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## **Phase II – Environmental Site Assessment**

Trail's Edge: Phase 4 (South) Southern Parcel (Commercial Zone) Part of 2284 Mer Bleue Road Ottawa, Ontario

**Prepared For** 

**Richcraft Group of Companies** 

January 8, 2021

Report: PE4999-1

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 Ottawa
 Kingston
 North Bay

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# **EXECUTIVE SUMMARY**

## Assessment

A Phase II ESA was conducted for a parcel of land (part of 2284 Mer Bleue Road) situated within the proposed Trail's Edge: Phase 4 (South) residential subdivision development, 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 for this assessment was conducted on September 29, October 19, and November 6, 2020. The field program consisted of drilling three (3) boreholes (BH4-20, BH5-20, and BH6-20), all of which were instrumented with groundwater monitoring wells, as well as excavating five (5) test pits (TP10-TP14). The boreholes were advanced to depths ranging from approximately 4.42 m to 5.94 m below ground surface and terminated within a layer of saturated native silty clay. The test pits were advanced to depths ranging from approximately 1.28 m to 2.61 m below ground surface and terminated within the underlying native soils.

Site soils generally consist of fill material (brown silty sand with crushed stone), underlain by stiff brown silty clay over top of soft grey silty clay. Bedrock was not encountered in any of the borehole or test pit locations.

Three (3) soil samples, recovered from BH4-20, BH5-20, and BH6-20, were submitted for laboratory analysis of: BTEX and PHCs (F<sub>1</sub>-F<sub>4</sub>). An additional five (5) soil samples, recovered from test pits TP10-TP14, were submitted for laboratory analysis of either: PHCs (F<sub>2</sub>-F<sub>4</sub>), PAHs, and/or metals. According to the analytical test results, the concentration of benzo[a]pyrene in soil sample TP14-G1 is in excess of the MECP Table 2 commercial standards.

Three (3) groundwater samples were recovered from the monitoring wells installed in BH4-20, BH5-20, and BH6-20 and submitted for laboratory analysis of VOCs and PHCs (F<sub>1</sub>-F<sub>4</sub>). According to the analytical test results, all detected parameter concentrations in the groundwater samples analyzed are in compliance with the selected MECP Table 2 commercial standards.

## Recommendations

PAH impacted soil/fill material was identified within the vicinity of TP14, located in the southwestern portion of the Phase II study area, requiring some remedial work. It is our understanding that the subject site is to be developed for commercial purposes in conjunction with the neighbouring residential subdivision. 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.

While in compliance with the site-specific standards, it should be noted that the concentration of PHCs and PAHs within the vicinity of TP12 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.

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

If the groundwater monitoring wells installed in BH4-20, BH5-20, and BH6-20 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 Richcraft Group of Companies (Richcraft), Paterson Group (Paterson) conducted a Phase II – Environmental Site Assessment (Phase II ESA) for a parcel of land situated within the proposed Trail's Edge: Phase 4 (South) residential subdivision development, in the City of Ottawa, Ontario. The purpose of this Phase II ESA has been to address the areas of potential environmental concern (APECs) identified on the subject site as a result the findings of the Phase I ESA Update, conducted by Paterson in September 2020.

## 1.1 Subject Site Information

Address:	Part of 2284 Mer Bleue Road, Ottawa, ON
Legal Description:	Block 198, Part of Lots 1, 2, and 3, Concession 3 (Ottawa Front), Part 1 of Registered Plan 4R-30034, Formerly the Geographic Township of Gloucester, in the City of Ottawa.
Location:	The subject site is located on the west side of Mer Bleue Road, south of Brian Coburn Boulevard, in the City of Ottawa, Ontario. Refer to Figure 1 – Key Plan for the site location.
Latitude and Longitude:	45° 26' 39" N, 75° 29' 54" W
Site Description	
Configuration:	Rectangular
Site Area:	8,750 m <sup>2</sup> (approximate)
Zoning:	DR – Development Reserve Zone.
Current Use:	The subject site is currently occupied with an abandoned metal workshop building and storage shed.
Services:	The subject site is located within a municipally serviced area.

## **1.2 Property Ownership**

The subject property is currently owned by Richcraft. Paterson was retained to complete this Phase II ESA by Mr. Patrick Gaudreault of Richcraft, whose offices are located at 2280 St. Laurent Boulevard, Suite 201, Ottawa, Ontario. Mr. Gaudreault can be contacted by telephone at 613-739-1111.

## **1.3 Current and Proposed Future Uses**

The subject site is currently occupied by an abandoned metal workshop building and a storage shed.

It is our understanding that the property is to be redeveloped for commercial purposes as part of the development of the subdivision.

## **1.4 Applicable Site Condition Standard**

The site condition standards for the subject properties were obtained from Table 2 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;
- Potable groundwater conditions;
- Commercial land use.

The commercial standards were selected based on the future land use of the subject sites. Grain size analysis was not conducted as part of this assessment. 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 located on the west side of Mer Bleue Road, south of Brian Coburn Boulevard, in the City of Ottawa, Ontario. The subject site is currently occupied with an abandoned metal workshop building and storage shed. The remainder of the subject site consists primarily of vacant grassland, apart from some gravel surfaced areas in the vicinity of the metal workshop building.

The subject site is at-grade with the adjacent roads as well as the neighbouring properties. The site topography is relatively flat, whereas the regional topography slopes very gently down towards the south, in the general direction of Mer Bleue Bog. Water drainage on the subject site occurs primarily via infiltration throughout the properties, as well as via surface run-off towards drainage ditches present along the adjacent roads.

# 3.0 SCOPE OF INVESTIGATION

## 3.1 Overview of Site Investigation

The subsurface investigation for this assessment was conducted on September 29, October 19, and November 6, 2020. The field program consisted of drilling three (3) boreholes (BH4-20, BH5-20, and BH6-20), all of which were instrumented with groundwater monitoring wells, as well as excavating five (5) test pits (TP10-TP14). The boreholes were advanced to depths ranging from approximately 4.42 m to 5.94 m below ground surface and terminated within a layer of saturated native silty clay. The test pits were advanced to depths ranging from approximately 1.28 m to 2.61 m below ground surface and terminated within the underlying native soils.

## 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 include the following:

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

## 3.3 Phase I ESA Conceptual Site Model

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

The subject site is currently occupied with a two (2) storey, slab-on-grade style metal workshop building (currently abandoned) as well as a storage shed.

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

No areas of natural and scientific interest are known to exist within the Phase I study area. The nearest named water body with respect to the subject site is Mer Bleue Bog, located approximately 2.25 km to the south.

#### **Geological and Hydrogeological Setting**

Based on the available mapping information, the bedrock within the area of the subject sites consist of interbedded limestone and shale of the Lindsay Formation, whereas the surficial geology consists of offshore marine deposits (clay and silt) with an overburden ranging from approximately 15 m to 50 m in thickness. Based on the regional topography, the groundwater is interpreted to be moving in a southerly direction towards Mer Bleue Bog.

#### Neighbouring Land Use

The neighbouring lands within the Phase I study area consist of a combination of residential and commercial properties or vacant land.

#### **Drinking Water Wells**

Based on the available MECP water well records, it is likely that some of the residential properties adjacent to Mer Bleue Road may still utilize private drinking water wells.

# Potentially Contaminating Activities and Areas of Potential Environmental Concern

Based on the findings of the Phase I ESA Update, two (2) potentially contaminating activities (PCAs), resulting in areas of potential environmental concern (APECs), were identified as pertaining to the subject sites:

- □ The presence and historical operation of an on-site metal workshop building, located within the central portion of the Phase II study area;
- □ The presence of fill material of unknown quality, located within the vicinity of the on-site metal workshop building;

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 aforementioned APECs are considered to be:

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

These CPCs have the potential to be present in the soil matrix and/or the groundwater situated beneath the subject sites.

#### 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.

# 4.0 INVESTIGATION METHOD

## 4.1 Subsurface Investigation

The subsurface investigation for this assessment was conducted on September 29, October 19, and November 6, 2020. The field program consisted of drilling three (3) boreholes (BH4-20, BH5-20, and BH6-20), all of which were instrumented with groundwater monitoring wells, as well as excavating five (5) test pits (TP10-TP14). The boreholes were advanced to depths ranging from approximately 4.42 m to 5.94 m below ground surface and terminated within a layer of saturated native silty clay. The test pits were advanced to depths ranging from approximately 1.28 m to 2.61 m below ground surface and terminated within the underlying native soils.

Under the full-time supervision of Paterson personnel, the boreholes were drilled using a track-mounted drill rig provided by George Downing Estate Drilling of Hawkesbury, Ontario; whereas the test pits were excavated using a backhoe provided by Quast Excavating of Ottawa, Ontario.

The locations of the boreholes and test pits are illustrated on Drawing PE4999-3 – 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.

Twenty-two (22) soil samples were obtained from the boreholes by means of auger and split spoon sampling, with samples taken at approximate 0.76 m intervals. The depths at which auger and split spoon samples were obtained from the boreholes are shown as "**AU**" and "**SS**", respectively, on the Soil Profile and Test Data Sheets, appended to this report.

Twelve (12) soil samples were obtained from the test pits by means of grab sampling. The depths at which grab samples were obtained from the test pits are shown as "G", on the Soil Profile and Test Data Sheets, appended to this report.

