-Aug-2020

Project Description: PE4614

	Client ID: Sample Date: Sample ID: MDL/Units	BH13-20-SS3 24-Aug-20 09:00 2035556-05 Soil	BH13-20-SS7 24-Aug-20 09:00 2035556-06 Soil	BH14-20-SS2 24-Aug-20 09:00 2035556-07 Soil	- - - -
1,4-Dichlorobenzene	0.05 ug/g dry	-	<0.05	-	-
1,1-Dichloroethane	0.05 ug/g dry	-	<0.05	-	-
1,2-Dichloroethane	0.05 ug/g dry	-	<0.05	-	-
1,1-Dichloroethylene	0.05 ug/g dry	-	<0.05	-	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	-	<0.05	-	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	-	<0.05	-	-
1,2-Dichloropropane	0.05 ug/g dry	-	<0.05	-	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	-	<0.05	-	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	-	<0.05	-	-
1,3-Dichloropropene, total	0.05 ug/g dry	-	<0.05	-	-
Ethylbenzene	0.05 ug/g dry	-	<0.05	-	-
Ethylene dibromide (dibromoethane, 1	0.05 ug/g dry	-	<0.05	-	-
Hexane	0.05 ug/g dry	-	<0.05	-	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	-	<0.50	-	-
Methyl Isobutyl Ketone	0.50 ug/g dry	-	<0.50	-	-
Methyl tert-butyl ether	0.05 ug/g dry	-	<0.05	-	-
Methylene Chloride	0.05 ug/g dry	-	<0.05	-	-
Styrene	0.05 ug/g dry	-	<0.05	-	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	-	<0.05	-	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	-	<0.05	-	-
Tetrachloroethylene	0.05 ug/g dry	-	<0.05	-	-
Toluene	0.05 ug/g dry	-	<0.05	-	-
1,1,1-Trichloroethane	0.05 ug/g dry	-	<0.05	-	-
1,1,2-Trichloroethane	0.05 ug/g dry	-	<0.05	-	-
Trichloroethylene	0.05 ug/g dry	-	<0.05	-	-
Trichlorofluoromethane	0.05 ug/g dry	-	<0.05	-	-
Vinyl chloride	0.02 ug/g dry	-	<0.02	-	-
m,p-Xylenes	0.05 ug/g dry	-	<0.05	-	-
o-Xylene	0.05 ug/g dry	-	<0.05	-	-
Xylenes, total	0.05 ug/g dry	-	<0.05	-	-
4-Bromofluorobenzene	Surrogate	-	111%	-	-
Dibromofluoromethane	Surrogate	-	98.5%	-	-
Toluene-d8	Surrogate	-	96.2%	-	-
Hydrocarbons			<b></b>	<b> </b>	<u>ـــــــ</u> ا
F2 PHCs (C10-C16)	4 ug/g dry	<4	-	-	-



Report Date: 02-Sep-2020 Order Date: 27-Aug-2020

Project Description: PE4614

	_				
	Client ID:	BH13-20-SS3	BH13-20-SS7	BH14-20-SS2	-
	Sample Date:	24-Aug-20 09:00	24-Aug-20 09:00	24-Aug-20 09:00	-
	Sample ID:	2035556-05	2035556-06	2035556-07	-
	MDL/Units	Soil	Soil	Soil	-
F3 PHCs (C16-C34)	8 ug/g dry	<8	-	-	-
F4 PHCs (C34-C50)	6 ug/g dry	<6	-	-	-



Dibromochloromethane

Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 30555

### Method Quality Control: Blank

Report Date: 02-Sep-2020

Order Date: 27-Aug-2020

Project Description: PE4614

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g						
F2 PHCs (C10-C16)	ND	4	ug/g						
F3 PHCs (C16-C34)	ND	8	ug/g						
F4 PHCs (C34-C50)	ND	6	ug/g						
F4G PHCs (gravimetric)	ND	50	ug/g						
Metals									
Antimony	ND	1.0	ug/g						
Arsenic	ND	1.0	ug/g						
Barium	ND	1.0	ug/g						
Beryllium	ND	0.5	ug/g						
Boron	ND	5.0	ug/g						
Cadmium	ND	0.5	ug/g						
Chromium (VI)	ND	0.2	ug/g						
Chromium	ND	5.0	ug/g						
Cobalt	ND	1.0	ug/g						
Copper	ND	5.0	ug/g						
Lead	ND	1.0	ug/g						
Mercury	ND	0.1	ug/g						
Molybdenum	ND	1.0	ug/g						
Nickel	ND	5.0	ug/g						
Selenium	ND	1.0	ug/g						
Silver	ND	0.3	ug/g						
Thallium	ND	1.0	ug/g						
Uranium	ND	1.0	ug/g						
Vanadium	ND	10.0	ug/g						
Zinc	ND	20.0	ug/g						
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g						
Acenaphthylene	ND	0.02	ug/g						
Anthracene	ND	0.02	ug/g						
Benzo [a] anthracene	ND	0.02	ug/g						
Benzo [a] pyrene	ND	0.02	ug/g						
Benzo [b] fluoranthene	ND	0.02	ug/g						
Benzo [g,h,i] perylene	ND	0.02	ug/g						
Benzo [k] fluoranthene	ND	0.02	ug/g						
Chrysene	ND	0.02	ug/g						
Dibenzo [a,h] anthracene	ND	0.02	ug/g						
Fluoranthene	ND	0.02	ug/g						
Fluorene	ND	0.02	ug/g						
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g						
1-Methylnaphthalene	ND	0.02	ug/g						
2-Methylnaphthalene	ND	0.02	ug/g						
Methylnaphthalene (1&2)	ND	0.04	ug/g						
Naphthalene	ND	0.01	ug/g						
Phenanthrene	ND	0.02	ug/g						
Pyrene	ND	0.02	ug/g			50 4 40			
Surrogate: 2-Fluorobiphenyl	1.26		ug/g		94.6	50-140			
Surrogate: Terphenyl-d14	1.57		ug/g		118	50-140			
Volatiles									
Acetone	ND	0.50	ug/g						
Benzene	ND	0.02	ug/g						
Bromodichloromethane	ND	0.05	ug/g						
Bromoform	ND	0.05	ug/g						
Bromomethane	ND	0.05	ug/g						
Carbon Tetrachloride	ND	0.05	ug/g						
Chlorobenzene	ND	0.05	ug/g						
Chloroform	ND	0.05	ug/g						
Dibromochloromethane	ND	0.05	ua/a						

OTTAWA • MISSISSAUGA • HAMILTON • CALGARY • KINGSTON • LONDON • NIAGARA • WINDSOR • RICHMOND HILL

ug/g

ND

0.05



### Method Quality Control: Blank

Report Date: 02-Sep-2020

Order Date: 27-Aug-2020

Project Description: PE4614

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Dichlorodifluoromethane	ND	0.05	ug/g						
1,2-Dichlorobenzene	ND	0.05	ug/g						
1,3-Dichlorobenzene	ND	0.05	ug/g						
1,4-Dichlorobenzene	ND	0.05	ug/g						
1,1-Dichloroethane	ND	0.05	ug/g						
1,2-Dichloroethane	ND	0.05	ug/g						
1,1-Dichloroethylene	ND	0.05	ug/g						
cis-1,2-Dichloroethylene	ND	0.05	ug/g						
trans-1,2-Dichloroethylene	ND	0.05	ug/g						
1,2-Dichloropropane	ND	0.05	ug/g						
cis-1,3-Dichloropropylene	ND	0.05	ug/g						
trans-1,3-Dichloropropylene	ND	0.05	ug/g						
1,3-Dichloropropene, total	ND	0.05	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Ethylene dibromide (dibromoethane, 1,2	ND	0.05	ug/g						
Hexane	ND	0.05	ug/g						
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g						
Methyl Isobutyl Ketone	ND	0.50	ug/g						
Methyl tert-butyl ether	ND	0.05	ug/g						
Methylene Chloride	ND	0.05	ug/g						
Styrene	ND	0.05	ug/g						
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g						
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g						
Tetrachloroethylene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
1,1,1-Trichloroethane	ND	0.05	ug/g						
1,1,2-Trichloroethane	ND	0.05	ug/g						
Trichloroethylene Trichlorofluoromethane	ND ND	0.05 0.05	ug/g						
		0.05	ug/g						
Vinyl chloride m,p-Xylenes	ND ND	0.02	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g						
Surrogate: 4-Bromofluorobenzene	3.56	0.05	ug/g <i>ug/g</i>		111	50-140			
-	3.30				102	50-140 50-140			
Surrogate: Dibromofluoromethane			ug/g						
Surrogate: Toluene-d8	3.79		ug/g		118	50-140			
Benzene	ND	0.02	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g		110	E0 140			
Surrogate: Toluene-d8	3.79		ug/g		118	50-140			



Client PO: 30555

Report Date: 02-Sep-2020 Order Date: 27-Aug-2020

Project Description: PE4614

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g dry	ND			NC	40	
F2 PHCs (C10-C16)	ND	4	ug/g dry	ND			NC	30	
F3 PHCs (C16-C34)	ND	8	ug/g dry	ND			NC	30	
F4 PHCs (C34-C50)	ND	6	ug/g dry	ND			NC	30	
Metals			. – .						
Antimony	1.4	1.0	ug/g dry	ND			NC	30	
Arsenic	10.0	1.0	ug/g dry	8.2			19.8	30	
Barium	198	1.0	ug/g dry	180			9.3	30	
Beryllium	0.9	0.5	ug/g dry	0.8			4.6	30	
Boron	10.7	5.0	ug/g dry	9.0			18.2	30	
Cadmium	ND	0.5	ug/g dry	ND			NC	30 25	
Chromium (VI)	ND 24-2	0.2	ug/g dry	ND			NC	35	
Chromium	34.2	5.0	ug/g dry	30.4			11.8 11.5	30 30	
Cobalt	10.7	1.0	ug/g dry	9.5 63.1			11.5 10.0	30 30	
Copper	69.8 230	5.0 1.0	ug/g dry ug/g dry	63.1 226			10.0 1 4	30 30	
Lead Mercury	230 1.58	1.0 0.1	ug/g dry ug/g dry	226 1 35			1.4 15.9	30 30	
Mercury Molybdenum	1.58 2.8	0.1 1.0	ug/g dry ug/g dry	1.35 2.2			15.9 24.2	30 30	
Nolybaenum Nickel	2.8 39.1	1.0 5.0	ug/g dry ug/g dry	2.2 34.8			24.2 11.7	30 30	
Selenium	39.1 1.1	5.0 1.0	ug/g ary ug/g dry	34.8 ND			NC	30 30	
Silver	1.1	0.3	ug/g dry ug/g dry	1.1			18.1	30 30	
Thallium	ND	0.3 1.0	ug/g dry ug/g dry	ND			NC	30	
Uranium	2.6	1.0	ug/g dry ug/g dry	2.4			10.5	30	
Vanadium	34.9	10.0	ug/g dry ug/g dry	30.5			13.5	30	
Zinc	193	20.0	ug/g dry ug/g dry	175			10.0	30	
Physical Characteristics		_0.0							
% Solids	95.5	0.1	% by Wt.	94.8			0.7	25	
Semi-Volatiles			-,						
Acenaphthene	ND	0.02	ug/g dry	ND			NC	40	
Acenaphthylene	0.044	0.02	ug/g dry	0.051			16.0	40	
Anthracene	0.034	0.02	ug/g dry	0.053			NC	40	
Benzo [a] anthracene	0.088	0.02	ug/g dry	0.130			38.9	40	
Benzo [a] pyrene	0.104	0.02	ug/g dry	0.144			32.3	40	
Benzo [b] fluoranthene	0.143	0.02	ug/g dry	0.187			27.1	40	
Benzo [g,h,i] perylene	0.082	0.02	ug/g dry	0.099			19.5	40	
Benzo [k] fluoranthene	0.063	0.02	ug/g dry	0.098			NC	40	
Chrysene	0.101	0.02	ug/g dry	0.132			26.6	40	
Dibenzo [a,h] anthracene	0.021	0.02	ug/g dry	0.027			24.3	40	
Fluoranthene	0.164	0.02	ug/g dry	0.272			NC	40	
	ND	0.02	ug/g dry	ND			NC	40	
Indeno [1,2,3-cd] pyrene	0.067	0.02	ug/g dry	0.092			31.9	40	
1-Methylnaphthalene	ND	0.02	ug/g dry	ND			NC	40 40	
2-Methylnaphthalene		0.02	ug/g dry ug/g dry	ND ND			NC NC	40 40	
Naphthalene	ND 0.083	0.01	ug/g dry ug/g dry	ND 0.150			NC NC	40 40	
Phenanthrene	0.083 0.151	0.02 0.02	ug/g dry ug/g dry	0.150 0.238			NC NC	40 40	
Pyrene Surrogate: 2-Fluorobiphenyl	0.151 1.52	0.02	ug/g dry <i>ug/g dry</i>	0.230	97.2	50-140	NC	+0	
Surrogate: 2-Fluorobiphenyi Surrogate: Terphenyl-d14	1.52 1.81		ug/g ary ug/g dry		97.2 116	50-140 50-140			
Volatiles			- 3. 5 Gry						
Acetone	ND	0.50	ug/g dry	ND			NC	50	
Benzene	ND	0.02	ug/g dry ug/g dry	ND			NC	50 50	
Bromodichloromethane	ND	0.02	ug/g dry ug/g dry	ND			NC	50 50	
Bromoform	ND	0.05	ug/g dry ug/g dry	ND			NC	50 50	
Bromomethane	ND	0.05	ug/g dry ug/g dry	ND			NC	50 50	
Carbon Tetrachloride	ND	0.05	ug/g dry	ND			NC	50	



### Method Quality Control: Duplicate

Report Date: 02-Sep-2020 Order Date: 27-Aug-2020

Project Description: PE4614

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
			-		%REU	LIIIII			110165
Chloroform	ND	0.05	ug/g dry	ND			NC	50	
Dibromochloromethane	ND	0.05	ug/g dry	ND			NC	50	
Dichlorodifluoromethane	ND	0.05	ug/g dry	ND			NC	50	
1,2-Dichlorobenzene	ND	0.05	ug/g dry	ND			NC	50	
1,3-Dichlorobenzene	ND	0.05	ug/g dry	ND			NC	50	
1,4-Dichlorobenzene	ND	0.05	ug/g dry	ND			NC	50	
1,1-Dichloroethane	ND	0.05	ug/g dry	ND			NC	50	
1,2-Dichloroethane	ND	0.05	ug/g dry	ND			NC	50	
1,1-Dichloroethylene	ND	0.05	ug/g dry	ND			NC	50	
cis-1,2-Dichloroethylene	ND	0.05	ug/g dry	ND			NC	50	
trans-1,2-Dichloroethylene	ND	0.05	ug/g dry	ND			NC	50	
1,2-Dichloropropane	ND	0.05	ug/g dry	ND			NC	50	
cis-1,3-Dichloropropylene	ND	0.05	ug/g dry	ND			NC	50	
trans-1,3-Dichloropropylene	ND	0.05	ug/g dry	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g dry	ND			NC	50	
Ethylene dibromide (dibromoethane, 1,2	ND	0.05	ug/g dry	ND			NC	50	
Hexane	ND	0.05	ug/g dry	ND			NC	50	
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g dry	ND			NC	50	
Methyl Isobutyl Ketone	ND	0.50	ug/g dry	ND			NC	50	
Methyl tert-butyl ether	ND	0.05	ug/g dry	ND			NC	50	
Methylene Chloride	ND	0.05	ug/g dry	ND			NC	50	
Styrene	ND	0.05	ug/g dry	ND			NC	50	
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g dry	ND			NC	50	
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g dry	ND			NC	50	
Tetrachloroethylene	ND	0.05	ug/g dry	ND			NC	50	
Toluene	ND	0.05	ug/g dry	ND			NC	50	
1,1,1-Trichloroethane	ND	0.05	ug/g dry	ND			NC	50	
1,1,2-Trichloroethane	ND	0.05	ug/g dry	ND			NC	50	
Trichloroethylene	ND	0.05	ug/g dry	ND			NC	50	
Trichlorofluoromethane	ND	0.05	ug/g dry	ND			NC	50	
Vinyl chloride	ND	0.02	ug/g dry	ND			NC	50	
m,p-Xylenes	ND	0.05	ug/g dry	ND			NC	50	
o-Xylene	ND	0.05	ug/g dry	ND			NC	50	
Surrogate: 4-Bromofluorobenzene	4.04		ug/g dry		110	50-140			
Surrogate: Dibromofluoromethane	3.84		ug/g dry		105	50-140			
Surrogate: Toluene-d8	4.31		ug/g dry		118	50-140			
Benzene	ND	0.02	ug/g dry	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g dry	ND			NC	50	
Toluene	ND	0.05	ug/g dry	ND			NC	50	
m,p-Xylenes	ND	0.05	ug/g dry	ND			NC	50	
o-Xylene	ND	0.05	ug/g dry	ND			NC	50	
Surrogate: Toluene-d8	4.31	0.00	ug/g dry		118	50-140			