Site soils generally consist of fill material (brown silty sand), underlain by stiff brown silty clay over top of soft grey silty clay. Bedrock was not encountered in any of the boreholes or test pits at the time of the field program.

## 4.3 Field Screening Measurements

All soil samples collected were subjected to a preliminary screening procedure, which included visual screening for colour and evidence of metals, as well as soil vapour screening with a Photo Ionization Detector.

The recovered soil samples were placed immediately into airtight plastic bags with nominal headspace. All lumps of soil inside the bags were broken by hand, and the soil was allowed to come to room temperature prior to conducting the vapour survey, ensuring consistency of readings between samples. To measure the soil vapours, the analyser probe was inserted into the nominal headspace above the sample. The sample was then agitated and manipulated gently by hand as the measurement was taken. The peak reading registered within the first 15 seconds was recorded as the vapour measurement. The parts per million (ppm) scale was used to measure concentrations of organic vapours.

Samples with the highest vapour readings for a given borehole were generally selected as candidates for laboratory analysis. The results of the vapour survey are presented on the Soil Profile and Test Data Sheets, appended to this report.

## 4.4 Groundwater Monitoring Well Installation

Three (3) groundwater monitoring wells were installed on the subject sites as part of this assessment. 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.

Table 1 Monitoring	g Well Construc	ction Deta	ails			
Well ID	Ground Surface Elevation (m ASL)	Total Depth (m BGS)	Screened Interval (m BGS)	Sand Pack (m BGS)	Bentonite Seal (m BGS)	Casing Type
BH4-20	88.24	5.94	2.94-5.94	2.44-5.94	0.10-2.44	Stick-Up
BH5-20	87.81	5.94	2.94-5.94	2.44-5.94	0.10-2.44	Flushmount
BH6-20	87.81	4.42	1.42-4.42	0.74-1.42	0.43-0.74	Stick-Up

## 4.5 Field Measurement of Water Quality Parameters

Groundwater sampling was conducted at BH4-20, BH5-20, and BH6-20 on October 2, and November 12, 2020. No water quality parameters were measured in the field at that time.

## 4.6 Groundwater Sampling

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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. Standing water was purged from each monitoring well prior to the recovery of the groundwater samples using dedicated sampling equipment. The samples were then stored in coolers to reduce possible analyte volatilization during their transportation. Further details of our standard operating procedure for groundwater sampling are provided in the Sampling and Analysis Plan, appended to this report.

### 4.7 Residue Management

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

## 4.8 Analytical Testing

The following soil and groundwater samples were submitted for laboratory analysis:

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Table 2									
Testing	Parameters for	or Su	bmit	ted S	oil S	amp	les		
	Parameters Analyzed								
Sample ID	Sample Depth & Stratigraphic Unit	ВТЕХ	PHCs F <sub>1</sub> -F <sub>4</sub>	PHCs F <sub>2</sub> -F <sub>4</sub>	PAHs	Metals <sup>1</sup>	Rationale		
BH4-20- SS5	3.05 – 3.66 m Grey Silty Clay	х	х						
BH5-20- SS4	2.29 – 2.90 m Grey Silty Clay	х	х				To assess for potential impacts resulting from a former metal workshop.		
BH6-20- SS3	1.52 – 2.13 m Grey Silty Clay	х	х						
TP10-G1	0.10 – 0.30 m Fill Material			х	х	х			
TP11-G1	0.30 – 0.50 m Fill Material			х	х	х			
TP12-G1	0.10 – 0.20 m Fill Material			х	х	х	To assess for potential impace resulting from the presence of material of unknown quality.		
TP13-G1	0.10 – 0.20 m Fill Material			х	х	х			
TP14-G1	0.10 – 0.20 m Fill Material			х	х	х			
1 – Including	Mercury and Hexava	lent Chr	omium						

Testing F			s Analyzed	•			
Sample ID	Screened Interval & Stratigraphic Unit	VOCs	PHCs F <sub>1</sub> -F <sub>4</sub>	Rationale			
BH4-20- GW1	2.94 – 5.94 m Grey Silty Clay	Х	Х				
BH5-20- GW1	2.94 – 5.94 m Grey Silty Clay	Х	Х	To assess for potential impacts resulting from			
BH6-20- GW1	1.42 – 4.42 m Grey Silty Clay	Х	Х	a former metal workshop.			
DUP1	2.94 – 5.94 m Grey Silty Clay	Х					
1 – Including N	Aercury and Hexava	ent Chromium					

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 Laboratory Accreditation (SCC/CALA) and is accredited and certified by the SCC/CALA for specific tests registered with the association.

## 4.9 Elevation Surveying

The ground surface elevations at each borehole and test pit location were surveyed using a GPS device by Paterson personnel and referenced to a geodetic datum.

## 4.10 Quality Assurance and Quality Control Measures

A summary of the quality assurance and quality control (QA/QC) measures, undertaken as part of this assessment, is provided in the Sampling and Analysis Plan in Appendix 1.

## 5.0 REVIEW AND EVALUATION

## 5.1 Geology

Generally, the subsurface profile encountered at the borehole locations consists of fill material, underlain by stiff brown silty clay over top of soft grey silty clay. The surficial fill material was observed to consist of brown silty sand with crushed stone and was noted to extend to depths ranging from approximately 0.23 m to 1.07 m below ground surface. No unusual visual or olfactory observations, as well as any deleterious substances, were identified within the fill material encountered on-site.

The subsurface profile encountered at the test pit locations consists of fill material underlain by native brown silty clay. The fill material was observed to consist of brown silty clay / silty sand, organics, as well as trace brick and asphalt, and was noted to extend to depths ranging from approximately 0.2 m to 0.8 m below ground surface.

Bedrock was not encountered in any of the boreholes or test pits.

Site geology details are provided in the Soil Profile and Test Data Sheets in Appendix 1.

## 5.2 Groundwater Elevations, Flow Direction, and Hydraulic Gradient

Groundwater levels were measured using an electronic water level meter at the monitoring wells installed in BH4-20, BH5-20, and BH6-20 on November 12, 2020. The groundwater levels are summarized below in Table 4.

Table 4 Groundwater Level Measurements								
Borehole Location	Ground Surface Elevation (m)	Water Level Depth (m below grade)	Water Level Elevation (m ASL)	Date of Measurement				
BH4-20	88.24	1.04	87.20					
BH5-20	87.81	0.37	87.44	November 12, 2020				
BH6-20	87.81	0.46	87.35					

The groundwater at the subject site was encountered within the native grey silty clay, at depths ranging from approximately 0.37 m to 1.04 m below the existing ground surface. No unusual visual or olfactory observations were noted in the groundwater samples recovered from the boreholes.

Using the groundwater elevations recorded during the November 12, 2020 sampling event, groundwater contour mapping was completed as part of this assessment. According to the mapped contour data, illustrated on Drawing PE4999-3 – Test Hole Location Plan in the appendix, the groundwater flow within the Phase II study area is interpreted to be in a northerly direction. A horizontal hydraulic gradient of approximately 0.007 m/m was also calculated as part of this assessment.

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

## 5.3 Fine/Coarse Soil Texture

Grain size analysis was not completed as part of this investigation. Coarse grained soil standards were chosen as a conservative approach.

### 5.4 Field Screening

Field screening of the soil samples collected during the drilling program resulted in organic vapour readings ranging from 0.1 ppm to 8.2 ppm. The organic vapour readings obtained from the field screening indicate that there is a negligible potential for the presence of volatile substances. Field screening results of each individual soil sample are provided on the Soil Profile and Test Data Sheets appended to this report.

## 5.5 Soil Quality

Three (3) soil samples, recovered from BH4-20, BH5-20, and BH6-20, were submitted for laboratory analysis of: BTEX and PHCs ( $F_1$ - $F_4$ ).

Five (5) soil samples, recovered from test pits TP10-TP14, were submitted for laboratory analysis of either: PHCs (F<sub>2</sub>-F<sub>4</sub>), PAHs, and/or metals.

The results of the analytical testing are presented below in Tables 5 to 8, as well as on the laboratory certificates of analysis included in Appendix 1.

Analytical 1 (BTEX & PH		4)	bil Samples (μg		MECP	MECP
			9, 2020	Nov. 6, 2020	Table 1	Table 2
Parameter	MDL (µg/g)	BH4-20-SS5	BH5-20-SS4	BH6-20-SS3	Background Soil Standards (µg/g)	Commercial Soil Standards (µg/g)
Benzene	0.05	nd	nd	nd	0.02	0.21
Ethylbenzene	0.05	nd	nd	nd	0.05	1.1
Toluene	0.05	nd	nd	nd	0.2	2.3
Xylenes	0.05	nd	nd	nd	0.05	3.1
PHCs F1	7	nd	nd	nd	25	55
PHCs F <sub>2</sub>	4	nd	nd	nd	10	98
PHCs F₃	8	nd	nd	nd	240	300
PHCs F <sub>4</sub>	6	nd	nd	nd	120	2,800
<ul> <li>nd – not</li> <li><u>Underlin</u></li> </ul>	ied – Value	ection Limit pove the MDL exceeds MECP Ta ed – value exceeds		standards		

No BTEX or PHC parameters were detected in the soil samples analyzed. The results are in compliance with the selected MECP Table 2 commercial standards as well as the MECP Table 1 standards.

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				Samples (	MECP Table 1	MECP Table 2		
Parameter	MDL (µg/g)	TP10- G1	TP11- G1	ct. 19, 202 TP12- G1	TP13- G1	TP14- G1	Background Soil Standards (µg/g)	Commercial Soil Standards (μg/g)
PHCs F <sub>2</sub>	4	nd	nd	nd	nd	nd	10	98
PHCs F <sub>3</sub>	8	nd	nd	40	nd	60	240	300
PHCs F <sub>4</sub>	6	nd	nd	122	nd	188	120	2,800
PHCs F <sub>4G</sub>	50	nt	nt	nt	nt	722	120	2,800

All detected PHC parameter concentrations in the soil samples analyzed are in compliance with the selected MECP Table 2 commercial standards.

The concentration of PHC  $F_4$  in sample TP12-G1, as well as the concentrations of PHC  $F_4$  and  $F_{4G}$  in sample TP14-G1 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.

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Table 7
Analytical Test Results – Soil (Test Pits
PAHs

TP10 -G1 nd nd nd nd nd nd nd nd nd nd	TP11 -G1 nd nd nd nd nd nd nd nd nd nd	<b>TP12</b> -G1 nd nd 0.05 0.18 0.22 0.14 0.06	<b>TP13</b> -G1 nd 0.02 0.07 0.16 0.16 0.18 0.10 0.09	TP14 -G1 0.04 0.09 0.36 0.41 0.79 0.34 0.44	Background Soil Standards (μg/g) 0.072 0.093 0.16 0.36 0.3 0.47 0.68 0.48	Commercial Soil Standards (μg/g) 7.9 0.15 0.67 0.5 0.3 0.78 6.6 0.78
nd nd nd nd nd nd nd nd nd nd	nd nd nd nd nd nd nd nd nd	nd nd 0.05 0.18 <u>0.50</u> 0.22 0.14	0.02 0.07 0.16 0.16 0.18 0.10 0.09	0.04 0.09 0.36 <b>0.41</b> <b>0.79</b> 0.34	0.093 0.16 0.36 0.3 0.47 0.68	0.15 0.67 0.5 0.3 0.78 6.6
nd nd nd nd nd nd nd nd nd	nd nd nd nd nd nd nd	nd 0.05 0.18 <u>0.50</u> 0.22 0.14	0.07 0.16 0.16 0.18 0.10 0.09	0.09 0.36 <u>0.41</u> <u>0.79</u> 0.34	0.16 0.36 0.3 0.47 0.68	0.67 0.5 0.3 0.78 6.6
nd nd nd nd nd nd nd	nd nd nd nd nd nd	0.05 0.18 <u>0.50</u> 0.22 0.14	0.16 0.16 0.18 0.10 0.09	0.36 0.41 0.79 0.34	0.36 0.3 0.47 0.68	0.5 0.3 0.78 6.6
nd nd nd nd nd nd	nd nd nd nd nd	0.18 0.50 0.22 0.14	0.16 0.18 0.10 0.09	0.41 0.79 0.34	0.3 0.47 0.68	0.3 0.78 6.6
nd nd nd nd nd	nd nd nd nd	0.50 0.22 0.14	0.18 0.10 0.09	0.34	0.47 0.68	0.78 6.6
nd nd nd nd	nd nd nd	0.22	0.10 0.09	0.34	0.68	6.6
nd nd nd	nd nd	0.14	0.09			
nd nd	nd			0.44	0 48	0.70
nd	-	0.06	0 1 0		0.10	0.70
-		0.00	0.16	0.42	2.8	7
	nd	0.04	0.03	0.08	0.1	0.1
nd	nd	0.11	0.35	0.84	0.56	0.69
nd	nd	nd	nd	nd	0.12	62
nd	nd	0.13	0.09	0.29	0.23	0.38
nd	nd	nd	nd	nd	0.59	0.99
nd	nd	nd	nd	nd	0.59	0.99
nd	nd	nd	nd	nd	0.59	0.99
nd	nd	nd	nd	nd	0.09	0.6
nd	nd	0.06	0.17	0.29	0.69	6.2
nd	nd	0.16	0.27	0.67	1	78
	nd nd nd nd	nd nd nd nd nd nd nd nd nd nd	nd nd nd nd nd nd nd nd nd nd nd 0.06 nd nd 0.16	nd         nd         nd         nd           nd         nd         0.06         0.17           nd         nd         0.16         0.27	nd         nd         nd         nd         nd           nd         nd         0.06         0.17         0.29           nd         nd         0.16         0.27         0.67	nd         nd         nd         nd         nd         0.59           nd         nd         nd         nd         nd         0.59           nd         nd         nd         nd         nd         0.59           nd         nd         nd         nd         nd         0.09           nd         nd         0.06         0.17         0.29         0.69           nd         nd         0.16         0.27         0.67         1

Bold and Underlined – value exceeds selected MECP standards

The concentrations of benzo[a]pyrene, benzo[b]fluoranthene, and fluoranthene detected in soil sample TP14-G1 is in excess of the selected MECP Table 2 commercial standards. All remaining PAH parameter concentrations detected in the soil samples analyzed are in compliance with the selected MECP Table 2 commercial standards.

The concentration of benzo[b]fluoranthene detected in soil sample TP12-G1, as well as the concentration of indeno[1,2,3-cd]pyrene detected in soil sample TP14-G1 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 these soils are ever to be removed from the property, they should be classified as contaminated and disposed of at an approved waste disposal site.

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				Samples ( ober 19, 2	MECP Table 1	MECP Table 2		
Parameter	MDL (µg/g)	TP10- G1	TP11- G1	TP12- G1	TP13- G1	TP14- G1	Background Soil Standards (µg/g)	Commercial Soil Standards (µg/g)
Antimony	1.0	nd	nd	nd	nd	nd	1.3	7.5
Arsenic	1.0	1.8	3.3	4.4	3.0	3.1	18	18
Barium	1.0	55.3	167	157	116	138	220	390
Beryllium	0.5	nd	0.6	0.6	nd	nd	2.5	4
Boron	5.0	nd	nd	6.2	6.2	5.4	36	120
Cadmium	0.5	nd	nd	nd	nd	nd	1.2	1.2
Chromium	5.0	15.4	69.2	69.1	48.2	53.5	70	160
Chromium (VI)	0.2	nd	nd	nd	nd	nd	0.66	8
Cobalt	1.0	3.6	13.7	12.0	9.0	10.9	21	22
Copper	5.0	5.4	26.8	22.8	17.2	23.3	92	140
Lead	1.0	4.0	7.8	12.8	11.6	15.1	120	120
Mercury	0.05	nd	nd	nd	nd	nd	0.27	0.27
Molybdenum	1.0	nd	nd	1.2	nd	nd	2	6.9
Nickel	5.0	9.3	38.5	35.5	25.0	31.1	82	100
Selenium	1.0	nd	nd	nd	nd	nd	1.5	2.4
Silver	0.3	nd	nd	nd	nd	nd	0.5	20
Thallium	1.0	nd	nd	nd	nd	nd	1	1
Uranium	1.0	nd	nd	2.4	1.3	nd	2.5	23
Vanadium	10.0	15.9	54.5	56.1	39.6	51.3	86	86
Zinc	20.0	nd	65.9	84.1	53.7	65.1	290	340

Bold and Underlined – value exceeds selected MECP standards

All detected metal concentrations in the soil samples analysed are in compliance with the selected MECP Table 2 commercial standards as well as the MECP Table 1 standards.

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Parameter	Maximum Concentration (µg/g)	Sample ID	Depth Interval (m BGS)
PHCs F <sub>3</sub>	60	TP14-G1	0.1 – 0.2 m
PHCs F <sub>4</sub>	<u>188</u>	TP14-G1	0.1 – 0.2 m
PHCs F <sub>4G</sub>	<u>722</u>	TP14-G1	0.1 – 0.2 m
Acenaphthylene	0.04	TP14-G1	0.1 – 0.2 m
Anthracene	0.09	TP14-G1	0.1 – 0.2 m
Benzo[a]anthracene	0.36	TP14-G1	0.1 – 0.2 m
Benzo[a]pyrene	<u>0.41</u>	TP14-G1	0.1 – 0.2 m
Benzo[b]fluoranthene	0.79	TP14-G1	0.1 – 0.2 m
Benzo[g,h,i]perylene	0.34	TP14-G1	0.1 – 0.2 m
Benzo[k]fluoranthene	0.44	TP14-G1	0.1 – 0.2 m
Chrysene	0.42	TP14-G1	0.1 – 0.2 m
Dibenzo[a,h]anthracene	0.08	TP14-G1	0.1 – 0.2 m
Fluoranthene	<u>0.84</u>	TP14-G1	0.1 – 0.2 m
Indeno[1,2,3-cd]pyrene	0.29	TP14-G1	0.1 – 0.2 m
Phenanthrene	0.29	TP14-G1	0.1 – 0.2 m
Pyrene	0.67	TP14-G1	0.1 – 0.2 m
Arsenic	4.4	TP12-G1	0.1 – 0.2 m
Barium	167	TP11-G1	0.3 – 0.5 m
Beryllium	0.6	TP11-G1 / TP12-G1	0.3 – 0.5 m / 0.1 – 0.2 r
Boron	6.2	TP12-G1 / TP13-G1	0.1 – 0.2 m / 0.1 – 0.2 r
Chromium	69.2	TP11-G1	0.3 – 0.5 m
Cobalt	13.7	TP11-G1	0.3 – 0.5 m
Copper	26.8	TP11-G1	0.3 – 0.5 m
Lead	15.1	TP14-G1	0.1 – 0.2 m
Molybdenum	1.2	TP12-G1	0.1 – 0.2 m
Nickel	38.5	TP11-G1	0.3 – 0.5 m
Uranium	2.4	TP12-G1	0.1 – 0.2 m
Vanadium	56.1	TP12-G1	0.1 – 0.2 m
Zinc	84.1	TP12-G1	0.1 – 0.2 m

All other parameter concentrations analyzed were below the laboratory detection limits. The laboratory certificates of analysis are provided in Appendix 1.