### Method Quality Control: Spike

Report Date: 02-Sep-2020

Order Date: 27-Aug-2020

Project Description: PE4614

Hydrocarbons           F1 PHCs (C6-C10)         202         7         ug/g         ND         101         80-120           F2 PHCs (C10-C16)         116         4         ug/g         ND         94.9         60-140           F3 PHCs (C16-C34)         307         8         ug/g         ND         102         60-140           F4 PHCs (C34-C50)         204         6         ug/g         ND         80-120           Metals         8         ug/g         ND         90.2         70-130           Arsenic         53.3         1.0         ug/g         3.3         100         70-130           Barium         126         1.0         ug/g         ND         95.8         70-130           Boron         46.3         5.0         ug/g         ND         94.1         70-130           Cadmium         47.2         0.5         ug/g         ND         94.1         70-130           Chromium (VI)         0.1         0.2         ug/g         ND         95.0         70-130         QM           Cobalt         53.3         1.0         ug/g         3.8         98.9         70-130         QM           Copper         75.4 </th <th></th>	
F2 PHCs (C10-C16)       116       4       ug'g       ND       94.9       60-140         F3 PHCs (C16-C34)       307       8       ug/g       ND       102       60-140         F4 PHCs (C34-C50)       204       6       ug/g       ND       107       60-140         F4 PHCs (gravimetric)       820       50       ug/g       ND       82.0       80-120         Matis          45.4       1.0       ug/g       ND       90.2       70-130         Arsenic       53.3       1.0       ug/g       3.3       100       70-130         Barium       126       1.0       ug/g       ND       95.8       70-130         Barium       48.2       0.5       ug/g       ND       95.8       70-130         Cadmium       47.2       0.5       ug/g       ND       94.1       70-130         Chromium (VI)       0.1       0.2       ug/g       ND       94.1       70-130         Chromium (VI)       0.3       0.9       ug/g       ND       94.1       70-130         Cobalt       53.3       1.0       ug/g       3.8       98.9       70-130 <td< td=""><td></td></td<>	
F3 PHCs (C16-C34)       307       8       ug/g       ND       102       60-140         F4 PHCs (C34-C50)       204       6       ug/g       ND       107       60-140         F4G PHCs (gravimetric)       820       50       ug/g       ND       82.0       80-120         Metals                 Antimony       45.4       1.0       ug/g       3.3       100       70-130       70-130         Barium       126       1.0       ug/g       72.0       107       70-130       70-130         Boron       46.3       5.0       ug/g       ND       95.8       70-130       70-130         Cadmium       47.2       0.5       ug/g       ND       95.4       70-130       70-130         Chromium (VI)       0.1       0.2       ug/g       ND       65.0       70-130       QN         Cobalt       53.3       1.0       ug/g       3.8       98.9       70-130       QN         Cobalt       53.3       1.0       ug/g       9.6       89.7       70-130       QN         Mercury       3.03       0.1       ug/g	
F4 PHCs (C34-C50)       204       6       ug/g       ND       107       60-140         F4G PHCs (gravimetric)       820       50       ug/g       ND       82.0       80-120         Metals         1.0       ug/g       ND       90.2       70-130         Arsenic       53.3       1.0       ug/g       3.3       100       70-130         Barium       126       1.0       ug/g       72.0       107       70-130         Beryllium       48.2       0.5       ug/g       ND       95.8       70-130         Boron       46.3       5.0       ug/g       ND       95.4       70-130         Cadmium       71       0.2       ug/g       ND       95.4       70-130         Coronium (V1)       0.1       0.2       ug/g       ND       95.0       70-130         Coper       75.4       5.0       ug/g       10.3       70-130       QM         Mercury       3.03       0.1       ug/g       9.6       89.7       70-130         Lead       135       1.0       ug/g       1.35       112       70-130         Molydenum       9.0       1.0	
F4G PHCs (gravimetric)         820         50         ug/g         ND         82.0         80-120           Metals         Antimony         45.4         1.0         ug/g         ND         90.2         70-130           Arsenic         53.3         1.0         ug/g         3.3         100         70-130           Barium         126         1.0         ug/g         72.0         107         70-130           Beryllium         48.2         0.5         ug/g         ND         95.8         70-130           Boron         46.3         5.0         ug/g         ND         85.4         70-130           Chromium (VI)         0.1         0.2         ug/g         ND         65.0         70-130         QM           Cobalt         5.3.3         1.0         ug/g         ND         65.0         70-130         QM           Cobalt         5.3.3         1.0         ug/g         3.8         98.9         70-130           Cobalt         5.3.3         1.0         ug/g         3.8         98.9         70-130           Lead         1.35         1.0         ug/g         1.35         1.12         70-130           Molybdenum	
Metals         Antimony         45.4         1.0         ug/g         ND         90.2         70-130           Arsenic         53.3         1.0         ug/g         3.3         100         70-130           Barium         126         1.0         ug/g         72.0         107         70-130           Beryllium         48.2         0.5         ug/g         ND         95.8         70-130           Boron         46.3         5.0         ug/g         ND         85.4         70-130           Cadmium         47.2         0.5         ug/g         ND         94.1         70-130           Chromium (VI)         0.1         0.2         ug/g         ND         65.0         70-130           Cobalt         53.3         1.0         ug/g         3.8         98.9         70-130           Copper         75.4         5.0         ug/g         1.35         100         70-130           Metaly         135         1.0         ug/g         1.35         112         70-130           Lead         135         1.0         ug/g         1.35         112         70-130           Molybdenum         49.0         1.0         ug	
Antimony45.41.0ug/gND90.270-130Arsenic53.31.0ug/g3.310070-130Barium1261.0ug/g72.010770-130Beryllium48.20.5ug/gND95.870-130Boron46.35.0ug/gND94.170-130Cadmium47.20.5ug/gND94.170-130Chromium (VI)0.10.2ug/gND65.070-130Cobalt53.31.0ug/g3.898.970-130Cobalt53.31.0ug/g3.898.970-130Lead1351.0ug/g90.689.770-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Arsenic53.31.0ug/g3.310070-130Barium1261.0ug/g72.010770-130Beryllium48.20.5ug/gND95.870-130Boron46.35.0ug/gND85.470-130Cadmium47.20.5ug/gND94.170-130Chromium (VI)0.10.2ug/gND65.070-130Cobalt53.31.0ug/g3.898.970-130Cobalt53.31.0ug/g3.898.970-130Copper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Arsenic53.31.0ug/g3.310070-130Barium1261.0ug/g72.010770-130Beryllium48.20.5ug/gND95.870-130Boron46.35.0ug/gND85.470-130Cadmium47.20.5ug/gND94.170-130Chromium (VI)0.10.2ug/gND65.070-130QNChromium63.85.0ug/g12.210370-130QNCobalt53.31.0ug/g3.898.970-130NDCopper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/gND96.370-130Selenium44.30.3ug/gND92.270-130	
Beryllium48.20.5ug/gND95.870-130Boron46.35.0ug/gND85.470-130Cadmium47.20.5ug/gND94.170-130Chromium (VI)0.10.2ug/gND65.070-130QNChromium63.85.0ug/g12.210370-130QNCobalt53.31.0ug/g3.898.970-130Copper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Mercury3.030.1ug/gND96.370-130Nolybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/gND92.270-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Boron46.35.0ug/gND85.470-130Cadmium47.20.5ug/gND94.170-130Chromium (VI)0.10.2ug/gND65.070-130QNChromium63.85.0ug/g12.210370-130QNCobalt53.31.0ug/g3.898.970-130Copper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Mercury3.030.1ug/gND96.370-130Nickel63.65.0ug/gND96.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Cadmium47.20.5ug/gND94.170-130Chromium (VI)0.10.2ug/gND65.070-130QNChromium63.85.0ug/g12.210370-130Cobalt53.31.0ug/g3.898.970-130Copper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Mercury3.030.1ug/g1.3511270-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/gND99.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Chromium (VI)         0.1         0.2         ug/g         ND         65.0         70-130         QM           Chromium         63.8         5.0         ug/g         12.2         103         70-130         QM           Cobalt         53.3         1.0         ug/g         3.8         98.9         70-130         Copper         Copper         75.4         5.0         ug/g         25.3         100         70-130         Copper         Lead         135         1.0         ug/g         90.6         89.7         70-130         Copper         Molybdenum         49.0         1.0         ug/g         ND         96.3         70-130         ND         ND         ND         10         ND         96.3         70-130         ND         ND         ND         10         ND         10         ND         10         ND         10         ND         10         10         ND         10         10         ND         10<	
Chromium63.85.0ug/g12.210370-130Cobalt53.31.0ug/g3.898.970-130Copper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Mercury3.030.1ug/g1.3511270-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Cobalt53.31.0ug/g3.898.970-130Copper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Mercury3.030.1ug/g1.3511270-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	<i>I</i> -05
Copper75.45.0ug/g25.310070-130Lead1351.0ug/g90.689.770-130Mercury3.030.1ug/g1.3511270-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Lead1351.0ug/g90.689.770-130Mercury3.030.1ug/g1.3511270-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Mercury3.030.1ug/g1.3511270-130Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Molybdenum49.01.0ug/gND96.370-130Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Nickel63.65.0ug/g13.999.370-130Selenium46.41.0ug/gND92.270-130Silver44.30.3ug/g0.487.770-130	
Selenium         46.4         1.0         ug/g         ND         92.2         70-130           Silver         44.3         0.3         ug/g         0.4         87.7         70-130	
Silver 44.3 0.3 ug/g 0.4 87.7 70-130	
Thallium 46.5 1.0 μα/α ND 92.8 70-130	
Uranium 48.4 1.0 ug/g ND 95.0 70-130	
Vanadium 63.8 10.0 ug/g 12.2 103 70-130	
Zinc 121 20.0 ug/g 69.9 102 70-130	
Semi-Volatiles	
Acenaphthene 0.210 0.02 ug/g ND 107 50-140	
Acenaphthylene 0.233 0.02 ug/g 0.051 93.0 50-140	
Anthracene 0.258 0.02 ug/g 0.053 105 50-140	
Benzo [a] anthracene 0.354 0.02 ug/g 0.130 115 50-140	
Benzo [a] pyrene 0.364 0.02 ug/g 0.144 113 50-140	
Benzo [b] fluoranthene 0.489 0.02 ug/g 0.187 154 50-140 QM	Л-06
Benzo [g,h,i] perylene 0.281 0.02 ug/g 0.099 93.2 50-140	
Benzo [k] fluoranthene 0.351 0.02 ug/g 0.098 129 50-140	
Chrysene 0.405 0.02 ug/g 0.132 140 50-140	
Dibenzo [a,h] anthracene 0.202 0.02 ug/g 0.027 89.5 50-140	
Fluoranthene 0.601 0.02 ug/g 0.272 169 50-140 QM	Л-06
Fluorene 0.199 0.02 ug/g ND 102 50-140	
Indeno [1,2,3-cd] pyrene 0.290 0.02 ug/g 0.092 101 50-140	
1-Methylnaphthalene 0.197 0.02 ug/g ND 101 50-140	
2-Methylnaphthalene 0.218 0.02 ug/g ND 112 50-140	
Naphthalene 0.208 0.01 ug/g ND 106 50-140	
	<i>I</i> -06
	<i>I</i> -06
Surrogate: 2-Fluorobiphenyl         1.60         ug/g         103         50-140	
Surrogate: Terphenyl-d14 1.91 ug/g 122 50-140	
Volatiles	