## 5.6 Groundwater Quality

Groundwater samples were recovered from the monitoring wells installed in BH4-20, BH5-20, and BH6-20 and submitted for laboratory analysis of: PHCs (F<sub>1</sub>-F<sub>4</sub>) and VOCs. The results of the analytical testing are presented below in Tables 10 and 11, as well as on the laboratory certificates of analysis included in Appendix 1.

		Grou	Indwater Samples (	(µg/L)	MECP Table 2
Parameter	MDL	Oct. 2	2, 2020	Nov. 12, 2020	Commercial Groundwater
Farameter	(µg/L)	BH4-20-GW1	BH5-20-GW1	BH6-20-GW1	Standards (µg/L)
PHC F1	25	nd	nd	nd	750
PHC F <sub>2</sub>	100	nd	nd	nd	150
PHC F <sub>3</sub>	100	nd	nd	nd	500
PHC F <sub>4</sub>	100	nd	nd	nd	500

No PHC parameters were detected in the groundwater samples analyzed. The results are in compliance with the selected MECP Table 2 commercial standards.

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Trail's Edge: Phase 4 (South) - Southern Parcel (Commercial Zone) Ottawa, Ontario

Analytical Test Results – Groundwate	Table 11		
	Analytical Test R	esults	- Groundwater
VOCs	VOCs		

		Grour	ndwater Samples	(ug/L)	MECP
Parameter	MDL (µg/L)	October	r 2, 2020	November 12, 2020	Table 2 Residential Groundwater Standards (µg/L)
	(µg/⊏)	BH4-20-GW1	BH5-20-GW1	BH6-20-GW1	
Acetone	5.0	nd	nd	nd	2,700
Benzene	0.5	nd	nd	nd	5
Bromodichloromethane	0.5	nd	nd	nd	16
Bromoform	0.5	nd	nd	nd	25
Bromomethane	0.5	nd	nd	nd	0.89
Carbon Tetrachloride	0.2	nd	nd	nd	0.79
Chlorobenzene	0.5	nd	nd	nd	30
Chloroform	0.5	nd	nd	nd	2.4
Dibromochloromethane	0.5	nd	nd	nd	25
Dichlorodifluoromethane	1.0	nd	nd	nd	590
1,2-Dichlorobenzene	0.5	nd	nd	nd	3
1,3-Dichlorobenzene	0.5	nd	nd	nd	59
1,4-Dichlorobenzene	0.5	nd	nd	nd	1
1,1-Dichloroethane	0.5	nd	nd	nd	5
1,2-Dichloroethane	0.5	nd	nd	nd	1.6
1,1-Dichloroethylene	0.5	nd	nd	nd	1.6
cis-1,2-Dichloroethylene	0.5	nd	nd	nd	1.6
trans-1,2-Dichloroethylene	0.5	nd	nd	nd	1.6
1,2-Dichloropropane	0.5	nd	nd	nd	5
1,3-Dichloropropene	0.5	nd	nd	nd	0.5
Ethylbenzene	0.5	nd	nd	nd	2.4
Ethylene Dibromide	0.2	nd	nd	nd	0.2
Hexane	1.0	nd	nd	nd	51
Methyl Ethyl Ketone	5.0	nd	nd	nd	1,800
Methyl Isobutyl Ketone	5.0	nd	nd	nd	640
Methyl tert-butyl ether	2.0	nd	nd	nd	15
Methylene Chloride	5.0	nd	nd	nd	20
Styrene	0.5	nd	nd	nd	5.4
1,1,1,2-Tetrachloroethane	0.5	nd	nd	nd	1.1
1,1,2,2-Tetrachloroethane	0.5	nd	nd	nd	1
Tetrachloroethylene	0.5	nd	nd	nd	1.6
Toluene	0.5	nd	nd	nd	24
1,1,1-Trichloroethane	0.5	nd	nd	nd	200
1,1,2-Trichloroethane	0.5	nd	nd	nd	4.7
Trichloroethylene	0.5	nd	nd	nd	1.6
Trichlorofluoromethane	1.0	nd	nd	nd	150
Vinvl Chloride	0.5	nd	nd	nd	0.5
Xylenes	0.5	nd	nd	nd	300
Notes: MDL – Method De nd – not detected Bold and Underli	above the	MDL			

No VOC parameters were detected in the groundwater samples analyzed. The results are in compliance with the selected MECP Table 2 commercial standards.

## 5.7 Quality Assurance and Quality Control Results

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.

As per the Sampling and Analysis Plan, a duplicate groundwater sample was obtained from the monitoring well installed in BH4-20 and analyzed for VOC parameters. No parameter concentrations were detected in both the original or the duplicate samples, and as such, the RPD results are considered to be acceptable. 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.

## 5.8 Phase II Conceptual Site Model

The following section has been prepared in accordance with the requirements of O.Reg. 153/04 amended by the Environmental Protection Act. Conclusions and recommendations are discussed in a subsequent section.

### Site Description

# Potentially Contaminating Activity and Areas of Potential Environmental Concern

As described in the Phase I ESA Update 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 sites:

#### Item 30: "Importation of Fill Material of Unknown Quality"

This PCA was identified as a result of the presence of fill material located within the vicinity of the on-site metal workshop.

#### Item 34: "Metal Fabrication"

This PCA was identified as a result of the historical operation of an on-site metal workshop.

#### Contaminants of Potential Concern

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

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

These CPCs have the potential to be present in the soil matrix and/or the groundwater situated beneath the subject sites.

#### Subsurface Structures and Utilities

Underground service locates were completed prior to the subsurface investigation. No underground utilities were identified on the subject site.

### Physical Setting

#### Site Stratigraphy

The stratigraphy of the subject sites generally consists of:

- □ Crushed stone (engineered fill); encountered at ground surface and extending to depths of approximately 0.10 m to 0.80 m below ground surface;
- □ Fill material consisting of brown silty sand; encountered at ground surface and extending to depths of approximately 0.38 m to 1.07 m below ground surface;
- □ Stiff native brown silty clay; encountered at depths ranging from approximately 0.23 m to 1.07 m below ground surface;
- □ Soft native grey silty clay; encountered at a depths ranging from approximately 1.98 m to 3.05 m below ground surface.

The site stratigraphy, from ground surface to the deepest aquifer or aquitard investigated, is provided in the Soil Profile and Test Data Sheets in Appendix 1.

#### Hydrogeological Characteristics

The groundwater beneath the subject site was typically encountered within the native grey silty clay, at depths ranging from approximately 0.37 m to 1.04 m below the existing ground surface. Based on the measured groundwater levels, the groundwater flow direction in the vicinity of the boreholes is interpreted to be in a northerly direction.

#### Approximate Depth to Bedrock

Bedrock was not encountered in any of the boreholes. According to the available mapping information, the bedrock is interpreted to lie at a depth of approximately 15 m to 50 m below ground level.

#### Approximate Depth to Water Table

The depth to the water table is approximately 0.37 m to 1.04 m below the existing ground surface.

#### Sections 41 and 43.1 of Ontario Regulation 153/04

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

Section 43.1 of the Regulation does not apply to the subject site, since the bedrock is not situated at a depth of less than 2 m below ground surface, and thus is not considered to be a shallow soil property.

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

The subject site is currently occupied with a two (2) storey, slab-on-grade style metal workshop building (currently abandoned) as well as a storage shed.

#### Fill Placement

Fill material, consisting of brown silty sand and silty clay was encountered within the test pit locations on the western portion of the study area. Engineered fill consisting of crushed stone was encountered within the borehole locations adjacent to the metal workshop building.

#### **Proposed Buildings and Other Structures**

It is our understanding that the land is to be redeveloped for commercial purposes as part of the development of the surrounding subdivision.

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

No areas of natural and scientific interest are known to exist within the Phase I study area. The nearest named water body with respect to the subject site is Mer Bleue Bog, located approximately 2.25 km to the south.

## **Environmental Condition**

#### Areas Where Contaminants are Present

According to the analytical test results, PAH impacted fill material was identified in TP14, located in the southwestern portion of the Phase II study area.

#### Types of Contaminants

According to the analytical test results, the concentration of PAHs detected in fill sample TP14-G1 is in excess of the MECP Table 2 commercial standards.

#### Contaminated Media

As noted above, the fill within the vicinity of TP14-G1 is in excess of the MECP Table 2 commercial standards.