### Order #: 2035556

Report Date: 02-Sep-2020

Order Date: 27-Aug-2020

Project Description: PE4614

### Method Quality Control: Spike

Action         12.8         0.50         ugig         ND         12.8         0.510           Bernendichkomethane         3.49         0.05         ugig         ND         84.8         60.130           Bromodichkomethane         3.79         0.05         ugig         ND         84.8         60.130           Bromodichkomethane         3.79         0.05         ugig         ND         78.8         60.130           Cathon Fibrachoride         3.19         0.05         ugig         ND         84.7         60.130           Cheroform         3.39         0.05         ugig         ND         86.2         60.130           Dicharodifucromethane         3.65         0.05         ugig         ND         86.7         60.130           1.3.Dicharodifucromethane         3.48         0.05         ugig         ND         86.6         60.130           1.3.Dicharodifucromethane         3.47         0.06         ugig         ND         86.8         60.130           1.3.Dicharodifucromethane         3.47         0.06         ugig         ND         75.8         60.130           1.3.Dicharodifucromethane         3.47         0.05         ugig         ND         75.8 <t< th=""><th>Analyte</th><th>Result</th><th>Reporting Limit</th><th>Units</th><th>Source Result</th><th>%REC</th><th>%REC Limit</th><th>RPD</th><th>RPD Limit</th><th>Notes</th></t<>	Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Bromoter Bromoter Bromoter3.390.05ug/gND8.4860-130Bromoter Bromoter Bromoter3.790.05ug/gND8.1150-140Carbon Tetrachloride3.190.05ug/gND8.7160-130Chinobertzene3.800.05ug/gND8.7260-130Dienkoordikorernethane3.800.05ug/gND8.7260-130Dienkoordikorernethane3.850.05ug/gND8.7460-1301.3-Dichtorbertzene3.840.05ug/gND8.7460-1301.3-Dichtorbertzene3.840.05ug/gND8.7460-1301.3-Dichtorbertzene3.740.05ug/gND8.7460-1301.3-Dichtorbertzene3.740.05ug/gND7.860-1301.3-Dichtorbertylene3.740.05ug/gND7.860-1301.3-Dichtorbertylene3.740.05ug/gND7.860-1301.3-Dichtorporphene3.740.05ug/gND7.860-1301.3-Dichtorporphene3.740.05ug/gND7.860-1301.3-Dichtorporphene3.740.05ug/gND7.860-1301.3-Dichtorporphene3.740.05ug/gND7.860-1301.3-Dichtorporphene3.740.05ug/gND6.960-1301.3-Dichtorporphene3.740.05	Acetone	12.8	0.50	ug/g	ND	128	50-140			
Bromorthame3.790.05ug/gND9.4360-130Bromorthame3.190.05ug/gND9.3150-140Chicorbentenchortoie4.130.05ug/gND10360-130Chicorbentenchortoine4.330.05ug/gND10260-130Dicordontrionethame4.880.05ug/gND8.4760-130Licolohortoinomethame3.840.05ug/gND8.4760-1301.2-Dichordoffuctorethame3.480.05ug/gND8.4760-1301.3-Dichortoinomethame3.470.05ug/gND8.4860-1301.3-Dichordoffuctorethy3.470.05ug/gND8.4860-1301.4-Dichortoinotethy3.470.05ug/gND7.560-1301.3-Dichordoffuctorethy3.470.05ug/gND7.560-1301.3-Dichortoinotethy3.470.05ug/gND7.560-1301.3-Dichortoinotethy3.470.05ug/gND7.560-1301.3-Dichortoinoty3.470.05ug/gND7.560-1301.3-Dichortoinoty3.470.05ug/gND6.1360-1301.3-Dichortoinoty3.670.05ug/gND7.560-1301.3-Dichortoinoty6.90.05ug/gND6.1360-1301.3-Dichortoinoty6.90.05ug/gND6.13 <td>Benzene</td> <td>3.41</td> <td>0.02</td> <td>ug/g</td> <td>ND</td> <td>85.2</td> <td>60-130</td> <td></td> <td></td> <td></td>	Benzene	3.41	0.02	ug/g	ND	85.2	60-130			
Brannentiane1.730.65ug'qND9.1460-140Carbon Teixachioride1.190.65ug'qND70.860-130Chiorabanzane3.190.55ug'qND10.260-130Dichorabitration3.580.55ug'qND80.460-130Dichorabitration3.580.55ug'qND80.460-1301.3-Dichorabenzane3.580.55ug'qND80.460-1301.3-Dichorabenzane3.560.55ug'qND81.760-1301.3-Dichorabenzane3.670.55ug'qND81.860-1301.1-Dichorabityene3.670.55ug'qND81.860-1301.1-Dichorabityene3.670.55ug'qND71.560-1301.1-Dichorabityene3.670.55ug'qND71.860-1301.2-Dichorabityene3.270.55ug'qND71.860-1301.2-Dichorabityene3.670.55ug'qND71.860-1301.2-Dichorabityene3.670.55ug'qND71.860-1301.2-Dichorabityene3.620.55ug'qND71.660-1301.2-Dichorabityene3.620.55ug'qND71.660-1301.2-Dichorabityene3.620.55ug'qND71.660-1301.2-Dichorabityene3.620.55ug'qND71.660-130 <td>Bromodichloromethane</td> <td>3.39</td> <td>0.05</td> <td>ug/g</td> <td>ND</td> <td>84.8</td> <td>60-130</td> <td></td> <td></td> <td></td>	Bromodichloromethane	3.39	0.05	ug/g	ND	84.8	60-130			
Carbon Fetrachionels3.190.05ug/gND7.8860-130Chiorobernzene4.130.05ug/gND10.4760-130Dinconchioromelhane3.890.05ug/gND06.250-1401.2-Dichorobernzene3.850.05ug/gND87.460-1301.2-Dichorobernzene3.860.05ug/gND88.760-1301.3-Dichorobernzene3.870.05ug/gND88.760-1301.4-Dichorobernzene3.670.05ug/gND88.860-1301.4-Dichorobernzene3.670.05ug/gND81.860-1301.4-Dichorobernyene3.660.05ug/gND75.560-1301.4-Dichorobernyene3.720.05ug/gND75.560-1301.2-Dichorobernyene3.720.05ug/gND75.560-1301.2-Dichoropropane3.120.05ug/gND75.560-1301.2-Dichoropropane2.740.05ug/gND76.060-1301.2-Dichoropropane2.740.05ug/gND76.060-1301.2-Dichoropropane2.740.05ug/gND76.060-1301.2-Dichoropropane2.740.05ug/gND76.060-1301.2-Dichoropropane2.740.05ug/gND76.060-1301.2-Dichoropropane3.720.05ug/gND60-130<	Bromoform	3.79	0.05	ug/g	ND	94.8	60-130			
Chiotocharma1.30.65ug'gND10.360-13Chiotocharma3.890.65ug'gND10.260-130Dichiotochironenthane3.860.05ug'gND0.6250-1401.2-Dichiroberbaren3.580.05ug'gND8.760-1301.3-Dichiroberbaren3.560.05ug'gND8.760-1301.4-Dichiroberbaren3.770.65ug'gND8.860-1301.1-Dichiroberbaren3.660.05ug'gND8.860-1301.1-Dichiroberbyten3.660.05ug'gND7.560-1301.1-Dichiroberbyten3.670.05ug'gND7.860-1301.2-Dichiroberbyten3.270.05ug'gND7.5560-1301.2-Dichiroberbyten3.220.05ug'gND7.5560-1301.2-Dichiroberbyten3.240.05ug'gND7.5560-1301.2-Dichiroberbyten2.640.05ug'gND61.761-1301.2-Dichiroberbyten2.640.05ug'gND61.761-1301.2-Dichiroberbyten6.740.55ug'gND61.761-1301.2-Dichiroberbyten6.840.55ug'gND61.761-1301.2-Dichiroberbyten7.640.55ug'gND61.761-1301.2-Dichiroberbyten7.640.55ug'gND61.761-1	Bromomethane	3.73	0.05	ug/g	ND	93.1	50-140			
Chieroform3.990.05ug'gND8.760-130Dibromochioromethane3.850.05ug'gND80.250-1401.2-Dichlorodhuzomethane3.860.05ug'gND87.460-1301.2-Dichlorodhuzomethane3.480.05ug'gND87.460-1301.3-Dichlorodhuzomethane3.480.05ug'gND87.460-1301.4-Dichlorodhuzomethane3.470.05ug'gND88.660-1301.1-Dichlorodhyden3.470.05ug'gND7.560-1301.1-Dichlorodhyden3.120.05ug'gND7.560-1301.1-Dichlorodhyden3.120.05ug'gND7.560-130is1.2.Dichlorophyden3.120.05ug'gND7.560-130is1.2.Dichlorophyden3.120.05ug'gND7.560-130is1.2.Dichlorophyden3.120.05ug'gND7.560-130is1.3.Dichlorophyden3.120.05ug'gND7.560-130istans.1.3.Dichlorophyden4.080.05ug'gND7.560-130istans.1.3.Dichlorophyden4.080.05ug'gND7.560-130istans.1.3.Dichlorophyden4.080.05ug'gND7.660-130istans.1.3.Dichlorophyden7.560-1301.160-130istans.1.3.Dichlorophyden7.660-1301.1 <td< td=""><td>Carbon Tetrachloride</td><td>3.19</td><td>0.05</td><td>ug/g</td><td>ND</td><td>79.8</td><td>60-130</td><td></td><td></td><td></td></td<>	Carbon Tetrachloride	3.19	0.05	ug/g	ND	79.8	60-130			
Dibriomochlanomethane4.980.05ug'gND10260-130Dichlacodifuciomethane3.580.05ug'gND8.6460-1301.3-Dichlorobenzene3.580.05ug'gND8.760-1301.4-Dichlorobenzene3.550.05ug'gND8.6860-1301.1-Dichloroethane3.770.05ug'gND8.6860-1301.2-Dichloroethane3.660.05ug'gND7.6560-1301.2-Dichloroethane3.660.05ug'gND7.6560-1301.2-Dichloroethyene3.660.05ug'gND7.5560-1301.2-Dichloroethyene3.670.05ug'gND7.5560-1301.2-Dichloropropyene3.270.05ug'gND7.5560-1301.2-Dichloropropyene3.680.05ug'gND7.5560-1301.2-Dichloropropyene2.740.05ug'gND7.560-1301.2-Dichloropropyene2.740.05ug'gND7.560-1301.2-Dichloropropyene2.740.05ug'gND7.560-1301.2-Dichloropropyene2.740.05ug'gND7.560-1301.2-Dichloropropyene2.740.05ug'gND7.560-1301.2-Dichloropropyene2.740.05ug'gND7.560-1301.2-Dichloropropyene2.740.05ug'gND <td>Chlorobenzene</td> <td>4.13</td> <td>0.05</td> <td>ug/g</td> <td>ND</td> <td>103</td> <td>60-130</td> <td></td> <td></td> <td></td>	Chlorobenzene	4.13	0.05	ug/g	ND	103	60-130			
Dichlaradihuacomethane3.850.65ug'gND96.250-1401,2-Dichlorobenzane3.480.65ug'gND87.160-1301,4-Dichlorobenzane3.470.65ug'gND86.760-1301,4-Dichlorobenzane3.470.65ug'gND86.760-1301,2-Dichlorobenzane3.470.65ug'gND81.860-1301,1-Dichlorobentylene3.600.05ug'gND79.160-1301,2-Dichlorobylene3.160.55ug'gND78.060-1301,2-Dichlorophylene3.120.55ug'gND78.060-1301,2-Dichlorophylene3.220.55ug'gND75.560-1301,2-Dichlorophylene3.220.55ug'gND10.260-1301,3-Dichlorophylene3.220.55ug'gND10.260-130Einylene dibornide (dibornoethane, 1,2-40.10.55ug'gND10.260-130Einylene dibornide (dibornoethane, 1,2-40.10.55ug'gND10.260-130Hexane6.970.55ug'gND10.260-130Heyne Chloride7.410.55ug'gND10.260-130Heyne Chloride6.970.55ug'gND10.260-130Hylene Chloride6.970.55ug'gND10.260-130Hylene Chloride6.970.55ug'g <td>Chloroform</td> <td>3.39</td> <td>0.05</td> <td>ug/g</td> <td>ND</td> <td>84.7</td> <td>60-130</td> <td></td> <td></td> <td></td>	Chloroform	3.39	0.05	ug/g	ND	84.7	60-130			
1.2-Dichlorobenzene3.680.05uggND89.460-1301.3-Dichlorobenzene3.480.05uggND88.760-1301.1-Dichlorobenzene3.470.05uggND81.860-1301.1-Dichlorobenzene3.470.05uggND81.860-1301.1-Dichlorobenzene3.680.05uggND81.860-1301.1-Dichlorobethree3.610.05uggND81.860-1301.1-Dichlorobethree3.620.05uggND81.860-1301.2-Dichlorobethree3.620.05uggND81.860-1301.2-Dichlorobethree3.620.05uggND81.860-1301.2-Dichloropophene3.620.05uggND61.560-1301.3-Dichloropophene2.740.05uggND60.560-130Entylene dichromoethane, 1.24.010.55uggND60.660-130Heinyl Ekone (Abulanne)6.870.50uggND81.860-130Heinyl Ekone (Abulanne)3.670.50uggND82.060-1301.1.1.2-Teichlorobethree3.670.50uggND81.960-1301.1.1.2-Teichlorobethree3.670.50uggND81.960-1301.1.1.2-Teichlorobethree3.600.50uggND81.960-1301.1.1.2-Teichlorobethree3.600.50u	Dibromochloromethane	4.08	0.05		ND	102	60-130			
1.2-Dichlorobenzene3.880.65ug/gND80.460-1301.3-Dichlorobenzene3.650.65ug/gND88.760-1301.1-Dichlorobenzene3.470.65ug/gND81.860-1301.1-Dichlorobenzene3.270.65ug/gND81.860-1301.2-Dichloroethylene3.270.65ug/gND81.860-1301.1-Dichloroethylene3.270.65ug/gND75.560-1301.2-Dichloroethylene3.270.65ug/gND75.560-1301.2-Dichloroethylene3.200.65ug/gND75.560-1301.2-Dichloroethylene3.200.65ug/gND75.560-1301.2-Dichloroethylene3.200.65ug/gND75.560-1301.3-Dichloropropane3.200.65ug/gND60.560-130Ethylene Chromoethane, 1.24.010.55ug/gND60.560-130Heithyl Ethylektone (2-Butanone)6.660.50ug/gND75.560-130Heithyl Ethylektone3.670.50ug/gND75.560-130Heithyl Ethylektone3.670.50ug/gND82.060-1301.1.1.2-Tichloroethylene3.670.50ug/gND61.61.1.2.2-Tetrachloroethane3.670.55ug/gND61.91.1.1.2-Tichloroethylene3.670.50ug/g	Dichlorodifluoromethane	3.85	0.05	ug/g	ND	96.2	50-140			
1,4-Dichlorobenzene3,550,05ug/gND88,760-1301,1-Dichloroethane3,470,05ug/gND86,660-1301,1-Dichloroethylene3,060,05ug/gND78,160-130(is-1,2-Dichloroethylene3,160,05ug/gND78,160-1301,2-Dichloroethylene3,120,05ug/gND78,060-1301,2-Dichloroethylene3,120,05ug/gND78,060-1301,3-Dichloroproprigne3,120,05ug/gND78,060-130trans-1,3-Dichloroproprigne2,740,05ug/gND71,560-130Enlylene dibromide (dibromoethane, 1,24,080,05ug/gND71,560-130Hexane2,860,05ug/gND71,560-130Hexane2,860,05ug/gND71,560-130Hethyle foltor7,810,05ug/gND71,660-130Hethyle foltor7,810,05ug/gND71,660-130Hethyle foltor3,280,05ug/gND71,760-130Hethyle foltoroethane4,05ug/gND10160-1301,1,2,2-Trachoroethane4,05ug/gND10160-1301,1,1,2-Trichoroethane3,160,05ug/gND10160-1301,1,1,2-Trichoroethane3,110,05ug/gND10160-130 </td <td>1,2-Dichlorobenzene</td> <td>3.58</td> <td>0.05</td> <td></td> <td>ND</td> <td>89.4</td> <td>60-130</td> <td></td> <td></td> <td></td>	1,2-Dichlorobenzene	3.58	0.05		ND	89.4	60-130			
1,4-Dichlorobenzene3,550,05ug/gND88,760-1301,1-Dichloroethane3,470,05ug/gND86,660-1301,1-Dichloroethylene3,060,05ug/gND78,160-130(is-1,2-Dichloroethylene3,160,05ug/gND78,160-1301,2-Dichloroethylene3,120,05ug/gND78,060-1301,2-Dichloroethylene3,120,05ug/gND78,060-1301,3-Dichloroproprigne3,120,05ug/gND78,060-130trans-1,3-Dichloroproprigne2,740,05ug/gND71,560-130Enlylene dibromide (dibromoethane, 1,24,080,05ug/gND71,560-130Hexane2,860,05ug/gND71,560-130Hexane2,860,05ug/gND71,560-130Hethyle foltor7,810,05ug/gND71,660-130Hethyle foltor7,810,05ug/gND71,660-130Hethyle foltor3,280,05ug/gND71,760-130Hethyle foltoroethane4,05ug/gND10160-1301,1,2,2-Trachoroethane4,05ug/gND10160-1301,1,1,2-Trichoroethane3,160,05ug/gND10160-1301,1,1,2-Trichoroethane3,110,05ug/gND10160-130 </td <td>1,3-Dichlorobenzene</td> <td>3.48</td> <td>0.05</td> <td></td> <td></td> <td>87.1</td> <td>60-130</td> <td></td> <td></td> <td></td>	1,3-Dichlorobenzene	3.48	0.05			87.1	60-130			
1.1-Dichloroethane3.470.05ug/gND86.60-1301.2-Dichloroethylene3.670.05ug/gND76.160-130cis-1.2-Dichloroethylene3.160.05ug/gND78.160-1301.1-Dichloroethylene3.120.05ug/gND78.160-1301.2-Dichloroethylene3.120.05ug/gND78.060-1301.2-Dichloropropylene3.120.05ug/gND75.560-1301.1-Dichloropropylene3.020.05ug/gND75.560-130Ethylenzene4.080.05ug/gND71.560-130Ethylenzene2.680.05ug/gND71.560-130Hexane2.680.05ug/gND71.560-130Methyl Ethyl Kotone (2-Butanone)6.860.50ug/gND60.650-140Methyl Ischulyl Kotone3.280.05ug/gND61.360-130Styrene3.280.05ug/gND61.360-1301.1,1.2-Tetrachloroethane4.740.05ug/gND61.360-1301.1,1.2-Tetrachloroethane3.140.05ug/gND61.360-1301.1,1.2-Tetrachloroethane4.05ug/gND61.360-1301.1,1.2-Tetrachloroethane4.05ug/gND10.160-1301.1,1.2-Tetrachloroethane3.110.05ug/gND10.2 <td>1,4-Dichlorobenzene</td> <td>3.55</td> <td>0.05</td> <td></td> <td></td> <td>88.7</td> <td>60-130</td> <td></td> <td></td> <td></td>	1,4-Dichlorobenzene	3.55	0.05			88.7	60-130			
1.2-Dichloroethane3.270.05ug/gND81.860-1301.1-Dichloroethylene3.060.05ug/gND75.560-130trans-1.2-Dichloroethylene3.270.05ug/gND81.860-1301.2-Dichloroptylene3.120.05ug/gND75.060-130trans-1.3-Dichloroptylene3.020.05ug/gND65.560-130Ethylenethane2.740.05ug/gND60.560-130Ethylenethane4.080.05ug/gND60.560-130Hexane2.860.05ug/gND60.560-130Hexane2.860.05ug/gND60.560-130Hethyl Ethylenet/2-Butanene6.970.05ug/gND60.750-140Methyl Ethylenet/2-Butanene6.970.05ug/gND61.750-140Methyl Etholoride7.810.50ug/gND61.760-130Syrene3.670.55ug/gND61.760-130Syrene3.660.50ug/gND10.160-130Tichloroethane1.420.55ug/gND61.30Tichloroethane3.600.55ug/gND61.30Tichloroethane3.600.55ug/gND61.30Tichloroethane3.610.55ug/gND61.30Tichloroethane3.610.55ug/gND										
1.1-Dichloroethylene3.060.05ug'gND76.560-130cis-1.2-Dichloroethylene3.160.05ug/gND78.060-1301.2-Dichloroethylene3.120.05ug/gND78.060-1301.2-Dichloropropane3.120.05ug/gND78.560-130tisn-1.3-Dichloropropane3.020.05ug/gND10260-130Ethylenzene4.080.05ug/gND10260-130Ethylenzene4.080.05ug/gND60.650-140Methyl Ethyl Ketone (2-Butanone)6.960.50ug/gND78.150-140Methyl Isobutyl Ketone (2-Butanone)6.960.50ug/gND78.150-140Methyl Isobutyl Ketone (2-Butanone)6.960.50ug/gND78.160-130Methyl Isobutyl Ketone (2-Butanone)6.960.50ug/gND78.160-130Methyl Isobutyl Ketone7.810.50ug/gND91.760-1301.1,1.2-Tetrachoroethane3.670.55ug/gND19.160-1301.1,1.2-Tetrachoroethane3.680.55ug/gND19.260-1301.1,1.2-Tetrachoroethane3.610.55ug/gND19.360-1301.1,1.2-Tetrachoroethane3.110.55ug/gND13.360-1301.1,1.2-Tetrachoroethane3.210.55ug/gND13.360-130										
cis-1.2-Dichloroethylene3.160.05u.g'gND7.9.160.130trans.1.2-Dichloroethylene3.270.05u.g'gND7.8.060.130cis-1.3-Dichloropropane3.020.05u.g'gND7.5.860.130cis-1.3-Dichloropropylene2.740.05u.g'gND6.5.560.130Ethylbenzene0.05u.g'gND10060.130Ethylbenzeh4.010.05u.g'gND7.1.560.130Ethylbenzeh6.660.05u.g'gND6.7.660.130Hexane0.660.50u.g'gND6.7.660.130Methyl Ethylketone (2-Butanone)6.670.50u.g'gND6.7.660.130Methyl Ethylether7.810.50u.g'gND6.7.660.130Methyl Ethylether6.670.55u.g'gND61.760.130Syrene3.670.55u.g'gND61.760.130Syrene3.670.55u.g'gND61.760.1301.1,2.2 Fetrachloroethylene3.660.55u.g'gND61.301.1,2.2 Fetrachloroethane3.610.55u.g'gND61.301.1,1.2 Fichloroethane3.640.55u.g'gND63.61.1,2.7 Fichloroethane3.640.55u.g'gND63.71.1,2.7 Fichloroethane3.640.55u.g'gND63.101.1,2.7 Fichloroethane<										
trans-1.2-Dichloroethylene3.270.05ug'gND81.860-1301.2-Dichloropropane3.120.05ug/gND78.060-130trans-1.3-Dichloropropylene2.740.05ug/gND68.560-130Ethylene dilormide (dibromoethane, 1.24.080.05ug/gND10260-130Hexane2.860.05ug/gND71.560-130Methyl Ethyl Ketone (2-Butanone)6.960.50ug/gND69.650-140Methyl Ethyl Ketone (2-Butanone)6.970.05ug/gND69.750-140Methyl Ich-Lulyl ether6.970.05ug/gND69.750-140Methyl Ichoroethane3.280.05ug/gND69.750-1401.1,1.2-Teitrachloroethane4.740.05ug/gND10160-1301.1,1.2-Teitrachloroethane3.130.05ug/gND78.360-1301.1,1.2-Teitrachloroethane3.140.05ug/gND78.360-1301.1,1.2-Teitrachloroethane3.130.05ug/gND78.360-1301.1,1.2-Teitrachloroethane3.140.05ug/gND78.360-1301.1,1.2-Teitrachloroethane3.140.05ug/gND83.150-1401.1,1.2-Teitrachloroethane3.140.05ug/gND77.760-1301.1,1.2-Teitrachloroethane3.140.05ug/gND83.1 <td></td>										
1.2-Dichloropropane3.120.05ug/gND78.060-130cis-1.3-Dichloropropylene3.020.05ug/gND75.560-130Ethylenzene4.080.05ug/gND10260-130Ethylenzene4.080.05ug/gND71.560-130Hexane2.860.05ug/gND71.560-130Methyl Ethyl Ketone (2-Butanone)6.960.65ug/gND78.150-140Methyl Isobutyl Ketone7.810.50ug/gND78.150-140Methyl Isobutyl Ketone7.810.50ug/gND82.060-130Methyl Isobutyl Ketone3.670.05ug/gND82.060-1301,1,2.2-Tetrachloroethane4.050.05ug/gND91.760-1301,1,1.2-Tetrachloroethane4.050.05ug/gND10160-1301,1,1.2-Tetrachloroethane3.660.05ug/gND10160-1301,1,1.2-Tetrachloroethane3.680.05ug/gND10160-1301,1,1.2-Tetrachloroethane3.640.05ug/gND17.560-1301,1,1.2-Tetrachloroethane3.610.05ug/gND17.660-1301,1,1.2-Tetrachloroethane3.620.05ug/gND17.660-1301,1,1.2-Tetrachloroethane3.610.05ug/gND17.760-1301,1,1.2-Tetrachloroethane <t< td=""><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></t<>										
cis-1.3-Dichloropropylene3.020.05u.g/gND75.560-130trans-1.3-Dichloropropylene2.740.05u.g/gND10260-130Ethylene dibromide (dibromoethane, 1.2.4.010.05u.g/gND10060-130Hexane2.860.05u.g/gND71.560-130Methyl Ketone (2-Butanone)6.960.50u.g/gND78.150-140Methyl Stohuly Ketone7.810.50u.g/gND82.060-130Methyl Letr-buly ether6.970.05u.g/gND82.060-130Styrene3.670.05u.g/gND91.760-1301,1,2.2-Tetrachloroethane4.050.05u.g/gND91.760-1301,1,2.2-Tetrachloroethane4.050.05u.g/gND10160-1301,1,2.2-Tetrachloroethane4.050.05u.g/gND10160-1301,1,2.2-Tetrachloroethane4.740.05u.g/gND10160-1301,1,1.2-Trichloroethane3.130.05u.g/gND82.060-1301,1,1.2-Trichloroethane3.140.05u.g/gND82.060-1301,1,1.2-Trichloroethane3.140.05u.g/gND82.060-1301,1,1.2-Trichloroethane3.140.05u.g/gND83.160-1301,1,2-Trichloroethane4.840.05u.g/gND83.160-130	-									
trans-1,3-Dichoropropylene       2,74       0.05       ug/g       ND       68.5       60-130         Ethylene dibromde (dibromethane, 1,2       4.01       0.05       ug/g       ND       102       60-130         Hexane       2.86       0.05       ug/g       ND       71.5       60-130         Methyl Ethyl Ketone (2-Butanone)       6.96       0.50       ug/g       ND       69.6       50-140         Methyl Isbutyl Ketone       7.81       0.50       ug/g       ND       69.7       50-140         Methyl Isbutyl Ketone       3.28       0.05       ug/g       ND       82.0       60-130         Styrene       3.67       0.05       ug/g       ND       81.7       60-130         1,1,2.2-Tetrachloroethane       4.05       0.05       ug/g       ND       81.0       60-130         1,1,2.2-Tetrachloroethane       3.66       0.05       ug/g       ND       81.0       60-130         1,1,2.2-Tetrachloroethane       3.60       0.05       ug/g       ND       81.0       60-130         1,1,2.2-Tetrachloroethane       3.13       0.05       ug/g       ND       78.3       60-130         1,1,2.2-Ticthoroethane       3.13       0.05										
Ethylbenzene       4.08       0.05       ug'g       ND       102       60-130         Ethylbene dibromide (dibromoethane, 1,2)       4.01       0.05       ug/g       ND       100       60-130         Hexane       2.86       0.50       ug/g       ND       69.6       50-140         Methyl Ethyl Ketone (2-Butanone)       6.96       0.50       ug/g       ND       69.7       50-140         Methyl Isobutyl Ketone       7.81       0.50       ug/g       ND       69.7       50-140         Methyl Isobutyl Ketone       7.81       0.50       ug/g       ND       69.7       50-140         Methyl Isobutyl Ketone       7.81       0.50       ug/g       ND       61.7       60-130         Styrene       3.67       0.05       ug/g       ND       101       60-130         1,1,2.7Etrachloroethane       4.74       0.05       ug/g       ND       101       60-130         1,1,2.7Etrachloroethane       3.67       0.05       ug/g       ND       107       60-130         1,1,1.7Trichloroethane       3.13       0.05       ug/g       ND       78.3       60-130         1,1,1.7Trichloroethane       3.41       0.05       ug/g </td <td></td>										
Ethylene dibromide (dibromoethane, 1,2.         4.01         0.05         ug/g         ND         100         60-130           Hexane         2.86         0.05         ug/g         ND         71.5         60-130           Methyl Ethyl Ketone (2-Butanone)         6.96         0.50         ug/g         ND         69.6         50-140           Methyl terbulyl ether         6.97         0.05         ug/g         ND         69.7         50-140           Methyl terbulyl ether         6.97         0.05         ug/g         ND         61.7         60-130           Styrene         3.67         0.05         ug/g         ND         61.7         60-130           1,1,2.2-Tetrachloroethane         4.05         0.05         ug/g         ND         61.7         60-130           1,1,2.2-Tetrachloroethane         4.74         0.05         ug/g         ND         119         60-130           1,1,2.2-Tetrachloroethane         3.67         0.05         ug/g         ND         107         60-130           1,1,2.7-Ticholroethane         3.13         0.05         ug/g         ND         130         60-130           1,1,2-Tricholroethane         3.13         0.05         ug/g         ND										
Hexane       2.86       0.05       ug/g       ND       71.5       60-130         Methyl Ethyl Ketone (2-Butanone)       6.96       0.50       ug/g       ND       69.6       50-140         Methyl Isobutyl Ketone       7.81       0.50       ug/g       ND       69.7       50-140         Methyl Isobutyl Ketone       6.97       0.05       ug/g       ND       82.0       60-130         Methyl Isobutyl Ketone       3.67       0.05       ug/g       ND       101       60-130         Styrene       3.67       0.05       ug/g       ND       101       60-130         1,1,2,2-Tetrachloroethane       4.74       0.05       ug/g       ND       107       60-130         Toluene       4.74       0.05       ug/g       ND       78.3       60-130         1,1,2-Teichoroethane       3.13       0.05       ug/g       ND       78.3       60-130         1,1,1-Trichloroethane       3.11       0.05       ug/g       ND       103       50-140         1,1,2-Trichloroethane       3.41       0.05       ug/g       ND       103       50-140         1,1,2-Trichloroethane       3.41       0.05       ug/g       ND										
Methyl Ethyl Ketone (2-Butanone)         6.96         0.50         ug/g         ND         69.6         50-140           Methyl Isobutyl Ketone         7.81         0.50         ug/g         ND         78.1         50-140           Methyl Isobutyl Ketone         6.97         0.05         ug/g         ND         69.7         50-140           Methyl Isobutyl Ketone         3.28         0.05         ug/g         ND         69.7         60-130           Styrene         3.67         0.05         ug/g         ND         101         60-130           1,1,2.2-Tetrachloroethane         4.05         0.05         ug/g         ND         197         60-130           1,1,2.2-Tetrachloroethane         4.74         0.05         ug/g         ND         107         60-130           1,1,2.2-Tetrachloroethane         3.67         0.05         ug/g         ND         107         60-130           1,1,2.2-Trichloroethane         3.13         0.05         ug/g         ND         77         60-130           1,1,2-Trichloroethane         3.11         0.05         ug/g         ND         77.7         60-130           1,1,2-Trichloroethane         3.11         0.05         ug/g         ND										
Methyl Isobutyl Ketone       7.81       0.50       ug/g       ND       78.1       50-140         Methyl tert-butyl ether       6.97       0.05       ug/g       ND       69.7       50-140         Methyl tert-butyl ether       3.28       0.05       ug/g       ND       82.0       60-130         Styrene       3.67       0.05       ug/g       ND       101       60-130         1,1,2-2-Fetrachloroethane       4.74       0.05       ug/g       ND       89.0       60-130         1,1,2-2-Fetrachloroethane       3.66       0.05       ug/g       ND       119       60-130         1,1,1-Trichloroethane       3.66       0.05       ug/g       ND       107       60-130         Toluene       4.28       0.05       ug/g       ND       78.3       60-130         1,1,1-Trichloroethane       3.11       0.05       ug/g       ND       77.7       60-130         1,1,1-Trichloroethane       3.11       0.05       ug/g       ND       103       50-140         Vinyl chloride       3.22       0.02       ug/g       ND       106       60-130         oxylenes       8.46       0.05       ug/g       ND       106<										
Methyl tert-butyl ether       6.97       0.05       ug'g       ND       69.7       50-140         Methylene Chloride       3.28       0.05       ug/g       ND       82.0       60-130         Styrene       3.67       0.05       ug/g       ND       91.7       60-130         1,1,2.2-Tetrachloroethane       4.05       0.05       ug/g       ND       101       60-130         Tetrachloroethane       4.74       0.05       ug/g       ND       89.0       60-130         Tetrachloroethane       3.56       0.05       ug/g       ND       89.0       60-130         Tetrachloroethane       3.56       0.05       ug/g       ND       107       60-130         Toluene       4.28       0.05       ug/g       ND       78.3       60-130         1,1,2-Trichloroethane       3.13       0.05       ug/g       ND       78.3       60-130         Trichloroethane       3.11       0.05       ug/g       ND       85.2       60-130         Trichloroethane       3.11       0.05       ug/g       ND       103       50-140         Vinyl chloride       3.32       0.02       ug/g       ND       83.1       50-										
Methylene Chloride         3.28         0.05         ug'g         ND         82.0         60-130           Styrene         3.67         0.05         ug/g         ND         91.7         60-130           1,1,1.2-Tetrachloroethane         4.05         0.05         ug/g         ND         101         60-130           1,1,2.2-Tetrachloroethane         4.05         0.05         ug/g         ND         89.0         60-130           Tetrachloroethylene         3.66         0.05         ug/g         ND         119         60-130           Tetrachloroethylene         3.66         0.05         ug/g         ND         89.0         60-130           Toluene         4.28         0.05         ug/g         ND         78.3         60-130           1,1,2-Trichloroethane         3.13         0.05         ug/g         ND         78.3         60-130           Trichlorofluoromethane         3.11         0.05         ug/g         ND         103         50-140           Vinyl chloride         3.32         0.02         ug/g         ND         133         60-130           Surrogate: <i>Heromofluorobenzene</i> 2.59         ug/g         ND         166         60-130										
Styrene       3.67       0.05       ug/g       ND       91.7       60-130         1,1,1,2-Tetrachloroethane       4.05       0.05       ug/g       ND       101       60-130         1,1,2,2-Tetrachloroethane       4.74       0.05       ug/g       ND       119       60-130         Tetrachloroethylene       3.56       0.05       ug/g       ND       89.0       60-130         Toluene       4.28       0.05       ug/g       ND       107       60-130         1,1,1-Trichloroethane       3.13       0.05       ug/g       ND       85.2       60-130         Trichloroethylene       3.41       0.05       ug/g       ND       85.2       60-130         Trichloroethane       3.11       0.05       ug/g       ND       83.1       50-140         Viny Ichloride       3.32       0.02       ug/g       ND       83.1       50-140         Surrogate: A-Bromofluorobenzene       8.46       0.05       ug/g       ND       106       60-130         Surrogate: Toluene-d8       8.46       0.05       ug/g       ND       106       60-130         Surrogate: Toluene-d8       2.97       ug/g       91.1       50-140										
1,1,1,2-Tetrachloroethane4.050.05ug'gND10160-1301,1,2,2-Tetrachloroethane4.740.05ug/gND11960-130Tetrachloroethylene3.560.05ug/gND89.060-130Toluene4.280.05ug/gND10760-1301,1,1-Trichloroethane3.130.05ug/gND78.360-1301,1,2-Trichloroethane3.410.05ug/gND77.760-1301,1,2-Trichloroethane3.110.05ug/gND77.760-130Trichlorofluoromethane4.110.05ug/gND10350-140Vinyl chloride3.220.02ug/gND10660-130o-Xylene4.530.05ug/gND11360-130Surrogate: Iblromofluoromethane2.59ug/gND11360-130Surrogate: Iblromofluoromethane3.410.02ug/gND11360-130o-Xylene4.530.05ug/gND11360-130Surrogate: Iblromofluoromethane3.67ug/g11150-140Surrogate: Iblromofluoromethane3.410.02ug/gND13360-130Surrogate: Iblromofluoromethane3.67ug/g11150-140Surrogate: Iblromofluoromethane3.67ug/g11150-140Surrogate: Iblromofluoromethane3.610.05ug/gND10260-130<	-									
1,1,2,2-Tetrachloroethane4.740.05ug'gND11960-130Tetrachloroethylene3.560.05ug/gND89.060-130Toluene4.280.05ug/gND10760-1301,1,1-Trichloroethane3.130.05ug/gND78.360-1301,1,2-Trichloroethane3.410.05ug/gND77.760-130Trichloroethylene3.110.05ug/gND77.760-130Trichloroethylene3.110.05ug/gND10350-140Trichloroethylene3.220.02ug/gND83.150-140m.p-Xylenes8.460.05ug/gND11360-130o-Xylene2.59ug/gND11360-130Surrogate: Toluene-d82.91ug/g71.150-140Benzene3.410.02ug/gND11360-130Surrogate: Toluene-d82.91ug/g71.150-140Benzene3.410.02ug/gND13660-130Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND10260-130Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND10260-130Toluene3.420.05ug/gND10260-130Toluene4.280.05u										
Tetrachloroethylene3.560.05ug/gND89.060-130Toluene4.280.05ug/gND10760-1301,1,1-Trichloroethane3.130.05ug/gND78.360-1301,1,2-Trichloroethane3.410.05ug/gND85.260-130Trichloroethylene3.110.05ug/gND77.760-130Trichloroethylene3.110.05ug/gND10350-140Vinyl chloride3.320.02ug/gND83.150-140wnp-Xylenes8.460.05ug/gND11360-130o-Xylene4.530.05ug/gND11360-130Surrogate:4.Bromfluoromethane3.57ug/g9.150-140Surrogate:Dibromfluoromethane3.410.02ug/gND132Benzene3.410.02ug/gND85.260-130Surrogate:Jubromfluoromethane3.57ug/g9.150-140Benzene3.410.02ug/gND85.260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10660-130Toluene4.530.05ug/gND10660-130<										
Toluene4.280.05ug/gND10760-1301,1,1-Trichloroethane3.130.05ug/gND78.360-1301,1,2-Trichloroethane3.410.05ug/gND85.260-130Trichloroethylene3.110.05ug/gND77.760-130Trichloroethylene3.110.05ug/gND10350-140Vinyl chloride3.320.02ug/gND83.150-140w.p-Xylenes8.460.05ug/gND10660-130o-Xylene2.59ug/gND11360-130Surrogate: 4-Bromofluorobenzene2.59ug/g91.150-140Surrogate: 70/uen-d82.97ug/g91.150-140Benzene3.410.02ug/gND10260-130Toluene4.280.05ug/gND10260-130m,p-Xylenes3.410.02ug/gND10260-130surrogate: Toluene-d82.97ug/g91.150-140Benzene3.410.02ug/gND10260-130Toluene4.280.05ug/gND10760-130m,p-Xylenes8.460.05ug/gND10660-130o-tylene4.530.05ug/gND10660-130o-tylene4.660.05ug/gND10660-130o-tylene4.530.05ug/g </td <td></td>										
1,1,1-Trichloroethane3.130.05ug/gND78.360-1301,1,2-Trichloroethane3.410.05ug/gND85.260-130Trichloroethylene3.110.05ug/gND77.760-130Trichloroethylene3.110.05ug/gND10350-140Vinyl chloride3.320.02ug/gND83.150-140m,p-Xylenes8.460.05ug/gND11360-130o-Xylene2.59ug/gND11360-130Surrogate: A-Bromofluorobenzene2.59ug/g71.150-140Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND85.260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10260-130o-Xylenes8.460.05ug/gND10760-130o-Xylenes8.460.05ug/gND10660-130o-Xylenes8.460.05ug/gND10660-130o-Xylenes8.460.05ug/gND10660-130o-Xylenes8.460.05ug/gND10660-130o-Xylenes8.460.05 <td>-</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	-									
1,1,2-Trichloroethane3.410.05ug'gND85.260-130Trichloroethylene3.110.05ug/gND77.760-130Trichlorofthare4.110.05ug/gND10350-140Vinyl chloride3.320.02ug/gND83.150-140m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130Surrogate: 4-Bromofluorobenzene2.59ug/gND11360-140Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND85.260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10260-130o-Xylenes6.460.05ug/gND10260-130Toluene4.280.05ug/gND10260-130m,p-Xylenes6.460.05ug/gND10760-130m,p-Xylenes6.460.05ug/gND10660-130o-Xylene4.530.05ug/gND10660-130o-Xylenes6.460.05ug/gND10660-130o-Xylenes6.4530.05ug/gND10660-130o-Xylene6.530.05ug/gND10660-130o-Xylene6.53 <td></td>										
Tichloroethylene3.110.05ug'gND77.760-130Trichlorofluoromethane4.110.05ug/gND10350-140Vinyl chloride3.320.02ug/gND83.150-140m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130Surrogate: 4-Bromofluorobenzene2.59ug/g80.950-140Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND85.260-130Toluene4.280.05ug/gND10260-130Toluene4.280.05ug/gND10260-130o-Xylenes8.460.05ug/gND10660-130corrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND10260-130Corrogate: Toluene-d80.05ug/gND10760-130Toluene4.280.05ug/gND10660-130o-Xylenes8.460.05ug/gND10660-130o-Xylenes8.460.05ug/gND10660-130o-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND10660-130o-Xylene4.530.05ug/gND </td <td></td>										
Trichlorofluoromethane4.110.05ug/gND10350-140Vinyl chloride3.320.02ug/gND83.150-140m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130Surrogate: 4-Bromofluorobenzene2.59ug/g80.950-140Surrogate: Dibromofluoromethane3.57ug/g11150-140Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND85.260-130Toluene4.080.05ug/gND10260-130m,p-Xylenes4.660.05ug/gND10660-130o-Xylene4.530.05ug/gND10260-130cover4.660.05ug/gND10660-130o-Xylenes8.460.05ug/gND10760-130o-Xylene4.530.05ug/gND10660-130o-Xylene4.530.05ug/gND10660-130o-Xylene4.530.05ug/gND10660-130										
Vinyl chloride3.320.02ug/gND83.150-140m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130Surrogate: 4-Bromofluorobenzene2.59ug/g80.950-140Surrogate: Dibromofluoromethane3.57ug/g11150-140Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND85.260-130Toluene4.080.05ug/gND10260-130Toluene4.280.05ug/gND10760-130m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND10660-130o-Xylene4.530.05ug/gND10660-130	-									
m,p-Xylenes       8.46       0.05       ug/g       ND       106       60-130         o-Xylene       4.53       0.05       ug/g       ND       113       60-130         Surrogate: 4-Bromofluorobenzene       2.59       ug/g       80.9       50-140         Surrogate: Dibromofluoromethane       3.57       ug/g       111       50-140         Surrogate: Toluene-d8       2.91       ug/g       91.1       50-140         Benzene       3.41       0.02       ug/g       ND       85.2       60-130         Ethylbenzene       4.08       0.05       ug/g       ND       102       60-130         Toluene       4.28       0.05       ug/g       ND       102       60-130         m,p-Xylenes       8.46       0.05       ug/g       ND       107       60-130         o-Xylene       8.46       0.05       ug/g       ND       106       60-130         o-Xylene       4.53       0.05       ug/g       ND       106       60-130										
o-Xylene4.530.05ug/gND11360-130Surrogate: 4-Bromofluorobenzene2.59ug/g80.950-140Surrogate: Dibromofluoromethane3.57ug/g11150-140Surrogate: Toluene-d82.91ug/g91.150-140Benzene3.410.02ug/gND85.260-130Ethylbenzene4.080.05ug/gND10260-130Toluene4.280.05ug/gND10760-130m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130	-			ug/g						
Surrogate: 4-Bromofluorobenzene       2.59       ug/g       80.9       50-140         Surrogate: Dibromofluoromethane       3.57       ug/g       111       50-140         Surrogate: Toluene-d8       2.91       ug/g       91.1       50-140         Benzene       3.41       0.02       ug/g       ND       85.2       60-130         Ethylbenzene       4.08       0.05       ug/g       ND       102       60-130         Toluene       4.28       0.05       ug/g       ND       107       60-130         m,p-Xylenes       8.46       0.05       ug/g       ND       106       60-130         o-Xylene       4.53       0.05       ug/g       ND       106       60-130		8.46	0.05	ug/g		106				
Surrogate:         Dibromofluoromethane         3.57         ug/g         111         50-140           Surrogate:         Toluene-d8         2.91         ug/g         91.1         50-140           Benzene         3.41         0.02         ug/g         ND         85.2         60-130           Ethylbenzene         4.08         0.05         ug/g         ND         102         60-130           Toluene         4.28         0.05         ug/g         ND         107         60-130           m,p-Xylenes         8.46         0.05         ug/g         ND         106         60-130           o-Xylene         4.53         0.05         ug/g         ND         106         60-130		4.53	0.05	ug/g	ND	113				
Surrogate: Toluene-d8         2.91         ug/g         91.1         50-140           Benzene         3.41         0.02         ug/g         ND         85.2         60-130           Ethylbenzene         4.08         0.05         ug/g         ND         102         60-130           Toluene         4.28         0.05         ug/g         ND         107         60-130           m,p-Xylenes         8.46         0.05         ug/g         ND         106         60-130           o-Xylene         4.53         0.05         ug/g         ND         113         60-130	5									
Benzene         3.41         0.02         ug/g         ND         85.2         60-130           Ethylbenzene         4.08         0.05         ug/g         ND         102         60-130           Toluene         4.28         0.05         ug/g         ND         107         60-130           m,p-Xylenes         8.46         0.05         ug/g         ND         106         60-130           o-Xylene         4.53         0.05         ug/g         ND         113         60-130	•									
Ethylbenzene4.080.05ug/gND10260-130Toluene4.280.05ug/gND10760-130m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130	•		0.05							
Toluene4.280.05ug/gND10760-130m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130										
m,p-Xylenes8.460.05ug/gND10660-130o-Xylene4.530.05ug/gND11360-130	-									
o-Xylene 4.53 0.05 ug/g ND 113 60-130										
Surrogate: Toluene-d8 2.91 ug/g 91.1 50-140	-		0.05		ND					
	Surrogate: Toluene-d8	2.91		ug/g		91.1	50-140			