According to the analytical test results, the groundwater beneath the subject site is not contaminated.

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

TP14 is located within the southwestern portion of the Phase II study area. The PAH contaminants identified in this location are the result of poor-quality fill material placed in this area.

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

As noted above, PAH impacted soil/fill material was identified within the southwestern portion of the subject site, in the vicinity of TP14. Based on their low mobility, it is anticipated that the PAH contaminants are contained within the soil/fill in this area of the subject site.

#### Discharge of Contaminants

The PAH impacted fill in the vicinity of TP14 is considered to have resulted from the importation and placement of poor-quality fill material.

#### **Potential for Vapour Intrusion**

Based on the non-volatile nature of the contaminants, as well as the slab-ongrade nature of the proposed commercial developments, there is no risk for any future vapour intrusion on the subject site.

#### **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.

The downward migration of PAH contaminants in the vicinity of TP14 is not suspected to have occurred based on the nature of the contaminants as well as their low concentration. In addition, any fluctuations in the groundwater level and groundwater flow are not considered to have affected any contaminant distribution at this location due to the depth of the water table well below the shallow fill material.

## 6.0 CONCLUSION

## Assessment

A Phase II ESA was conducted for a parcel of land (part of 2284 Mer Bleue Road) situated within the proposed Trail's Edge: Phase 4 (South) residential subdivision development, 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 for this assessment was conducted on September 29, October 19, and November 6, 2020. The field program consisted of drilling three (3) boreholes (BH4-20, BH5-20, and BH6-20), all of which were instrumented with groundwater monitoring wells, as well as excavating five (5) test pits (TP10-TP14). The boreholes were advanced to depths ranging from approximately 4.42 m to 5.94 m below ground surface and terminated within a layer of saturated native silty clay. The test pits were advanced to depths ranging from approximately 1.28 m to 2.61 m below ground surface and terminated within the underlying native soils.

Site soils generally consist of fill material (brown silty sand with crushed stone), underlain by stiff brown silty clay over top of soft grey silty clay. Bedrock was not encountered in any of the borehole or test pit locations.

Three (3) soil samples, recovered from BH4-20, BH5-20, and BH6-20, were submitted for laboratory analysis of: BTEX and PHCs ( $F_1$ - $F_4$ ). An additional five (5) soil samples, recovered from test pits TP10-TP14, were submitted for laboratory analysis of either: PHCs ( $F_2$ - $F_4$ ), PAHs, and/or metals. According to the analytical test results, the concentration of benzo[a]pyrene in soil sample TP14-G1 is in excess of the MECP Table 2 commercial standards.

Three (3) groundwater samples were recovered from the monitoring wells installed in BH4-20, BH5-20, and BH6-20 and submitted for laboratory analysis of VOCs and PHCs ( $F_1$ - $F_4$ ). According to the analytical test results, all detected parameter concentrations in the groundwater samples analyzed are in compliance with the selected MECP Table 2 commercial standards.

## Recommendations

PAH impacted soil/fill material was identified within the vicinity of TP14, located in the southwestern portion of the Phase II study area, requiring some remedial work. It is our understanding that the subject site is to be developed for commercial purposes in conjunction with the neighbouring residential subdivision. 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 fill, the latter of which will require disposal at an approved waste disposal facility.

While in compliance with the site-specific standards, it should be noted that the concentration of PHCs and PAHs within the vicinity of TP12 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.

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

If the groundwater monitoring wells installed in BH4-20, BH5-20, and BH6-20 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 in general accordance with O.Reg. 153/04, as amended, 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 Richcraft Group of Companies. Permission and notification from the Richcraft Group of Companies and Paterson Group will be required prior to the release of 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

## **DRAWING PE4999-3 – TEST HOLE LOCATION PLAN**

DRAWING PE4999-4 – ANALYTICAL TESTING PLAN – SOIL (PAHS)

DRAWING PE4999-4A – CROSS SECTION A-A' – SOIL (PAHS)

DRAWING PE4999-5 – ANALYTICAL TESTING PLAN – SOIL (BTEX, PHCS, METALS)

DRAWING PE4999-5A – CROSS SECTION A-A' – SOIL (BTEX, PHCS, METALS)

DRAWING PE4999-6 – ANALYTICAL TESTING PLAN – GROUNDWATER

DRAWING PE4999-6A – CROSS SECTION A-A' – GROUNDWATER

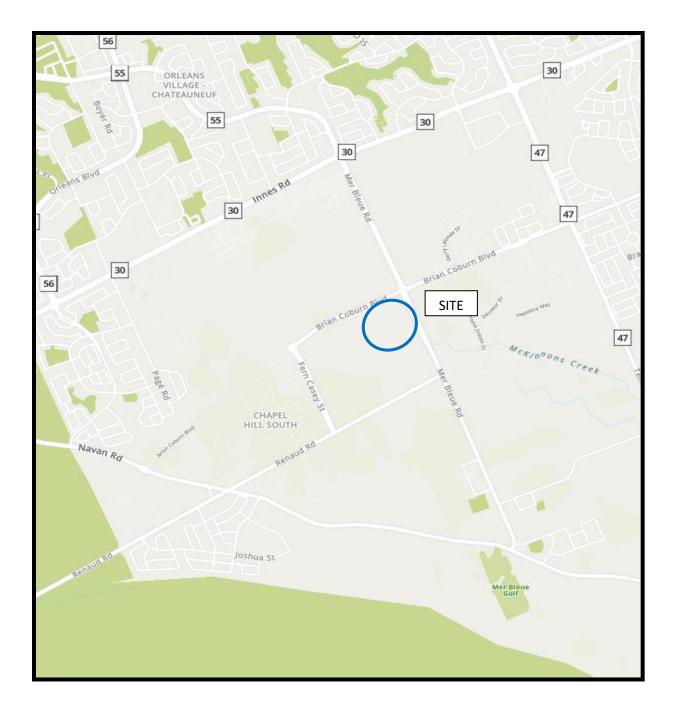
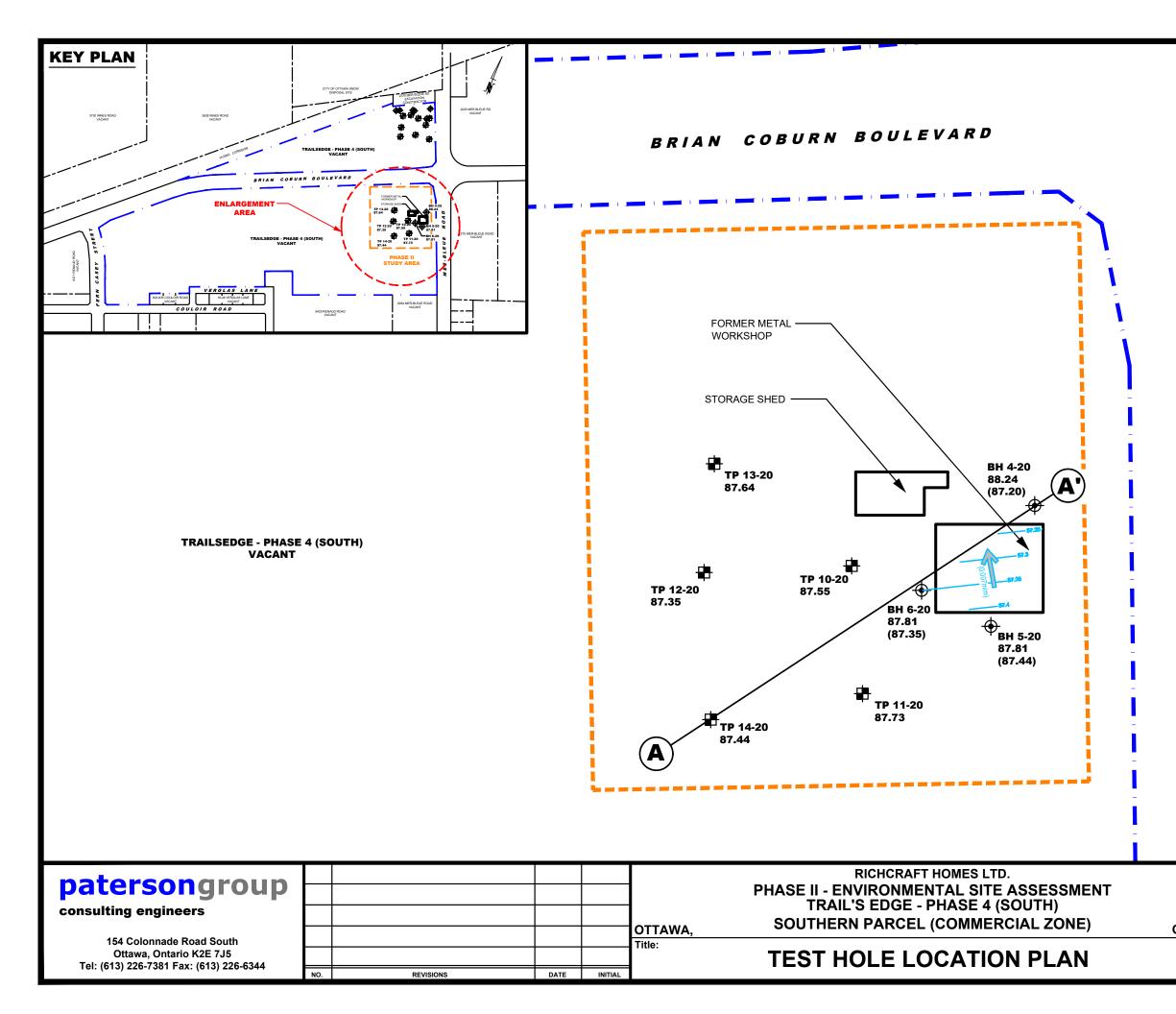
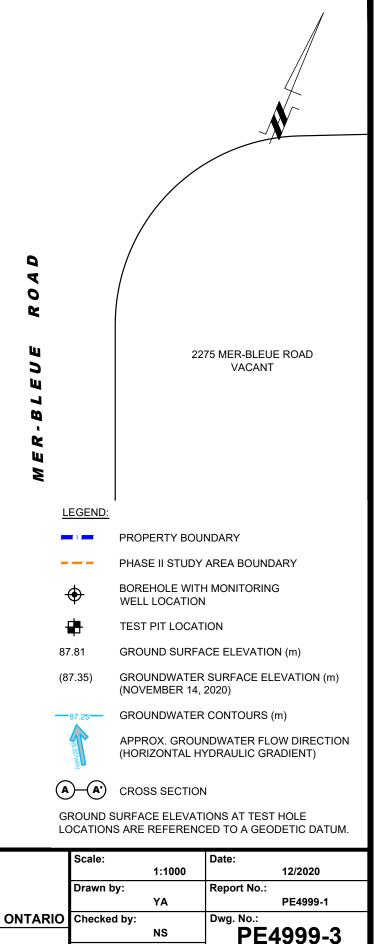