#### Sample Qualifiers :

1: Elevated detection limits due to the nature of the sample matrix.

2: GC-FID signal did not return to baseline by C50

### QC Qualifiers :

- QM-05 : The spike recovery was outside acceptance limits for the matrix spike due to matrix interference.
- QM-06 : Due to noted non-homogeneity of the QC sample matrix, the spike recoveries were out side the accepted range. Batch data accepted based on other QC.

### Sample Data Revisions

None

### Work Order Revisions / Comments:

None

### Other Report Notes:

n/a: not applicable ND: Not Detected MDL: Method Detection Limit Source Result: Data used as source for matrix and duplicate samples %REC: Percent recovery. RPD: Relative percent difference. NC: Not Calculated

Soil results are reported on a dry weight basis when the units are denoted with 'dry'. Where %Solids is reported, moisture loss includes the loss of volatile hydrocarbons.

#### CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.

- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.
- When reported, data for F4G has been processed using a silica gel cleanup.

	Paracel ID: 2035556 Laurent Blvd. ario K1G 4J8 )-1947 paracellabs.com Hlabs.com	er Ch
Client Name: PATERSON	Project Ref: PE4614	
Contact Name: Marx D'Army	Quote #:	Т
Address: 154 Colonnade Road	<sup>PO #:</sup> 30555	🗆 1 day
134 Coronnace Icoac	E-mail: mdarcy @patersong roup. ca	2 day
Telephone: 613 - 226-7381	hickory apartersent roup. Ca	Date Requir

Address: 154 Colonnade Telephone: 613 - 226-738		- 		PO#: 30555 E-mail: Mdarcy@patersong-cup.ca									1 day     1 day     2 day     Date Required:			] 3 day Regular			
Regulation 153/04	Other I	Regulation		Aatrix 1	[vne:	s (Soil/S	Sed ) GW (Gr	ound Water)						-					
Table 1 Res/Park Med/Fine	🗆 REG 558	D PWQO	1		rface \	Nater)	SS (Storm/San	itary Sewer)						Rei	quired	Analys	iis		
Table 2 Ind/Comm Coarse Coarse	CCME				P (f	Paint) A	(Air) O (Oth	er)		1			Τ	Τ	Γ	Γ			
🛿 Table 3 🗌 Agri/Other	🗆 SU - Sani	SU - Storm			s				BTEX										
Table	Mun:			au	Containers		Sample	Taken	F1-F4+BTEX			by ICP			>				
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Sample ID/Locatio	n Name		Matrix	Air	t of		Date	Time	PHCs	vocs	PAHs	Metals	Hg 1	B (H	K				2
1 B1411 - 20 - AUI			S		1	AUG	24/2020						Τ	Τ	$\overline{}$	<b></b>			,
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Chain of Custody (Env.) xlsx						Re	vision 3.0												

Chain Of Custody (Lab Use Only)

Page <u></u>of <u></u>

Nº 125769



RELIABLE.

# Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Mark D'Arcy

Client PO: 26843 Project: PE4614 Custody: 122294

Report Date: 10-Jun-2019 Order Date: 4-Jun-2019

Order #: 1923235

This Certificate of Analysis contains analytical data applicable to the following samples as submitted:

Paracel ID	Client ID
1923235-01	BH2-GW1
1923235-02	BH3-GW1
1923235-03	BH4-GW1

Approved By:

Dale Robertson, BSc Laboratory Director

Any use of these results implies your agreement that our total liability in connection with this work, however arising, shall be limited to the amount paid by you for this work, and that our employees or agents shall not under any circumstances be liable to you in connection with this work.



# **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
PHC F1	CWS Tier 1 - P&T GC-FID	7-Jun-19	7-Jun-19
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	5-Jun-19	7-Jun-19
REG 153: PAHs by GC-MS	EPA 625 - GC-MS, extraction	10-Jun-19	10-Jun-19
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	7-Jun-19	7-Jun-19

Report Date: 10-Jun-2019 Order Date: 4-Jun-2019



Order #: 1923235

Report Date: 10-Jun-2019 Order Date: 4-Jun-2019

51del Date: 4-5dl1-2013

	Client ID: Sample Date: Sample ID:	BH2-GW1 03-Jun-19 09:00 1923235-01	BH3-GW1 03-Jun-19 12:00 1923235-02	BH4-GW1 03-Jun-19 09:00 1923235-03	- - -
	MDL/Units	Water	Water	Water	-
Volatiles	"				
Acetone	5.0 ug/L	<5.0	<5.0	<5.0	-
Benzene	0.5 ug/L	<0.5	<0.5	<0.5	-
Bromodichloromethane	0.5 ug/L	<0.5	<0.5	<0.5	-
Bromoform	0.5 ug/L	<0.5	<0.5	<0.5	-
Bromomethane	0.5 ug/L	<0.5	<0.5	<0.5	-
Carbon Tetrachloride	0.2 ug/L	<0.2	<0.2	<0.2	-
Chlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
Chloroform	0.5 ug/L	<0.5	<0.5	<0.5	-
Dibromochloromethane	0.5 ug/L	<0.5	<0.5	<0.5	-
Dichlorodifluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	-
1,2-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,3-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,4-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,1-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-
1,2-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-
1,1-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	-
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	-
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,2-Dichloropropane	0.5 ug/L	<0.5	<0.5	<0.5	-
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	-
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,3-Dichloropropene, total	0.5 ug/L	<0.5	<0.5	<0.5	-
Ethylbenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
Ethylene dibromide (dibromoethar	0.2 ug/L	<0.2	<0.2	<0.2	-
Hexane	1.0 ug/L	<1.0	<1.0	<1.0	-
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	<5.0	<5.0	-
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	<5.0	<5.0	-
Methyl tert-butyl ether	2.0 ug/L	<2.0	<2.0	<2.0	-
Methylene Chloride	5.0 ug/L	<5.0	<5.0	<5.0	-
Styrene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-
Tetrachloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	-
Toluene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-



Order #: 1923235

Report Date: 10-Jun-2019 Order Date: 4-Jun-2019

	Client ID: Sample Date: Sample ID:	BH2-GW1 03-Jun-19 09:00 1923235-01	BH3-GW1 03-Jun-19 12:00 1923235-02	BH4-GW1 03-Jun-19 09:00 1923235-03	-
	MDL/Units	Water	Water	Water	-
1,1,2-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-
Trichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	-
Trichlorofluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	-
Vinyl chloride	0.5 ug/L	<0.5	<0.5	<0.5	-
m,p-Xylenes	0.5 ug/L	<0.5	<0.5	<0.5	-
o-Xylene	0.5 ug/L	<0.5	<0.5	<0.5	-
Xylenes, total	0.5 ug/L	<0.5	<0.5	<0.5	-
4-Bromofluorobenzene	Surrogate	113%	107%	110%	-
Dibromofluoromethane	Surrogate	91.1%	92.9%	91.8%	-
Toluene-d8	Surrogate	109%	101%	111%	-
Hydrocarbons			1	1	<b>1</b>
F1 PHCs (C6-C10)	25 ug/L	<25	<25	<25	-
F2 PHCs (C10-C16)	100 ug/L	<100	<100	<100	-
F3 PHCs (C16-C34)	100 ug/L	<100	<100	<100	-
F4 PHCs (C34-C50)	100 ug/L	<100	<100	<100	-
Semi-Volatiles					
Acenaphthene	0.05 ug/L	<0.05	<0.05	<0.05	-
Acenaphthylene	0.05 ug/L	<0.05	<0.05	<0.05	-
Anthracene	0.01 ug/L	0.05	<0.01	<0.01	-
Benzo [a] anthracene	0.01 ug/L	<0.01	<0.01	<0.01	-
Benzo [a] pyrene	0.01 ug/L	0.03	<0.01	<0.01	-
Benzo [b] fluoranthene	0.05 ug/L	<0.05	<0.05	<0.05	-
Benzo [g,h,i] perylene	0.05 ug/L	<0.05	<0.05	<0.05	-
Benzo [k] fluoranthene	0.05 ug/L	<0.05	<0.05	<0.05	-
Chrysene	0.05 ug/L	<0.05	<0.05	<0.05	-
Dibenzo [a,h] anthracene	0.05 ug/L	<0.05	<0.05	<0.05	-
Fluoranthene	0.01 ug/L	0.07	<0.01	<0.01	-
Fluorene	0.05 ug/L	<0.05	<0.05	<0.05	-
Indeno [1,2,3-cd] pyrene	0.05 ug/L	<0.05	<0.05	<0.05	-
1-Methylnaphthalene	0.05 ug/L	<0.05	<0.05	<0.05	-
2-Methylnaphthalene	0.05 ug/L	<0.05	<0.05	<0.05	-
Methylnaphthalene (1&2)	0.10 ug/L	<0.10	<0.10	<0.10	-
Naphthalene	0.05 ug/L	<0.05	<0.05	<0.05	-
Phenanthrene	0.05 ug/L	<0.05	<0.05	<0.05	-
Pyrene	0.01 ug/L	0.06	<0.01	<0.01	-
2-Fluorobiphenyl	Surrogate	88.9%	95.4%	90.4%	-
Terphenyl-d14	Surrogate	109%	121%	115%	-



Order #: 1923235

Report Date: 10-Jun-2019

Order Date: 4-Jun-2019

Project Description: PE4614

# Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
F2 PHCs (C10-C16)	ND	100	ug/L						
F3 PHCs (C16-C34)	ND	100	ug/L						
F4 PHCs (C34-C50)	ND	100	ug/L						
Semi-Volatiles									
Acenaphthene	ND	0.05	ug/L						
Acenaphthylene	ND	0.05	ug/L						
Anthracene	ND	0.01	ug/L						
Benzo [a] anthracene	ND	0.01	ug/L						
Benzo [a] pyrene	ND	0.01	ug/L						
Benzo [b] fluoranthene	ND	0.05	ug/L						
Benzo [g,h,i] perylene	ND	0.05	ug/L						
Benzo [k] fluoranthene	ND	0.05	ug/L						
Chrysene	ND	0.05	ug/L						
Dibenzo [a,h] anthracene	ND	0.05	ug/L						
Fluoranthene	ND	0.01	ug/L						
Fluorene	ND ND	0.05 0.05	ug/L						
Indeno [1,2,3-cd] pyrene 1-Methylnaphthalene	ND	0.05	ug/L ug/L						
2-Methylnaphthalene	ND	0.05	ug/L						
Methylnaphthalene (1&2)	ND	0.00	ug/L						
Naphthalene	ND	0.05	ug/L						
Phenanthrene	ND	0.05	ug/L						
Pyrene	ND	0.01	ug/L						
Surrogate: 2-Fluorobiphenyl	19.3	0.01	ug/L		96.4	50-140			
Surrogate: Terphenyl-d14	23.5		ug/L		117	50-140			
Volatiles	2010		ug/ =			00 1 10			
		5.0							
Acetone Benzene	ND ND	5.0 0.5	ug/L ug/L						
Bromodichloromethane	ND	0.5	ug/L						
Bromoform	ND	0.5	ug/L						
Bromomethane	ND	0.5	ug/L						
Carbon Tetrachloride	ND	0.2	ug/L						
Chlorobenzene	ND	0.5	ug/L						
Chloroform	ND	0.5	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dichlorobenzene	ND	0.5	ug/L						
1,3-Dichlorobenzene	ND	0.5	ug/L						
1,4-Dichlorobenzene	ND	0.5	ug/L						
1,1-Dichloroethane	ND	0.5	ug/L						
1,2-Dichloroethane	ND	0.5	ug/L						
1,1-Dichloroethylene	ND	0.5	ug/L						
cis-1,2-Dichloroethylene	ND	0.5	ug/L						
trans-1,2-Dichloroethylene 1.2-Dichloropropane	ND ND	0.5	ug/L						
cis-1,3-Dichloropropylene	ND ND	0.5 0.5	ug/L ug/L						
trans-1,3-Dichloropropylene	ND	0.5 0.5	ug/L ug/L						
1,3-Dichloropropene, total	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Ethylene dibromide (dibromoethane,	ND	0.2	ug/L						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Isobutyl Ketone	ND	5.0	ug/L						
Methyl tert-butyl ether	ND	2.0	ug/L						
Methylene Chloride	ND	5.0	ug/L						
Styrene	ND	0.5	ug/L						
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L						



Report Date: 10-Jun-2019 Order Date: 4-Jun-2019

Older Date: 4-Juli-2019

Project Description: PE4614

# Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Tetrachloroethylene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
1,1,1-Trichloroethane	ND	0.5	ug/L						
1,1,2-Trichloroethane	ND	0.5	ug/L						
Trichloroethylene	ND	0.5	ug/L						
Trichlorofluoromethane	ND	1.0	ug/L						
Vinyl chloride	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
Xylenes, total	ND	0.5	ug/L						
Surrogate: 4-Bromofluorobenzene	85.8		ug/L		107	50-140			
Surrogate: Dibromofluoromethane	55.0		ug/L		68.8	50-140			
Surrogate: Toluene-d8	84.7		ug/L		106	50-140			



Order #: 1923235

Report Date: 10-Jun-2019

Order Date: 4-Jun-2019

Project Description: PE4614

# Method Quality Control: Duplicate

AnalyteResultLimitUnitsResult%RECLimitRPDLimitNotesHydrocarbons F1 PHCs (C6-C10)ND25ug/LND30VolatilesAcetoneND5.0ug/LND30BenzeneND0.5ug/LND30Bromodichloromethane5.530.5ug/LND30BromodichloromethaneND0.5ug/LND30BromodichloromethaneND0.5ug/LND30Carbon TetrachlorideND0.2ug/LND30Carbon TetrachlorideND0.2ug/LND30ChlorobenzeneND0.5ug/LND30Chloroform16.70.5ug/LND30DibromochloromethaneND0.5ug/LND30DichlorodifluoromethaneND0.5ug/LND301.3-DichlorobenzeneND0.5ug/LND301.4-DichlorobenzeneND0.5ug/LND301.4-DichlorobenzeneND0.5ug/LND301.4-DichlorobenaeND0.5ug/LND301.4-DichloroethaneND0.5ug/LND301.1-DichloroethaneND0.5ug/LND301.1-DichloroethaneND0.5ug/LND301.1-DichloroethaneND
F1         PHCs (C6-C10)         ND         25         ug/L         ND         30           Volatiles
Volatiles         ND         5.0         ug/L         ND         30           Benzene         ND         0.5         ug/L         ND         30           Bromodichloromethane         5.53         0.5         ug/L         ND         30           Bromoform         ND         0.5         ug/L         ND         30           Bromoform         ND         0.5         ug/L         ND         30           Bromomethane         ND         0.5         ug/L         ND         30           Bromomethane         ND         0.5         ug/L         ND         30           Carbon Tetrachloride         ND         0.5         ug/L         ND         30           Chlorobenzene         ND         0.5         ug/L         ND         30           Chloroform         16.7         0.5         ug/L         ND         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene<
Acetone         ND         5.0         ug/L         ND         30           Benzene         ND         0.5         ug/L         ND         30           Bromodichloromethane         5.53         0.5         ug/L         4.15         28.5         30           Bromoform         ND         0.5         ug/L         ND         30         30           Bromoferm         ND         0.5         ug/L         ND         30         30           Bromomethane         ND         0.5         ug/L         ND         30         30           Carbon Tetrachloride         ND         0.2         ug/L         ND         30         30           Chlorobenzene         ND         0.5         ug/L         ND         30         30           Chloroform         16.7         0.5         ug/L         ND         30         30           Dichlorodifluoromethane         ND         0.5         ug/L         ND         30         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30         30
Benzene         ND         0.5         ug/L         ND         30           Bromodichloromethane         5.53         0.5         ug/L         4.15         28.5         30           Bromoform         ND         0.5         ug/L         ND         30           Bromoform         ND         0.5         ug/L         ND         30           Bromomethane         ND         0.5         ug/L         ND         30           Carbon Tetrachloride         ND         0.2         ug/L         ND         30           Chlorobenzene         ND         0.5         ug/L         ND         30           Chloroform         16.7         0.5         ug/L         ND         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30 </td
Bromodichloromethane         5.53         0.5         ug/L         4.15         28.5         30           Bromoform         ND         0.5         ug/L         ND         30           Bromomethane         ND         0.5         ug/L         ND         30           Carbon Tetrachloride         ND         0.2         ug/L         ND         30           Chlorobenzene         ND         0.5         ug/L         ND         30           Chloroform         16.7         0.5         ug/L         ND         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND
Bromoform         ND         0.5         ug/L         ND         30           Bromomethane         ND         0.5         ug/L         ND         30           Carbon Tetrachloride         ND         0.2         ug/L         ND         30           Chlorobenzene         ND         0.5         ug/L         ND         30           Chloroform         16.7         0.5         ug/L         ND         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30
Bromomethane         ND         0.5         ug/L         ND         30           Carbon Tetrachloride         ND         0.2         ug/L         ND         30           Chlorobenzene         ND         0.5         ug/L         ND         30           Chloroform         16.7         0.5         ug/L         ND         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
Carbon Tetrachloride         ND         0.2         ug/L         ND         30           Chlorobenzene         ND         0.5         ug/L         ND         30           Chloroform         16.7         0.5         ug/L         14.2         16.6         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         N
Chlorobenzene         ND         0.5         ug/L         ND         30           Chloroform         16.7         0.5         ug/L         14.2         16.6         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         1.0         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
Chloroform         16.7         0.5         ug/L         14.2         16.6         30           Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         1.0         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
Dibromochloromethane         ND         0.5         ug/L         ND         30           Dichlorodifluoromethane         ND         1.0         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
Dichlorodifluoromethane         ND         1.0         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
1,3-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,2-Dichlorobenzene         ND         0.5         ug/L         ND         30
1,4-Dichlorobenzene         ND         0.5         ug/L         ND         30           1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
1,1-Dichloroethane         ND         0.5         ug/L         ND         30           1,2-Dichloroethane         ND         0.5         ug/L         ND         30
1,2-Dichloroethane ND 0.5 ug/L ND 30
1,1-Dichloroethylene         ND         0.5         ug/L         ND         30           cis-1,2-Dichloroethylene         ND         0.5         ug/L         ND         30
trans-1,2-Dichloroethylene ND 0.5 ug/L ND 30
1,2-Dichloropropane ND 0.5 ug/L ND 30
cis-1,3-Dichloropropylene ND 0.5 ug/L ND 30
trans-1,3-Dichloropropylene ND 0.5 ug/L ND 30
Ethylbenzene ND 0.5 ug/L ND 30
Ethylene dibromide (dibromoethane, ND 0.2 ug/L ND 30
Hexane ND 1.0 ug/L ND 30
Methyl Ethyl Ketone (2-Butanone) ND 5.0 ug/L ND 30
Methyl Isobutyl Ketone ND 5.0 ug/L ND 30
Methyl tert-butyl ether ND 2.0 ug/L ND 30
Methylene Chloride ND 5.0 ug/L ND 30
Styrene ND 0.5 ug/L ND 30
1,1,1,2-Tetrachloroethane ND 0.5 ug/L ND 30
1,1,2,2-Tetrachloroethane ND 0.5 ug/L ND 30
Tetrachloroethylene ND 0.5 ug/L ND 30
Toluene ND 0.5 ug/L ND 30
1,1,1-Trichloroethane ND 0.5 ug/L ND 30
1,1,2-Trichloroethane ND 0.5 ug/L ND 30
Trichloroethylene ND 0.5 ug/L ND 30
Trichlorofluoromethane ND 1.0 ug/L ND 30
Vinyl chloride ND 0.5 ug/L ND 30
m,p-Xylenes ND 0.5 ug/L ND 30
o-Xylene ND 0.5 ug/L ND 30
Surrogate: 4-Bromofluorobenzene 81.7 ug/L 102 50-140
Surrogate: Dibromofluoromethane 79.6 ug/L 99.5 50-140
Surrogate: Toluene-d8 80.9 ug/L 101 50-140



#### Certificate of Analysis **Client: Paterson Group Consulting Engineers** Client PO: 26843

## Method Quality Control: Spike

Report Date: 10-Jun-2019

Order Date: 4-Jun-2019

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1660	25	ug/L		82.8	68-117			
F2 PHCs (C10-C16)	1290	100	ug/L		80.4	60-140			
F3 PHCs (C16-C34)	3210	100	ug/L		81.8	60-140			
F4 PHCs (C34-C50)	2200	100	ug/L		88.9	60-140			
Semi-Volatiles									
Acenaphthene	3.94	0.05	ug/L		78.8	50-140			
Acenaphthylene	3.64	0.05	ug/L		72.7	50-140			
Anthracene	4.21	0.01	ug/L		84.1	50-140			
Benzo [a] anthracene	3.95	0.01	ug/L		78.9	50-140			
Benzo [a] pyrene	3.30	0.01	ug/L		66.0	50-140			
Benzo [b] fluoranthene	5.26	0.05	ug/L		105	50-140			
Benzo [g,h,i] perylene	3.15	0.05	ug/L		63.1	50-140			
Benzo [k] fluoranthene	4.96	0.05	ug/L		99.3	50-140			
Chrysene	4.49	0.05	ug/L		89.8	50-140			
Dibenzo [a,h] anthracene	3.42	0.05	ug/L		68.5	50-140			
Fluoranthene	3.95	0.01	ug/L		79.0	50-140			
Fluorene	3.87	0.05	ug/L		77.4	50-140			
Indeno [1,2,3-cd] pyrene	3.48	0.05	ug/L		69.5	50-140			
1-Methylnaphthalene	4.40	0.05	ug/L		88.0	50-140			
2-Methylnaphthalene	4.81	0.05	ug/L		96.1	50-140			
Naphthalene	3.90	0.05	ug/L		78.1	50-140			
Phenanthrene	3.88	0.05	ug/L		77.7	50-140			
Pyrene	4.04	0.01	ug/L		80.7	50-140			
Surrogate: 2-Fluorobiphenyl	19.9		ug/L		99.7	50-140			
Volatiles									
Acetone	56.6	5.0	ug/L		56.6	50-140			
Benzene	42.2	0.5	ug/L		105	60-130			
Bromodichloromethane	29.2	0.5	ug/L		73.0	60-130			
Bromoform	25.1	0.5	ug/L		62.8	60-130			
Bromomethane	33.6	0.5	ug/L		84.1	50-140			
Carbon Tetrachloride	28.4	0.2	ug/L		71.0	60-130			
Chlorobenzene	36.3	0.5	ug/L		90.8	60-130			
Chloroform	28.3	0.5	ug/L		70.7	60-130			
Dibromochloromethane	27.2	0.5	ug/L		68.0	60-130			
Dichlorodifluoromethane	24.1	1.0	ug/L		60.2	50-140			
1,2-Dichlorobenzene	44.7	0.5	ug/L		112	60-130			
1,3-Dichlorobenzene	47.2	0.5	ug/L		118	60-130			
1,4-Dichlorobenzene	39.3	0.5	ug/L		98.3	60-130			
1,1-Dichloroethane	30.0	0.5	ug/L		75.1	60-130			
1,2-Dichloroethane	28.3	0.5	ug/L		70.8	60-130			
1,1-Dichloroethylene	33.1	0.5	ug/L		82.7	60-130			
cis-1,2-Dichloroethylene	42.5	0.5	ug/L		106	60-130			
trans-1,2-Dichloroethylene	34.6	0.5	ug/L		86.6	60-130			
1,2-Dichloropropane	40.4	0.5	ug/L		101	60-130			
cis-1,3-Dichloropropylene	46.0	0.5	ug/L		115	60-130			
trans-1,3-Dichloropropylene	43.0	0.5	ug/L		107	60-130			
Ethylbenzene	40.9	0.5	ug/L		102	60-130			
Ethylene dibromide (dibromoethane,	38.1	0.2	ug/L		95.3	60-130			
Hexane	49.8	1.0	ug/L		124	60-130			
Methyl Ethyl Ketone (2-Butanone)	86.5	5.0	ug/L		86.5	50-140			
	88.9	5.0	ug/L		88.9	50-140			



### Order #: 1923235

Report Date: 10-Jun-2019

Order Date: 4-Jun-2019

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Project Description: PE4614

## Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Methyl tert-butyl ether	85.3	2.0	ug/L		85.3	50-140			
Methylene Chloride	28.5	5.0	ug/L		71.2	60-130			
Styrene	46.1	0.5	ug/L		115	60-130			
1,1,1,2-Tetrachloroethane	29.3	0.5	ug/L		73.3	60-130			
1,1,2,2-Tetrachloroethane	38.2	0.5	ug/L		95.5	60-130			
Tetrachloroethylene	38.4	0.5	ug/L		96.0	60-130			
Toluene	36.8	0.5	ug/L		92.0	60-130			
1,1,1-Trichloroethane	28.2	0.5	ug/L		70.6	60-130			
1,1,2-Trichloroethane	36.8	0.5	ug/L		92.0	60-130			
Trichloroethylene	34.6	0.5	ug/L		86.6	60-130			
Trichlorofluoromethane	26.1	1.0	ug/L		65.4	60-130			
Vinyl chloride	29.1	0.5	ug/L		72.7	50-140			
m,p-Xylenes	83.6	0.5	ug/L		105	60-130			
o-Xylene	39.0	0.5	ug/L		97.4	60-130			



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#### Qualifier Notes:

None

Sample Data Revisions

None

#### Work Order Revisions / Comments:

None

#### **Other Report Notes:**

n/a: not applicable ND: Not Detected MDL: Method Detection Limit Source Result: Data used as source for matrix and duplicate samples %REC: Percent recovery. RPD: Relative percent difference.

#### CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.

- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.
- When reported, data for F4G has been processed using a silica gel cleanup.

	GPARACEL				EL WO: 1923235			Head Office 300-2319 St. Laurent Blvd. Ottawa, Ontario K1G 4J8 p: 1-800-749-1947 e: paracel@paracellabs.com					Chain of Custody (Lab Use Only) • N= 122294				
LADURATURIES	LID.		_								4			Paj	ge [	of ]	
Client Name: Paterson			_	Project Reference	PE4	014								Turn	aroun	d Time	e:
Contact Name: Mark D'AYCY				Quote #									011	Day		D31	Day
Address:				PO# 268	43											1	
154 Colonnade St. S.	Limui Address						-D2I	÷.		PRe	gular						
Telephone: (613)226-7381				malarcy@patersongroup.cg					Date	Requir	red:						
Criteria: DO. Reg. 153/04 (As Amended) Table	RSC Filing	10. Rej	g. 558/00	D DPWQO D	CCME II SI	UB (Sto	(m		UB (	Sanita	ry) M	unicipality:		D(	Other:		
Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface	Water) 55 (Storm 3	anitary ?	Sewer) P	(Paint) A (Air) O (	Other)	Re	quire	ed A	naly	ses							
Paracel Order Number:			2			X					T				1	T	T
1923235	i,	Air Volume	of Containers	Sample	Taken	PHCs F1-F4+BTE			s by ICP		VS)						
Sample ID/Location Name	Matrix	Air )	# of	Date	Time	HCs	vocs	PAHs	Metals	Hg	B (HWS)						
BH2-GW	GW		4	Jun3/19	AM	1	1	1	~	-			-	-	-	-	-
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Chain of Custody (Env) - Rev 0.7 Feb. 2016



RELIABLE.

# Certificate of Analysis

## **Paterson Group Consulting Engineers**

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Mark D'Arcy

Client PO: 27107 Project: PE4614 Custody: 122847

Report Date: 25-Jul-2019 Order Date: 19-Jul-2019

Order #: 1929690

This Certificate of Analysis contains analytical data applicable to the following samples as submitted:

Paracel ID	Client ID
1929690-01	BH3-GW2
1929690-02	BH7-GW1
1929690-03	BH9-GW1
1929690-04	BH10-GW1
1929690-05	DUP1

Approved By:

Mark Frata

Mark Foto, M.Sc. Lab Supervisor

Any use of these results implies your agreement that our total liability in connection with this work, however arising, shall be limited to the amount paid by you for this work, and that our employees or agents shall not under any circumstances be liable to you in connection with this work.



Order #: 1929690 Report Date: 25-Jul-2019

Order Date: 19-Jul-2019

Project Description: PE4614

## **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
PHC F1	CWS Tier 1 - P&T GC-FID	23-Jul-19	24-Jul-19
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	24-Jul-19	25-Jul-19
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	23-Jul-19	24-Jul-19



Order #: 1929690

Report Date: 25-Jul-2019 Order Date: 19-Jul-2019

	Client ID: Sample Date:	BH3-GW2 19-Jul-19 12:30	BH7-GW1 19-Jul-19 09:45	BH9-GW1 19-Jul-19 11:15	BH10-GW1 19-Jul-19 10:30
Г	Sample ID: MDL/Units	1929690-01 Water	1929690-02 Water	1929690-03 Water	1929690-04 Water
Volatiles	MDL/Units	Water	Water	Water	Water
Acetone	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Benzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromodichloromethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromoform	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromomethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Carbon Tetrachloride	0.2 ug/L	<0.2	<0.2	<0.2	<0.2
Chlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Chloroform	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Dibromochloromethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Dichlorodifluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
1,2-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,3-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,4-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,2-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,2-Dichloropropane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,3-Dichloropropene, total	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Ethylbenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Ethylene dibromide (dibromoethan	0.2 ug/L	<0.2	<0.2	<0.2	<0.2
Hexane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Methyl tert-butyl ether	2.0 ug/L	<2.0	<2.0	<2.0	<2.0
Methylene Chloride	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Styrene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Tetrachloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Toluene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,1-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5



Order #: 1929690

Report Date: 25-Jul-2019 Order Date: 19-Jul-2019

	r				
	Client ID:	BH3-GW2	BH7-GW1	BH9-GW1	BH10-GW1
	Sample Date:	19-Jul-19 12:30 1929690-01	19-Jul-19 09:45 1929690-02	19-Jul-19 11:15 1929690-03	19-Jul-19 10:30 1929690-04
	Sample ID:	Water	Water	Water	Water
	MDL/Units		1		
1,1,2-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Trichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Trichlorofluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
Vinyl chloride	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
m,p-Xylenes	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
o-Xylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Xylenes, total	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
4-Bromofluorobenzene	Surrogate	110%	124%	115%	114%
Dibromofluoromethane	Surrogate	78.3%	76.6%	83.3%	76.4%
Toluene-d8	Surrogate	113%	115%	108%	116%
Hydrocarbons					
F1 PHCs (C6-C10)	25 ug/L	-	<25	-	<25
F2 PHCs (C10-C16)	100 ug/L	-	<100	-	<100
F3 PHCs (C16-C34)	100 ug/L	-	<100	-	<100
F4 PHCs (C34-C50)	100 ug/L	-	<100	-	<100



Report Date: 25-Jul-2019

Order Date: 19-Jul-2019

	Client ID: Sample Date:	DUP1 19-Jul-19 12:30	-	-	-
-	Sample ID:	1929690-05	-	-	-
	MDL/Units	Water	-	-	-
Volatiles	<b>5</b> 0				[]
Acetone	5.0 ug/L	<5.0	-	-	-
Benzene	0.5 ug/L	<0.5	-	-	-
Bromodichloromethane	0.5 ug/L	<0.5	-	-	-
Bromoform	0.5 ug/L	<0.5	-	-	-
Bromomethane	0.5 ug/L	<0.5	-	-	-
Carbon Tetrachloride	0.2 ug/L	<0.2	-	-	-
Chlorobenzene	0.5 ug/L	<0.5	-	-	-
Chloroform	0.5 ug/L	<0.5	-	-	-
Dibromochloromethane	0.5 ug/L	<0.5	-	-	-
Dichlorodifluoromethane	1.0 ug/L	<1.0	-	-	-
1,2-Dichlorobenzene	0.5 ug/L	<0.5	-	-	-
1,3-Dichlorobenzene	0.5 ug/L	<0.5	-	-	-
1,4-Dichlorobenzene	0.5 ug/L	<0.5	-	-	-
1,1-Dichloroethane	0.5 ug/L	<0.5	-	-	-
1,2-Dichloroethane	0.5 ug/L	<0.5	-	-	-
1,1-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
1,2-Dichloropropane	0.5 ug/L	<0.5	-	-	-
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	-	-	-
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	-	-	-
1,3-Dichloropropene, total	0.5 ug/L	<0.5	-	-	-
Ethylbenzene	0.5 ug/L	<0.5	-	-	-
Ethylene dibromide (dibromoethar	0.2 ug/L	<0.2	-	-	-
Hexane	1.0 ug/L	<1.0	-	-	-
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	-	-	-
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	-	-	-
Methyl tert-butyl ether	2.0 ug/L	<2.0	-	-	-
Methylene Chloride	5.0 ug/L	<5.0	-	-	-
Styrene	0.5 ug/L	<0.5	-	-	-
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	-	-	-
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	-	-	-
Tetrachloroethylene	0.5 ug/L	<0.5	-	-	-
Toluene	0.5 ug/L	<0.5	-	-	-



## Order #: 1929690

Report Date: 25-Jul-2019 Order Date: 19-Jul-2019

	_				
	Client ID:	DUP1	-	-	-
	Sample Date:	19-Jul-19 12:30	-	-	-
	Sample ID:	1929690-05	-	-	-
	MDL/Units	Water	-	-	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	-	-	-
1,1,2-Trichloroethane	0.5 ug/L	<0.5	-	-	-
Trichloroethylene	0.5 ug/L	<0.5	-	-	-
Trichlorofluoromethane	1.0 ug/L	<1.0	-	-	-
Vinyl chloride	0.5 ug/L	<0.5	-	-	-
m,p-Xylenes	0.5 ug/L	<0.5	-	-	-
o-Xylene	0.5 ug/L	<0.5	-	-	-
Xylenes, total	0.5 ug/L	<0.5	-	-	-
4-Bromofluorobenzene	Surrogate	117%	-	-	-
Dibromofluoromethane	Surrogate	78.0%	-	-	-
Toluene-d8	Surrogate	110%	-	-	-



Order #: 1929690

Report Date: 25-Jul-2019

Order Date: 19-Jul-2019

Project Description: PE4614

## Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
F2 PHCs (C10-C16)	ND	100	ug/L						
F3 PHCs (C16-C34)	ND	100	ug/L						
F4 PHCs (C34-C50)	ND	100	ug/L						
Volatiles									
Acetone	ND	5.0	ug/L						
Benzene	ND	0.5	ug/L						
Bromodichloromethane	ND	0.5	ug/L						
Bromoform	ND	0.5	ug/L						
Bromomethane	ND	0.5	ug/L						
Carbon Tetrachloride	ND	0.2	ug/L						
Chlorobenzene	ND	0.5	ug/L						
Chloroform	ND	0.5	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dichlorobenzene	ND	0.5	ug/L						
1,3-Dichlorobenzene	ND	0.5	ug/L						
1,4-Dichlorobenzene	ND	0.5	ug/L						
1,1-Dichloroethane	ND	0.5	ug/L						
1,2-Dichloroethane	ND	0.5	ug/L						
1,1-Dichloroethylene	ND	0.5	ug/L						
cis-1,2-Dichloroethylene	ND	0.5	ug/L						
trans-1,2-Dichloroethylene	ND ND	0.5 0.5	ug/L						
1,2-Dichloropropane cis-1,3-Dichloropropylene	ND	0.5	ug/L						
trans-1,3-Dichloropropylene	ND	0.5	ug/L ug/L						
1,3-Dichloropropene, total	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Ethylene dibromide (dibromoethane	ND	0.2	ug/L						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Isobutyl Ketone	ND	5.0	ug/L						
Methyl tert-butyl ether	ND	2.0	ug/L						
Methylene Chloride	ND	5.0	ug/L						
Styrene	ND	0.5	ug/L						
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L						
Tetrachloroethylene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
1,1,1-Trichloroethane	ND	0.5	ug/L						
1,1,2-Trichloroethane	ND	0.5	ug/L						
Trichloroethylene	ND	0.5	ug/L						
Trichlorofluoromethane	ND	1.0	ug/L						
Vinyl chloride	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
Xylenes, total	ND	0.5	ug/L						
Surrogate: 4-Bromofluorobenzene	96.4		ug/L		120	50-140			
Surrogate: Dibromofluoromethane	61.4		ug/L		76.7	50-140			
Surrogate: Toluene-d8	92.0		ug/L		115	50-140			



Order #: 1929690

Report Date: 25-Jul-2019

Order Date: 19-Jul-2019

Project Description: PE4614

## Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
-		-	01110	ricoult	/on ieo	Liint		Linin	
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L	ND				30	
Volatiles									
Acetone	ND	5.0	ug/L	ND				30	
Benzene	ND	0.5	ug/L	ND				30	
Bromodichloromethane	ND	0.5	ug/L	ND				30	
Bromoform	ND	0.5	ug/L	ND				30	
Bromomethane	ND	0.5	ug/L	ND				30	
Carbon Tetrachloride	ND	0.2	ug/L	ND				30	
Chlorobenzene	ND	0.5	ug/L	ND				30	
Chloroform	ND	0.5	ug/L	ND				30	
Dibromochloromethane	ND	0.5	ug/L	ND				30	
Dichlorodifluoromethane	ND	1.0	ug/L	ND				30	
1,2-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,1-Dichloroethane	ND	0.5	ug/L	ND				30	
1,2-Dichloroethane	ND	0.5	ug/L	ND				30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND				30	
cis-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
trans-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
1,2-Dichloropropane	ND	0.5	ug/L	ND				30	
cis-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30	
trans-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30	
Ethylbenzene	ND	0.5	ug/L	ND				30	
Ethylene dibromide (dibromoethane	ND	0.2	ug/L	ND				30	
Hexane	ND	1.0	ug/L	ND				30	
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L	ND				30	
Methyl Isobutyl Ketone	ND	5.0	ug/L	ND				30	
Methyl tert-butyl ether	ND	2.0	ug/L	ND				30	
Methylene Chloride	ND	5.0	ug/L	ND				30	
Styrene 1,1,1,2-Tetrachloroethane	ND ND	0.5 0.5	ug/L	ND ND				30 30	
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L	ND				30	
Tetrachloroethylene	ND	0.5	ug/L ug/L	ND				30	
Toluene	ND	0.5		ND				30	
1,1,1-Trichloroethane	ND	0.5	ug/L ug/L	ND				30	
1,1,2-Trichloroethane	ND	0.5	ug/L	ND				30	
Trichloroethylene	ND	0.5	ug/L	ND				30	
Trichlorofluoromethane	ND	1.0	ug/L	ND				30	
Vinyl chloride	ND	0.5	ug/L	ND				30	
m,p-Xylenes	ND	0.5	ug/L	ND				30	
o-Xylene	ND	0.5	ug/L	ND				30	
Surrogate: 4-Bromofluorobenzene	91.8	0.0	ug/L		115	50-140			
Surrogate: Dibromofluoromethane	61.6		ug/L		77.0	50-140			
Surrogate: Toluene-d8	91.0		ug/L		114	50-140			
Currogate. Tolucite do	31.0		uy/L		114	50.140			