FIGURE 1 KEY PLAN

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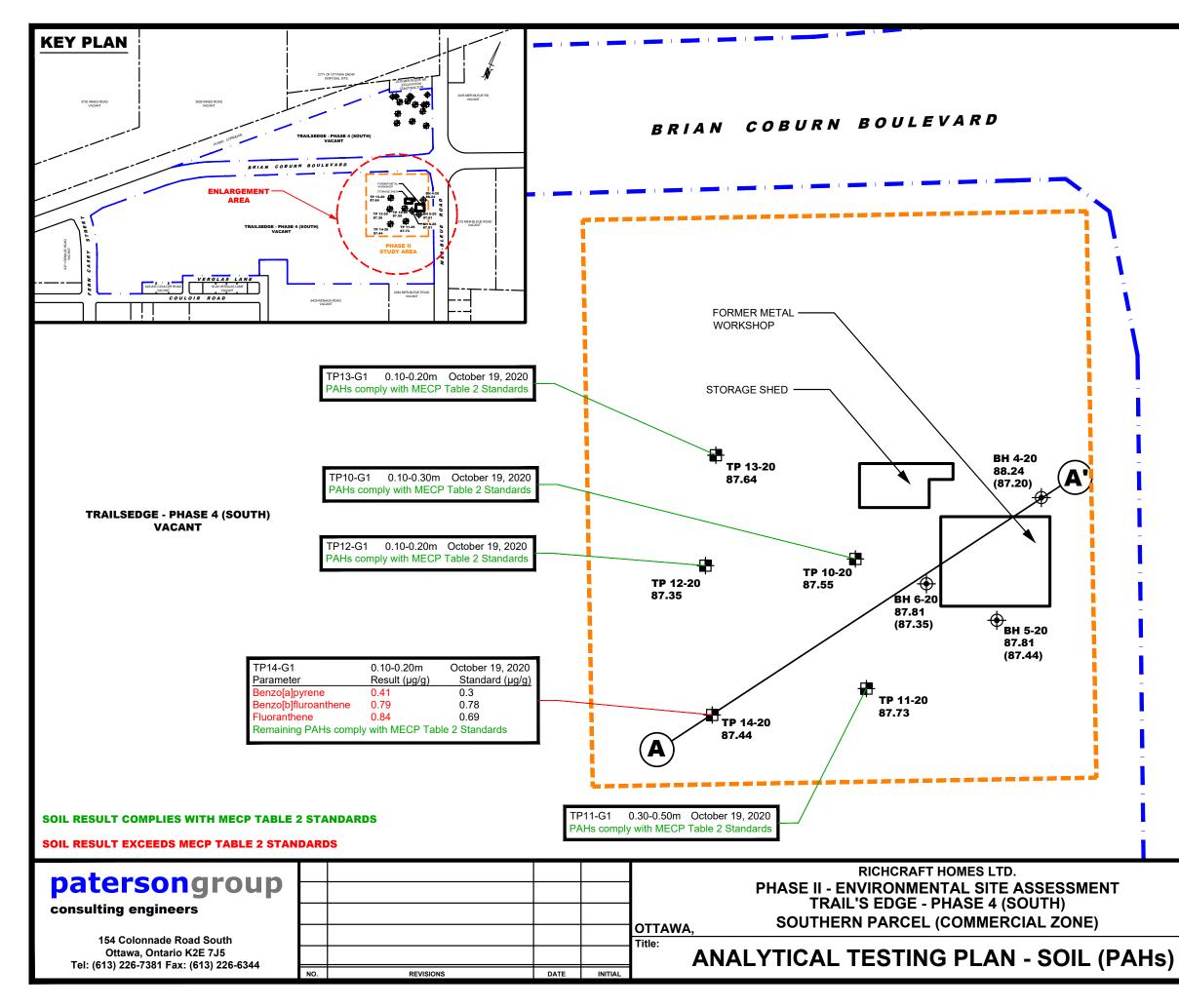


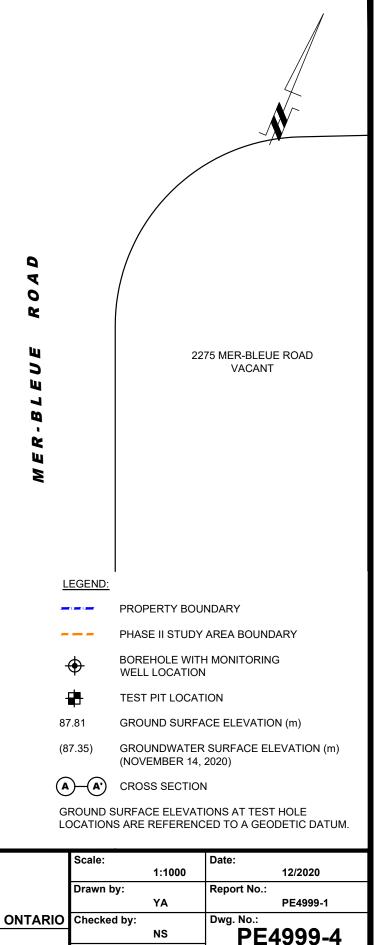
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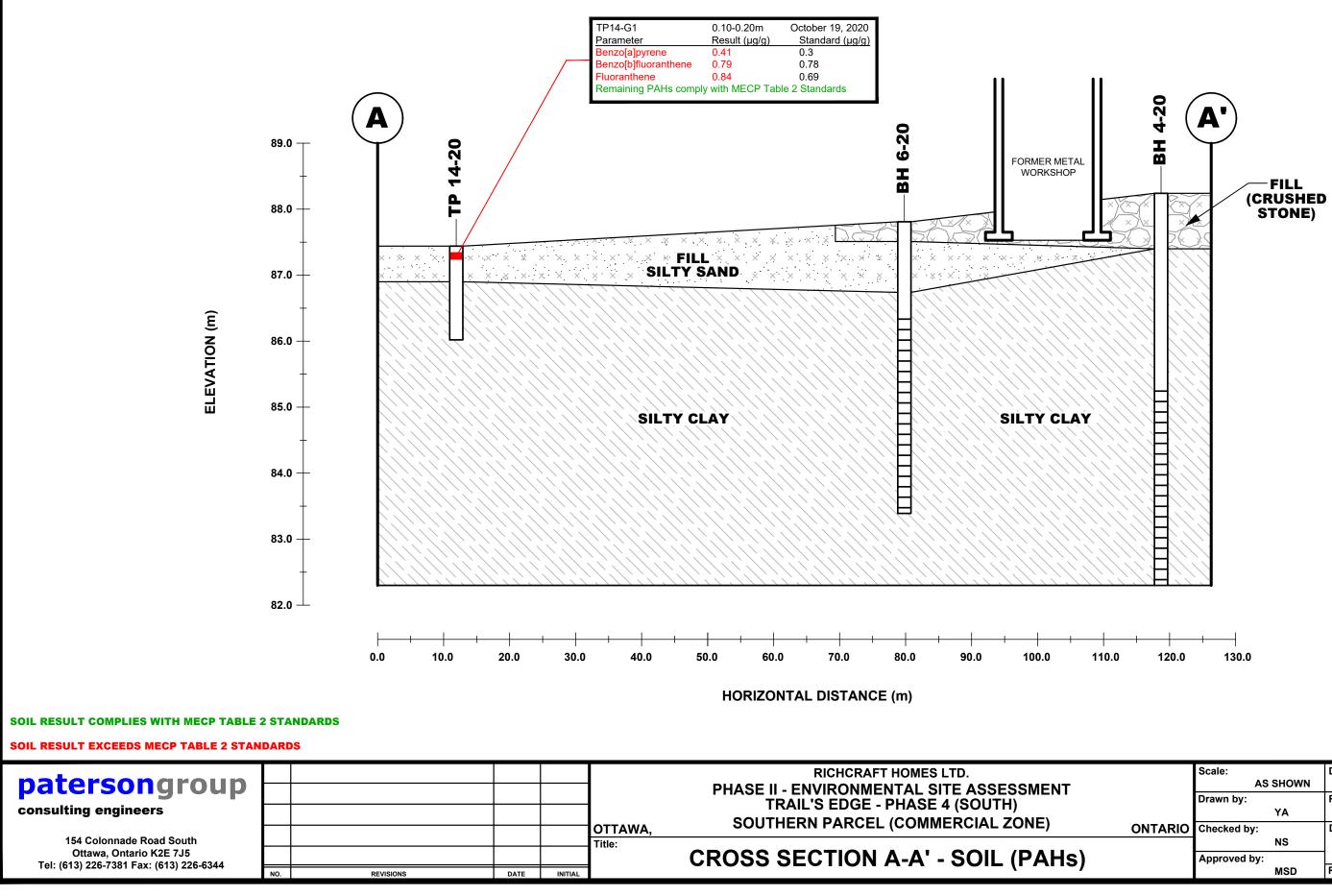


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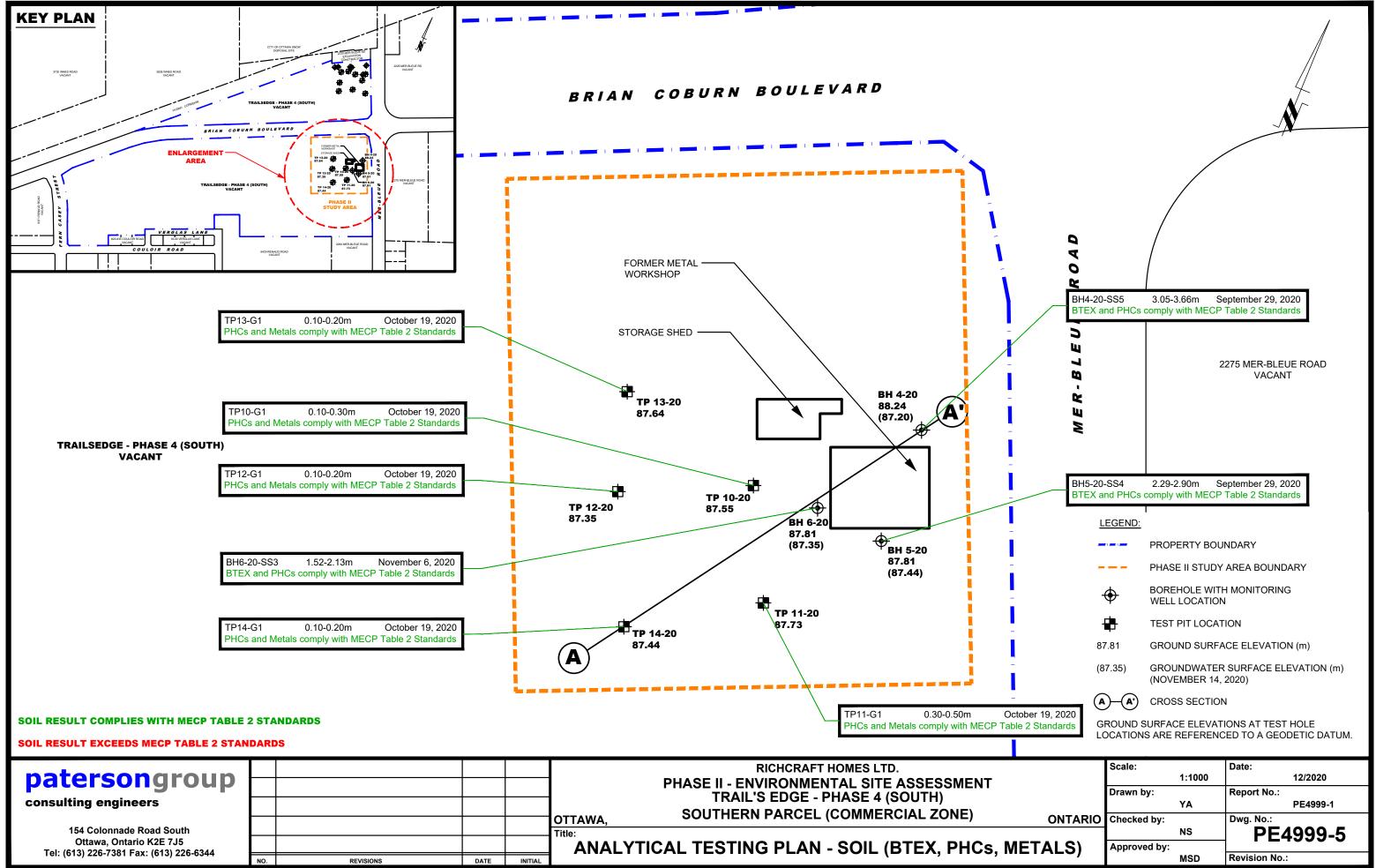
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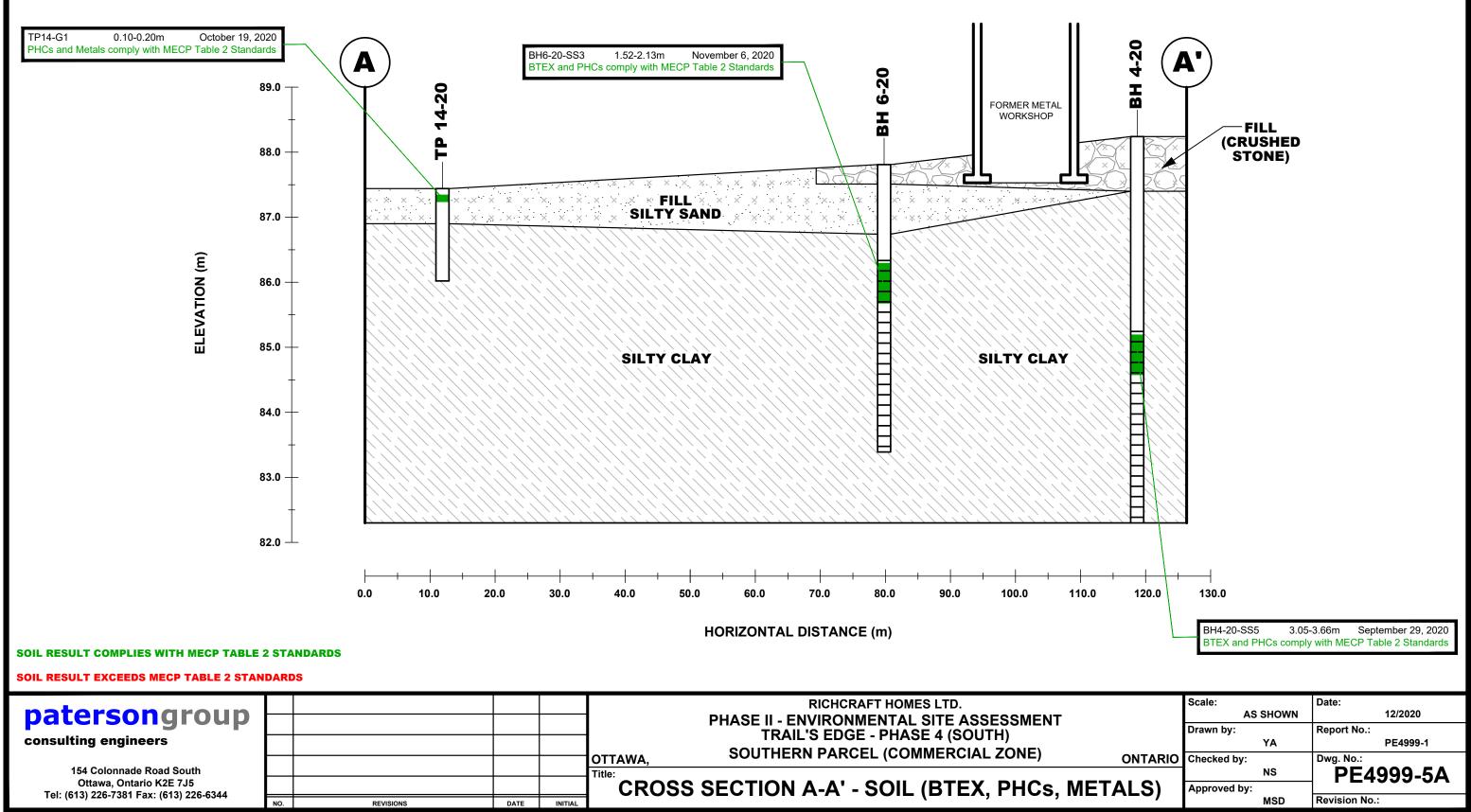
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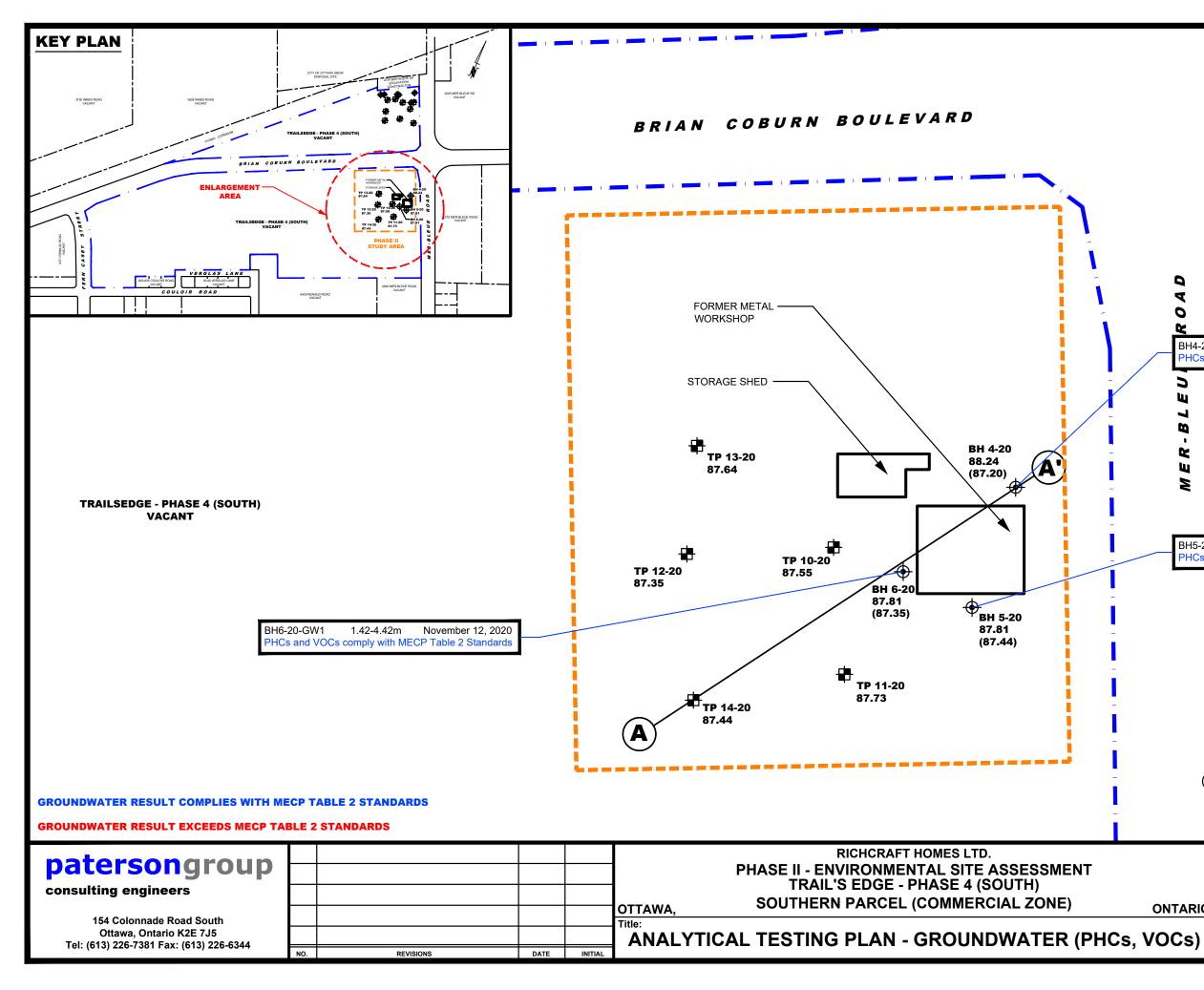
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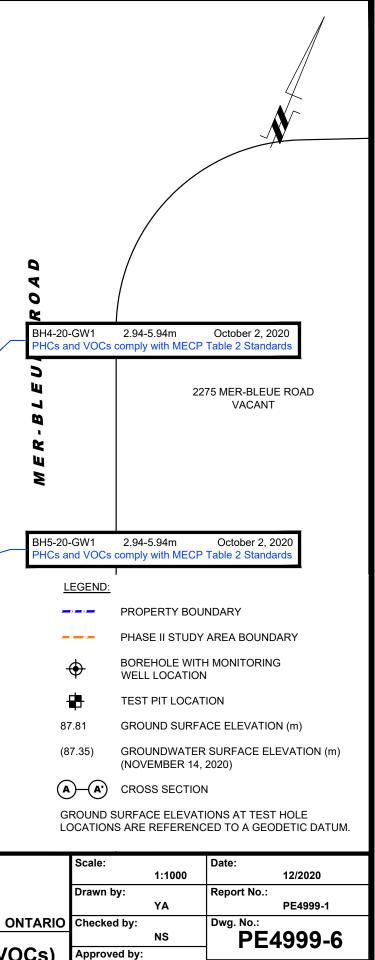