## Method Quality Control: Spike

Report Date: 25-Jul-2019

Order Date: 19-Jul-2019

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	2060	25	ug/L		103	68-117			
F2 PHCs (C10-C16)	1720	100	ug/L		108	60-140			
F3 PHCs (C16-C34)	3890	100	ug/L		99.3	60-140			
F4 PHCs (C34-C50)	2600	100	ug/L		105	60-140			
Volatiles									
Acetone	68.5	5.0	ug/L		68.5	50-140			
Benzene	31.4	0.5	ug/L		78.5	60-130			
Bromodichloromethane	24.3	0.5	ug/L		60.7	60-130			
Bromoform	34.6	0.5	ug/L		86.5	60-130			
Bromomethane	32.0	0.5	ug/L		80.1	50-140			
Carbon Tetrachloride	28.4	0.2	ug/L		71.0	60-130			
Chlorobenzene	37.2	0.5	ug/L		92.9	60-130			
Chloroform	26.6	0.5	ug/L		66.4	60-130			
Dibromochloromethane	39.6	0.5	ug/L		99.0	60-130			
Dichlorodifluoromethane	30.0	1.0	ug/L		74.9	50-140			
1,2-Dichlorobenzene	43.8	0.5	ug/L		110	60-130			
1,3-Dichlorobenzene	48.1	0.5	ug/L		120	60-130			
1,4-Dichlorobenzene	39.4	0.5	ug/L		98.5	60-130			
1,1-Dichloroethane	32.8	0.5	ug/L		82.0	60-130			
1,2-Dichloroethane	27.4	0.5	ug/L		68.6	60-130			
1,1-Dichloroethylene	33.2	0.5	ug/L		83.0	60-130			
cis-1,2-Dichloroethylene	28.0	0.5	ug/L		70.0	60-130			
trans-1,2-Dichloroethylene	30.7	0.5	ug/L		76.8	60-130			
1,2-Dichloropropane	26.9	0.5	ug/L		67.2	60-130			
cis-1,3-Dichloropropylene	26.8	0.5	ug/L		66.9	60-130			
trans-1,3-Dichloropropylene	28.4	0.5	ug/L		70.9	60-130			
Ethylbenzene	43.0	0.5	ug/L		107	60-130			
Ethylene dibromide (dibromoethane	40.2	0.2	ug/L		100	60-130			
Hexane	31.7	1.0	ug/L		79.3	60-130			
Methyl Ethyl Ketone (2-Butanone)	78.0	5.0	ug/L		78.0	50-140			
Methyl Isobutyl Ketone	62.8	5.0	ug/L		62.8	50-140			
Methyl tert-butyl ether	90.0	2.0	ug/L		90.0	50-140			
Methylene Chloride	32.7	5.0	ug/L		81.8	60-130			
Styrene	39.8	0.5	ug/L		99.6	60-130			
1,1,1,2-Tetrachloroethane	47.0	0.5	ug/L		117	60-130			
1,1,2,2-Tetrachloroethane	44.5	0.5	ug/L		111	60-130			
Tetrachloroethylene	43.6	0.5	ug/L		109	60-130			
Toluene	44.5	0.5	ug/L		111	60-130			
1,1,1-Trichloroethane	26.6	0.5	ug/L		66.6	60-130			
1,1,2-Trichloroethane	29.4	0.5	ug/L		73.4	60-130			
Trichloroethylene	38.0	0.5	ug/L		94.9	60-130			
Trichlorofluoromethane	29.6	1.0	ug/L		74.0	60-130			
Vinyl chloride	45.3	0.5	ug/L		113	50-140			
m,p-Xylenes	74.0	0.5	ug/L		92.4	60-130			
o-Xylene	39.1	0.5	ug/L		97.8	60-130			
Surrogate: 4-Bromofluorobenzene	61.5		ug/L		76.9	50-140			



#### Qualifier Notes:

None

Sample Data Revisions

None

#### Work Order Revisions / Comments:

None

#### **Other Report Notes:**

n/a: not applicable ND: Not Detected MDL: Method Detection Limit Source Result: Data used as source for matrix and duplicate samples %REC: Percent recovery. RPD: Relative percent difference.

#### CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.

- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.
- When reported, data for F4G has been processed using a silica gel cleanup.

GPARACEL	DE	SPC			Paracel II							urent Blvd. K1G 4J8 I7 pellabs.com		(Lab U	f Custody se Outy) 22847
Client Name: Poterson Contact Name: Mark D'Ary				Project Referen	ice: PE	40	14	/			_			Turnarou	of
Address: Telephone: $227 - 7387$ Criteria: $\Box$ O. Reg. 153/04 (As Amended) Table $\Box$ RSC Fi	iling 🛛	O. Reg	g. 558/0(	PO # Email Address:	27/07		orm)	O S	UB	Sanita	V) N	funicinativa	- 0 2 D	ay Required:	🗆 3 Day
Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) SS						T	quire				y) iv	runcipality:		Other	
3 13119 - GWI 4 13110 - GWI	A B B B Matrix A B C C Matrix	Air Volume	UNDUND# of Containers	Date	Тіте 12:30рм 9:45ам 11:15ал 10:30ам	V PHCs F1-F4+BTEX	Vocs	PAHs	Metals by ICP	Hg CrVI	B (HWS)				
6 7 8			đ		12:30pm		✓								
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RELIABLE.

# Certificate of Analysis

## **Paterson Group Consulting Engineers**

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Mark D'Arcy

Client PO: 27660 Project: PE4614 Custody: 122874

Report Date: 1-Aug-2019 Order Date: 31-Jul-2019

Order #: 1931355

This Certificate of Analysis contains analytical data applicable to the following samples as submitted:

Paracel ID **Client ID** 1931355-01 BH8B-GW1

Approved By:

Mark Frata

Mark Foto, M.Sc. Lab Supervisor

Any use of these results implies your agreement that our total liability in connection with this work, however arising, shall be limited to the amount paid by you for this work, and that our employees or agents shall not under any circumstances be liable to you in connection with this work.



Order #: 1931355

Report Date: 01-Aug-2019 Order Date: 31-Jul-2019

Project Description: PE4614

### **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	1-Aug-19	1-Aug-19



Report Date: 01-Aug-2019

Order Date: 31-Jul-2019

г	Client ID: Sample Date: Sample ID:	BH8B-GW1 31-Jul-19 10:10 1931355-01 Water		- - -	- - -
Volatiles	MDL/Units	water	-	-	-
Acetone	5.0 ug/L	<5.0	-	-	_
Benzene	0.5 ug/L	<0.5	_	_	_
Bromodichloromethane	0.5 ug/L	<0.5	_	-	_
Bromoform	0.5 ug/L	<0.5	_	-	_
Bromomethane	0.5 ug/L	<0.5	-	_	_
Carbon Tetrachloride	0.2 ug/L	<0.2	_	-	-
Chlorobenzene	0.5 ug/L	<0.5	_	-	-
Chloroform	0.5 ug/L	<0.5	_	_	_
Dibromochloromethane	0.5 ug/L	<0.5	_	-	
Dichlorodifluoromethane	1.0 ug/L	<1.0	_	_	_
1,2-Dichlorobenzene	0.5 ug/L	<0.5	_	_	_
1,3-Dichlorobenzene	0.5 ug/L	<0.5	-	-	-
1,4-Dichlorobenzene	0.5 ug/L	<0.5	_	_	_
1,1-Dichloroethane	0.5 ug/L	<0.5	_	_	
1,2-Dichloroethane	0.5 ug/L	<0.5	_	_	_
1,1-Dichloroethylene	0.5 ug/L	<0.5	_	_	_
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
1,2-Dichloropropane	0.5 ug/L	<0.5	-	-	-
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	-	-	-
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	-	-	-
1,3-Dichloropropene, total	0.5 ug/L	<0.5	-	-	-
Ethylbenzene	0.5 ug/L	<0.5	-	-	-
Ethylene dibromide (dibromoethan	0.2 ug/L	<0.2	-	-	-
Hexane	1.0 ug/L	<1.0	-	-	-
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	-	-	-
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	-	-	-
Methyl tert-butyl ether	2.0 ug/L	<2.0	-	-	-
Methylene Chloride	5.0 ug/L	<5.0	-	-	-
Styrene	0.5 ug/L	<0.5	-	-	-
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	-	-	-
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	-	-	-
Tetrachloroethylene	0.5 ug/L	<0.5	-	-	-
Toluene	0.5 ug/L	<0.5	-	-	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	-	-	-



## Order #: 1931355

Report Date: 01-Aug-2019 Order Date: 31-Jul-2019

	Client ID:		-	-	-
	Sample Date:	31-Jul-19 10:10	-	-	-
	Sample ID:	1931355-01	-	-	-
	MDL/Units	Water	-	-	-
1,1,2-Trichloroethane	0.5 ug/L	<0.5	-	-	-
Trichloroethylene	0.5 ug/L	<0.5	-	-	-
Trichlorofluoromethane	1.0 ug/L	<1.0	-	-	-
Vinyl chloride	0.5 ug/L	<0.5	-	-	-
m,p-Xylenes	0.5 ug/L	<0.5	-	-	-
o-Xylene	0.5 ug/L	<0.5	-	-	-
Xylenes, total	0.5 ug/L	<0.5	-	-	-
4-Bromofluorobenzene	Surrogate	108%	-	-	-
Dibromofluoromethane	Surrogate	97.8%	-	-	-
Toluene-d8	Surrogate	116%	-	-	-



Order #: 1931355

Report Date: 01-Aug-2019

Order Date: 31-Jul-2019

Project Description: PE4614

## Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Volatiles									
Acetone	ND	5.0	ug/L						
Benzene	ND	0.5	ug/L						
Bromodichloromethane	ND	0.5	ug/L						
Bromoform	ND	0.5	ug/L						
Bromomethane	ND	0.5	ug/L						
Carbon Tetrachloride	ND	0.2	ug/L						
Chlorobenzene	ND	0.5	ug/L						
Chloroform	ND	0.5	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dichlorobenzene	ND	0.5	ug/L						
1,3-Dichlorobenzene	ND	0.5	ug/L						
1,4-Dichlorobenzene	ND	0.5	ug/L						
1,1-Dichloroethane	ND	0.5	ug/L						
1,2-Dichloroethane	ND	0.5	ug/L						
1,1-Dichloroethylene	ND	0.5	ug/L						
cis-1,2-Dichloroethylene	ND	0.5	ug/L						
trans-1.2-Dichloroethylene	ND	0.5	ug/L						
1,2-Dichloropropane	ND	0.5	ug/L						
cis-1,3-Dichloropropylene	ND	0.5	ug/L						
trans-1,3-Dichloropropylene	ND	0.5	ug/L						
1,3-Dichloropropene, total	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Ethylene dibromide (dibromoethane	ND	0.2	ug/L						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Isobutyl Ketone	ND	5.0	ug/L						
Methyl tert-butyl ether	ND	2.0	ug/L						
Methylene Chloride	ND	5.0	ug/L						
Styrene	ND	0.5	ug/L						
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L						
Tetrachloroethylene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
1,1,1-Trichloroethane	ND	0.5	ug/L						
1,1,2-Trichloroethane	ND	0.5	ug/L						
Trichloroethylene	ND	0.5	ug/L						
Trichlorofluoromethane	ND	1.0	ug/L						
Vinyl chloride	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
Xylenes, total	ND	0.5	ug/L						
Surrogate: 4-Bromofluorobenzene	83.4		ug/L		104	50-140			
Surrogate: Dibromofluoromethane	81.0		ug/L		101	50-140			
Surrogate: Toluene-d8	78.1		ug/L		97.6	50-140			
Surregale. Toldene-do	70.1		uy/L		37.0	50-140			



Order #: 1931355

Report Date: 01-Aug-2019 Order Date: 31-Jul-2019

Project Description: PE4614

Method Quality Control: Duplicate

Amelida		Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Volatiles									
Acetone	ND	5.0	ug/L	ND				30	
Benzene	ND	0.5	ug/L	ND				30	
Bromodichloromethane	ND	0.5	ug/L	ND				30	
Bromoform	ND	0.5	ug/L	ND				30	
Bromomethane	ND	0.5	ug/L	ND				30	
Carbon Tetrachloride	ND	0.2	ug/L	ND				30	
Chlorobenzene	ND	0.5	ug/L	ND				30	
Chloroform	ND	0.5	ug/L	ND				30	
Dibromochloromethane	ND	0.5	ug/L	ND				30	
Dichlorodifluoromethane	ND	1.0	ug/L	ND				30	
1,2-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,1-Dichloroethane	ND	0.5	ug/L	ND				30	
1,2-Dichloroethane	ND	0.5	ug/L	ND				30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND				30	
cis-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
trans-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
1,2-Dichloropropane	ND	0.5	ug/L	ND				30	
cis-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30	
trans-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30	
Ethylbenzene	ND	0.5	ug/L	ND				30	
Ethylene dibromide (dibromoethane	ND	0.2	ug/L	ND				30	
Hexane	ND	1.0	ug/L	ND				30	
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L	ND				30	
Methyl Isobutyl Ketone	ND	5.0	ug/L	ND				30	
Methyl tert-butyl ether	ND	2.0	ug/L	ND				30	
Methylene Chloride	ND	5.0	ug/L	ND				30	
Styrene	ND	0.5	ug/L	ND				30	
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L	ND				30	
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L	ND				30	
Tetrachloroethylene	ND	0.5	ug/L	ND				30	
Toluene	ND	0.5	ug/L	ND				30	
1,1,1-Trichloroethane	ND	0.5	ug/L	ND				30	
1,1,2-Trichloroethane	ND	0.5	ug/L	ND				30	
Trichloroethylene	ND	0.5	ug/L	ND				30	
Trichlorofluoromethane	ND	1.0	ug/L	ND				30	
Vinyl chloride	ND	0.5	ug/L	ND				30	
m,p-Xylenes	ND	0.5	ug/L	ND				30	
o-Xylene	ND	0.5	ug/L	ND				30	
Surrogate: 4-Bromofluorobenzene	75.6		ug/L		94.4	50-140			
Surrogate: Dibromofluoromethane	92.1		ug/L		115	50-140			
Surrogate: Toluene-d8	71.2		ug/L		89.0	50-140			



## Method Quality Control: Spike

Report Date: 01-Aug-2019

Order Date: 31-Jul-2019

Analyte	Result	Reporting Limit	Units	Source %REC Result	%REC Limit	RPD	RPD Limit	Notes
Volatiles								
Acetone	58.6	5.0	ug/L	58.6	50-140			
Benzene	41.9	0.5	ug/L	105	60-130			
Bromodichloromethane	31.3	0.5	ug/L	78.2	60-130			
Bromoform	26.3	0.5	ug/L	65.8	60-130			
Bromomethane	26.6	0.5	ug/L	66.6	50-140			
Carbon Tetrachloride	30.6	0.2	ug/L	76.4	60-130			
Chlorobenzene	29.4	0.5	ug/L	73.5	60-130			
Chloroform	29.6	0.5	ug/L	74.0	60-130			
Dibromochloromethane	27.7	0.5	ug/L	69.2	60-130			
Dichlorodifluoromethane	27.3	1.0	ug/L	68.3	50-140			
1,2-Dichlorobenzene	38.8	0.5	ug/L	96.9	60-130			
1,3-Dichlorobenzene	38.1	0.5	ug/L	95.3	60-130			
1,4-Dichlorobenzene	32.2	0.5	ug/L	80.6	60-130			
1,1-Dichloroethane	35.5	0.5	ug/L	88.8	60-130			
1,2-Dichloroethane	32.4	0.5	ug/L	81.1	60-130			
1,1-Dichloroethylene	34.6	0.5	ug/L	86.4	60-130			
cis-1,2-Dichloroethylene	38.1	0.5	ug/L	95.2	60-130			
trans-1,2-Dichloroethylene	33.3	0.5	ug/L	83.2	60-130			
1,2-Dichloropropane	40.2	0.5	ug/L	101	60-130			
cis-1,3-Dichloropropylene	28.6	0.5	ug/L	71.5	60-130			
trans-1,3-Dichloropropylene	33.5	0.5	ug/L	83.8	60-130			
Ethylbenzene	30.9	0.5	ug/L	77.2	60-130			
Ethylene dibromide (dibromoethane	29.7	0.2	ug/L	74.3	60-130			
Hexane	40.8	1.0	ug/L	102	60-130			
Methyl Ethyl Ketone (2-Butanone)	89.0	5.0	ug/L	89.0	50-140			
Methyl Isobutyl Ketone	80.1	5.0	ug/L	80.1	50-140			
Methyl tert-butyl ether	97.6	2.0	ug/L	97.6	50-140			
Methylene Chloride	37.8	5.0	ug/L	94.4	60-130			
Styrene	29.8	0.5	ug/L	74.4	60-130			
1,1,1,2-Tetrachloroethane	36.1	0.5	ug/L	90.3	60-130			
1,1,2,2-Tetrachloroethane	28.3	0.5	ug/L	70.8	60-130			
Tetrachloroethylene	28.4	0.5	ug/L	71.1	60-130			
Toluene	31.6	0.5	ug/L	78.9	60-130			
1,1,1-Trichloroethane	31.6	0.5	ug/L	78.9	60-130			
1,1,2-Trichloroethane	31.1	0.5	ug/L	77.8	60-130			
Trichloroethylene	36.0	0.5	ug/L	90.1	60-130			
Trichlorofluoromethane	26.8	1.0	ug/L	66.9	60-130			
Vinyl chloride	53.6	0.5	ug/L	134	50-140			
m,p-Xylenes	59.7	0.5	ug/L	74.6	60-130			
o-Xylene	30.4	0.5	ug/L	76.0	60-130			
Surrogate: 4-Bromofluorobenzene	62.2		ug/L	77.8	50-140			



#### **Qualifier Notes:**

None

Sample Data Revisions None

#### Work Order Revisions / Comments:

None

#### **Other Report Notes:**

n/a: not applicable ND: Not Detected MDL: Method Detection Limit Source Result: Data used as source for matrix and duplicate samples %REC: Percent recovery. RPD: Relative percent difference. Report Date: 01-Aug-2019 Order Date: 31-Jul-2019 Project Description: PE4614

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LABORATORIES LTD.				Project Reference:	PEHL	14								1		aroun	d Time:
Client Name: Paterson Group Contact Name: Mark D'Arcy				Quote #	110-10	7								X	Day		□ 3 Day
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Criteria: 0. Reg. 153/04 (As Amended) Table 🗆 RSC I	iling 🛛	O. Reg	. 558/00	DPWQO DO	CME USUB	(Sto	rm)	ds	JB (S	Sanita	ary)	Mur	nicipality:		0	Other:	1
Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) S						Req	uire	d A	naly	ses		П					<del>.</del>
Paracel Order Number: 1931355	ńx	Air Volume	of Containers	Sample	Taken	PHCs F1-F4+BTEX	2s	ls	als by ICP		_	(HWS)					
Sample ID/Location Name	Matrix	Air	*	Date	Time	-	VOCs	PAHs	Metals	Τġ	CrVI	B ()		+	-	-	
1 BH8B-GW	GW		2	July 31/19	10:10 AM		Λ			-	+	-		-	+	-	
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Date/Time: July 31, 2019	Tempe	rature:	/	d	PH Tempe	rature	: 16	.8	°C		2.7.2.2		pHV	/erified [	By:		NA

Chain of Custody (Env) - Rev 0.7 Feb. 2016



RELIABLE.

300 - 2319 St. Laurent Blvd Ottawa, ON, K1G 4J8 1-800-749-1947 www.paracellabs.com

## Certificate of Analysis

#### **Paterson Group Consulting Engineers**

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Mark D'Arcy

Client PO: 30732 Project: PE4641 Custody: 128122

Report Date: 11-Sep-2020 Order Date: 4-Sep-2020

Order #: 2036662

This Certificate of Analysis contains analytical data applicable to the following samples as submitted:

Paracel ID	Client ID
2036662-01	BH3-GW3
2036662-02	BH8B-GW1
2036662-03	BH11-20-GW1
2036662-04	BH13-20-GW1
2036662-05	DUP
2036662-04	BH13-20-GW1

Approved By:

Dale Robertson, BSc Laboratory Director

Any use of these results implies your agreement that our total liability in connection with this work, however arising, shall be limited to the amount paid by you for this work, and that our employees or agents shall not under any circumstances be liable to you in connection with this work.



Report Date: 11-Sep-2020 Order Date: 4-Sep-2020

Order #: 2036662

Project Description: PE4641

#### **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
PHC F1	CWS Tier 1 - P&T GC-FID	8-Sep-20	9-Sep-20
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	10-Sep-20	10-Sep-20
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	8-Sep-20	9-Sep-20



Client PO: 30732

Order #: 2036662

Report Date: 11-Sep-2020 Order Date: 4-Sep-2020

Project Description: PE4641

Г	Client ID: Sample Date: Sample ID: MDL/Units	BH3-GW3 03-Sep-20 09:00 2036662-01 Water	BH8B-GW1 03-Sep-20 09:00 2036662-02 Water	BH11-20-GW1 03-Sep-20 09:00 2036662-03 Water	BH13-20-GW1 03-Sep-20 09:00 2036662-04 Water
Volatiles			•	<u>!</u>	
Acetone	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Benzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromodichloromethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromoform	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromomethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Carbon Tetrachloride	0.2 ug/L	<0.2	<0.2	<0.2	<0.2
Chlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Chloroform	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Dibromochloromethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Dichlorodifluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
1,2-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,3-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,4-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,2-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,2-Dichloropropane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,3-Dichloropropene, total	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Ethylbenzene	0.5 ug/L	<0.5	<0.5	0.6	<0.5
Ethylene dibromide (dibromoethane, 1,2-)	0.2 ug/L	<0.2	<0.2	<0.2	<0.2
Hexane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Methyl tert-butyl ether	2.0 ug/L	<2.0	<2.0	<2.0	<2.0
Methylene Chloride	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Styrene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Tetrachloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Toluene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,1-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5

## PARACEL LABORATORIES LTD.

#### Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 30732

Report Date: 11-Sep-2020 Order Date: 4-Sep-2020

Order #: 2036662

Project Description: PE4641

	Client ID:	BH3-GW3	BH8B-GW1	BH11-20-GW1	BH13-20-GW1
	Sample Date:	03-Sep-20 09:00	03-Sep-20 09:00	03-Sep-20 09:00	03-Sep-20 09:00
	Sample ID:	2036662-01	2036662-02	2036662-03	2036662-04
	MDL/Units	Water	Water	Water	Water
1,1,2-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Trichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Trichlorofluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
Vinyl chloride	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
m,p-Xylenes	0.5 ug/L	<0.5	<0.5	0.7	<0.5
o-Xylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Xylenes, total	0.5 ug/L	<0.5	<0.5	0.7	<0.5
4-Bromofluorobenzene	Surrogate	103%	104%	104%	103%
Dibromofluoromethane	Surrogate	96.3%	95.9%	99.3%	98.6%
Toluene-d8	Surrogate	104%	105%	104%	104%
Hydrocarbons			•		
F1 PHCs (C6-C10)	25 ug/L	-	-	<25	<25
F2 PHCs (C10-C16)	100 ug/L	-	-	<100	<100
F3 PHCs (C16-C34)	100 ug/L	-	-	<100	<100
F4 PHCs (C34-C50)	100 ug/L	-	-	<100	<100



Report Date: 11-Sep-2020

Order Date: 4-Sep-2020

Project Description: PE4641

-	Client ID: Sample Date: Sample ID:	DUP 03-Sep-20 09:00 2036662-05			- -
No bella a	MDL/Units	Water	-	-	-
Volatiles Acetone	5.0 ug/L	<5.0			
	0.5 ug/L		-	-	-
Benzene	0.5 ug/L	<0.5	-	-	-
Bromodichloromethane	0.5 ug/L	<0.5	-	-	-
Bromoform	0.5 ug/L	<0.5	-	-	-
Bromomethane	0.3 ug/L	<0.5	-	-	-
Carbon Tetrachloride	0.2 ug/L	<0.2	-	-	-
Chlorobenzene	0.5 ug/L	<0.5	-	-	-
Chloroform	0.5 ug/L	<0.5	-	-	-
Dibromochloromethane	0.5 ug/L 1.0 ug/L	<0.5	-	-	-
Dichlorodifluoromethane	-	<1.0	-	-	-
1,2-Dichlorobenzene	0.5 ug/L	<0.5	-	-	-
1,3-Dichlorobenzene	0.5 ug/L	<0.5	-	-	-
1,4-Dichlorobenzene	0.5 ug/L	<0.5	-	-	-
1,1-Dichloroethane	0.5 ug/L	<0.5	-	-	-
1,2-Dichloroethane	0.5 ug/L	<0.5	-	-	-
1,1-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	-	-	-
1,2-Dichloropropane	0.5 ug/L	<0.5	-	-	-
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	-	-	-
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	-	-	-
1,3-Dichloropropene, total	0.5 ug/L	<0.5	-	-	-
Ethylbenzene	0.5 ug/L	<0.5	-	-	-
Ethylene dibromide (dibromoethane, 1	0.2 ug/L	<0.2	-	-	-
Hexane	1.0 ug/L	<1.0	-	-	-
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	-	-	-
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	-	-	-
Methyl tert-butyl ether	2.0 ug/L	<2.0	-	-	-
Methylene Chloride	5.0 ug/L	<5.0	-	-	-
Styrene	0.5 ug/L	<0.5	-	-	-
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	-	-	-
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	-	-	-
Tetrachloroethylene	0.5 ug/L	<0.5	-	-	-
Toluene	0.5 ug/L	<0.5	-	-	-



Report Date: 11-Sep-2020 Order Date: 4-Sep-2020

	Client ID	DUP	-	_	_
	Sample Date:	03-Sep-20 09:00	-	-	-
	Sample ID:	2036662-05	-	-	-
	MDL/Units	Water	-	-	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	-	-	-
1,1,2-Trichloroethane	0.5 ug/L	<0.5	-	-	-
Trichloroethylene	0.5 ug/L	<0.5	-	-	-
Trichlorofluoromethane	1.0 ug/L	<1.0	-	-	-
Vinyl chloride	0.5 ug/L	<0.5	-	-	-
m,p-Xylenes	0.5 ug/L	<0.5	-	-	-
o-Xylene	0.5 ug/L	<0.5	-	-	-
Xylenes, total	0.5 ug/L	<0.5	-	-	-
4-Bromofluorobenzene	Surrogate	104%	-	-	-
Dibromofluoromethane	Surrogate	98.2%	-	-	-
Toluene-d8	Surrogate	104%	-	-	-



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Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 30732

#### Method Quality Control: Blank

Report Date: 11-Sep-2020

Order Date: 4-Sep-2020

Project Description: PE4641

		Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
F2 PHCs (C10-C16)	ND	100	ug/L						
F3 PHCs (C16-C34)	ND	100	ug/L						
F4 PHCs (C34-C50)	ND	100	ug/L						
Volatiles									
Acetone	ND	5.0	ug/L						
Benzene	ND	0.5	ug/L						
Bromodichloromethane	ND	0.5	ug/L						
Bromoform	ND	0.5	ug/L						
Bromomethane	ND	0.5	ug/L						
Carbon Tetrachloride	ND	0.2	ug/L						
Chlorobenzene	ND	0.5	ug/L						
Chloroform	ND	0.5	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dichlorobenzene	ND	0.5	ug/L						
1,3-Dichlorobenzene	ND	0.5	ug/L						
1,4-Dichlorobenzene	ND	0.5	ug/L						
1,1-Dichloroethane	ND	0.5	ug/L						
1,2-Dichloroethane	ND	0.5	ug/L						
1,1-Dichloroethylene	ND	0.5	ug/L						
cis-1,2-Dichloroethylene	ND	0.5	ug/L						
trans-1,2-Dichloroethylene	ND	0.5	ug/L						
1,2-Dichloropropane	ND	0.5	ug/L						
cis-1,3-Dichloropropylene	ND	0.5	ug/L						
trans-1,3-Dichloropropylene	ND	0.5	ug/L						
1,3-Dichloropropene, total	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Ethylene dibromide (dibromoethane, 1,2	ND	0.2	ug/L						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Isobutyl Ketone	ND	5.0	ug/L						
Methyl tert-butyl ether	ND	2.0	ug/L						
Methylene Chloride	ND	5.0	ug/L						
Styrene	ND	0.5	ug/L						
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L						
Tetrachloroethylene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
1,1,1-Trichloroethane	ND	0.5	ug/L						
1,1,2-Trichloroethane	ND	0.5	ug/L						
Trichloroethylene	ND	0.5	ug/L						
Trichlorofluoromethane	ND	1.0	ug/L						
Vinyl chloride	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
Xylenes, total	ND	0.5	ug/L						
Surrogate: 4-Bromofluorobenzene	79.4	0.0	ug/L		99.2	50-140			
Surrogate: Dibromofluoromethane	67.4		ug/L		84.2	50-140			
-			-		04.2 106				
Surrogate: Toluene-d8	84.6		ug/L		106	50-140			



#### Method Quality Control: Duplicate

Report Date: 11-Sep-2020 Order Date: 4-Sep-2020

Project Description: PE4641

Analista		Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L	ND			NC	30	
Volatiles			-						
Acetone	ND	5.0	ug/L	ND			NC	30	
Benzene	ND	0.5	ug/L	ND			NC	30	
Bromodichloromethane	ND	0.5	ug/L	ND			NC	30	
Bromoform	ND	0.5	ug/L	ND			NC	30	
Bromomethane	ND	0.5	ug/L	ND			NC	30	
Carbon Tetrachloride	ND	0.2	ug/L	ND			NC	30	
Chlorobenzene	ND	0.5	ug/L	ND			NC	30	
Chloroform	ND	0.5	ug/L	ND			NC	30	
Dibromochloromethane	ND	0.5	ug/L	ND			NC	30	
Dichlorodifluoromethane	ND	1.0	ug/L	ND			NC	30	
1,2-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,1-Dichloroethane	ND	0.5	ug/L	ND			NC	30	
1,2-Dichloroethane	ND	0.5	ug/L	ND			NC	30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND			NC	30	
cis-1,2-Dichloroethylene	2.01	0.5	ug/L	2.00			0.5	30	
trans-1,2-Dichloroethylene	ND	0.5	ug/L	ND			NC	30	
1,2-Dichloropropane	ND	0.5	ug/L	ND			NC	30	
cis-1,3-Dichloropropylene	ND	0.5	ug/L	ND			NC	30	
trans-1,3-Dichloropropylene	ND	0.5	ug/L	ND			NC	30	
Ethylbenzene	ND	0.5	ug/L	ND			NC	30	
Ethylene dibromide (dibromoethane, 1,2	ND	0.2	ug/L	ND			NC	30	
Hexane	ND	1.0	ug/L	ND			NC	30	
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L	ND			NC	30	
Methyl Isobutyl Ketone	ND	5.0	ug/L	ND			NC	30	
Methyl tert-butyl ether	ND	2.0	ug/L	ND			NC	30	
Methylene Chloride	ND	5.0	ug/L	ND			NC	30 30	
Styrene 1,1,1,2-Tetrachloroethane	ND ND	0.5 0.5	ug/L	ND ND			NC NC	30 30	
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L	ND			NC	30	
Tetrachloroethylene	ND	0.5	ug/L ug/L	ND			NC	30	
Toluene	ND	0.5	ug/L	ND			NC	30	
1,1,1-Trichloroethane	ND	0.5	ug/L	ND			NC	30	
1,1,2-Trichloroethane	ND	0.5	ug/L	ND			NC	30	
Trichloroethylene	ND	0.5	ug/L	ND			NC	30	
Trichlorofluoromethane	ND	1.0	ug/L	ND			NC	30	
Vinyl chloride	ND	0.5	ug/L	ND			NC	30	
m,p-Xylenes	ND	0.5	ug/L	ND			NC	30	
o-Xylene	ND	0.5	ug/L	ND			NC	30	
Surrogate: 4-Bromofluorobenzene	81.1		ug/L		101	50-140			
Surrogate: Dibromofluoromethane	75.8		ug/L		94.8	50-140			
Surrogate: Toluene-d8	83.6		ug/L		104	50-140			
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#### Method Quality Control: Spike

Report Date: 11-Sep-2020 Order Date: 4-Sep-2020

Project Description: PE4641

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1780	25	ug/L	ND	89.1	68-117			
F2 PHCs (C10-C16)	1660	100	ug/L	ND	104	60-140			
F3 PHCs (C16-C34)	4100	100	ug/L	ND	105	60-140			
F4 PHCs (C34-C50)	2700	100	ug/L	ND	109	60-140			
Volatiles									
Acetone	91.0	5.0	ug/L	ND	91.0	50-140			
Benzene	32.3	0.5	ug/L	ND	80.7	60-130			
Bromodichloromethane	28.2	0.5	ug/L	ND	70.6	60-130			
Bromoform	29.1	0.5	ug/L	ND	72.8	60-130			
Bromomethane	31.8	0.5	ug/L	ND	79.5	50-140			
Carbon Tetrachloride	30.4	0.2	ug/L	ND	76.0	60-130			
Chlorobenzene	34.2	0.5	ug/L	ND	85.5	60-130			
Chloroform	31.6	0.5	ug/L	ND	79.0	60-130			
Dibromochloromethane	30.6	0.5	ug/L	ND	76.5	60-130			
Dichlorodifluoromethane	30.4	1.0	ug/L	ND	76.0	50-140			
1,2-Dichlorobenzene	33.3	0.5	ug/L	ND	83.2	60-130			
1,3-Dichlorobenzene	34.9	0.5	ug/L	ND	87.2	60-130			
1,4-Dichlorobenzene	34.4	0.5	ug/L	ND	85.9	60-130			
1,1-Dichloroethane	32.4	0.5	ug/L	ND	80.9	60-130			
1,2-Dichloroethane	30.6	0.5	ug/L	ND	76.5	60-130			
1,1-Dichloroethylene	31.0	0.5	ug/L	ND	77.4	60-130			
cis-1,2-Dichloroethylene	34.4	0.5	ug/L	ND	86.1	60-130			
trans-1,2-Dichloroethylene	31.5	0.5	ug/L	ND	78.8	60-130			
1,2-Dichloropropane	31.6	0.5	ug/L	ND	79.1	60-130			
cis-1,3-Dichloropropylene	27.8	0.5	ug/L	ND	69.6	60-130			
trans-1,3-Dichloropropylene	27.2	0.5	ug/L	ND	68.0	60-130			
Ethylbenzene	34.2	0.5	ug/L	ND	85.6	60-130			
Ethylene dibromide (dibromoethane, 1,2-	28.8	0.2	ug/L	ND	71.9	60-130			
Hexane	39.8	1.0	ug/L	ND	99.4	60-130			
Methyl Ethyl Ketone (2-Butanone)	68.2	5.0	ug/L	ND	68.2	50-140			
Methyl Isobutyl Ketone	69.5	5.0	ug/L	ND	69.5	50-140			
Methyl tert-butyl ether	72.7	2.0	ug/L	ND	72.7	50-140			
Methylene Chloride	34.0	5.0	ug/L	ND	85.0	60-130			
Styrene	32.5	0.5	ug/L	ND	81.3	60-130			
1,1,1,2-Tetrachloroethane	33.0	0.5	ug/L	ND	82.6	60-130			
1,1,2,2-Tetrachloroethane	28.6	0.5	ug/L	ND	71.6	60-130			
Tetrachloroethylene	33.2	0.5	ug/L	ND	82.9	60-130			
Toluene	33.8	0.5	ug/L	ND	84.6	60-130			
1,1,1-Trichloroethane	30.1	0.5	ug/L	ND	75.3	60-130			
1,1,2-Trichloroethane	28.4	0.5	ug/L	ND	71.0	60-130			
Trichloroethylene	31.3	0.5	ug/L	ND	78.3	60-130			
Trichlorofluoromethane	32.8	1.0	ug/L	ND	82.1	60-130			
Vinyl chloride	30.8	0.5	ug/L	ND	77.1	50-140			
m,p-Xylenes	69.4	0.5	ug/L	ND	86.8	60-130			
o-Xylene	34.6	0.5	ug/L	ND	86.4	60-130			
Surrogate: 4-Bromofluorobenzene	82.6		ug/L		103	50-140			
Surrogate: Dibromofluoromethane	77.0		ug/L		96.3	50-140			
Surrogate: Toluene-d8	83.6		ug/L		104	50-140			



#### Qualifier Notes:

Login Qualifiers :

Container(s) - Bottle and COC sample ID don't match - reads GW2 instead of GW1 Applies to samples: BH8B-GW1

#### Sample Data Revisions

None

#### Work Order Revisions / Comments:

None

#### **Other Report Notes:**

n/a: not applicable ND: Not Detected MDL: Method Detection Limit Source Result: Data used as source for matrix and duplicate samples %REC: Percent recovery. RPD: Relative percent difference. NC: Not Calculated

#### CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.

- F1 range corrected for BTEX.

- F2 to F3 ranges corrected for appropriate PAHs where available.

- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.

- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.

- When reported, data for F4G has been processed using a silica gel cleanup.

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Sample ID/Location Name		Matrix	Air	to #	Date	Time	PHCs	VOCS	PAHs	Metals	ВН	Ş	B (H							
1 BH3-GW3		Gw	>	2	SEPT 03/20			$\setminus$												
2 BH8B-GWI		GW	$\langle \rangle$	Z	SEPT 03/20			$\mathbf{V}$												
3 BH11-20-GWTL		GW	1	3	SEPT 03/20		$\mathbf{k}$	$\overline{}$					_					-		
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