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		YA	PE4999-1
ONTARIO	Checked by:		Dwg. No.:
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TALS)	Approved by:		
.,,		MSD	Revision No.:

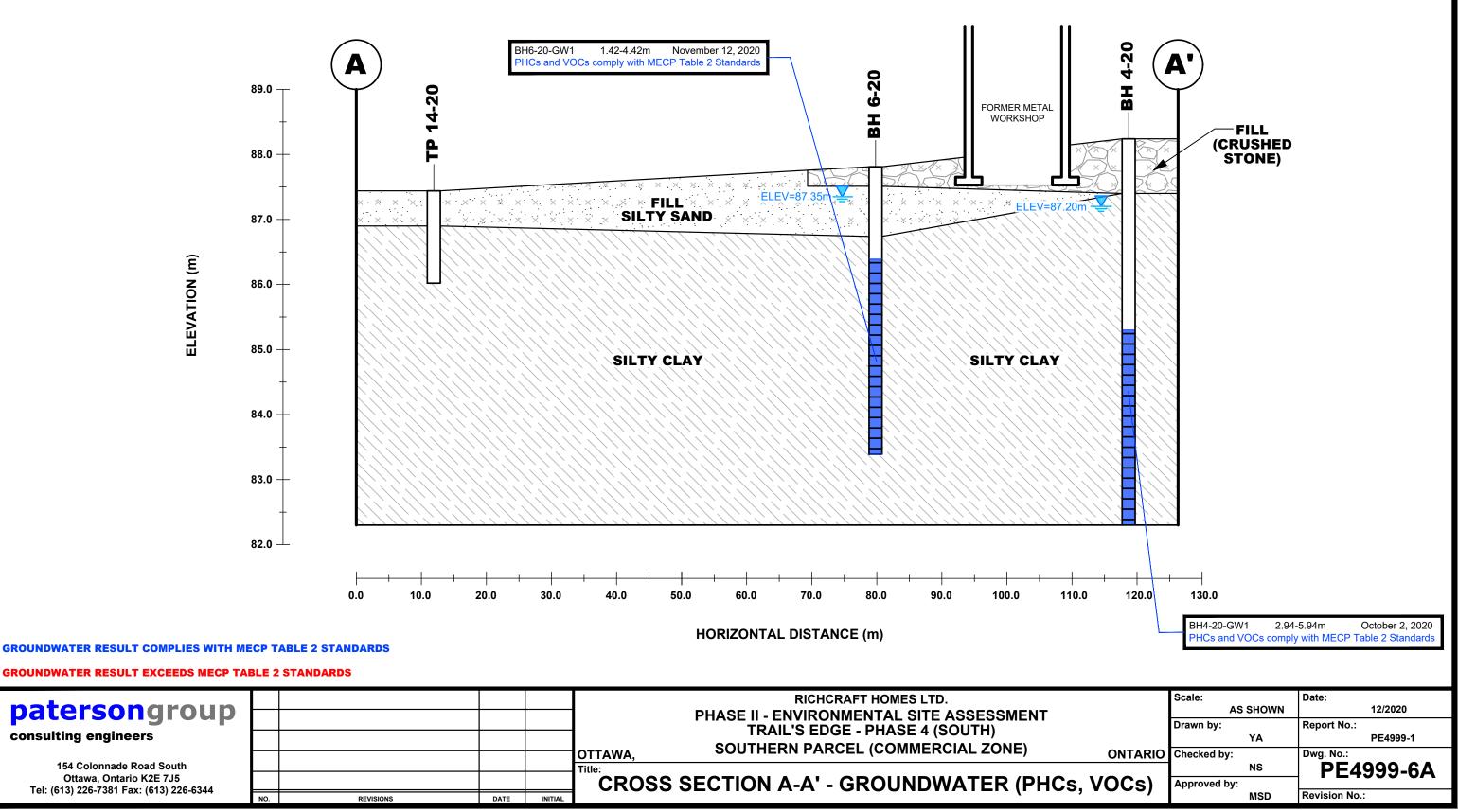




**Revision No.:** 

MSD

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	Scale:	Date:
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	Drawn by:	Report No.:
	YA	PE4999-1
ONTARIO	Checked by:	Dwg. No.:
	NS	<b>PE4999-6A</b>
VOCs)	Approved by:	
,	MSD	Revision No.:

# **APPENDIX 1**

# SAMPLING AND ANALYSIS PLAN

# SOIL PROFILE AND TEST DATA SHEETS

# SYMBOLS AND TERMS

# LABORATORY CERTIFICATES OF ANALYSIS

Geotechnical Engineering

Environmental Engineering

Hydrogeology

Geological Engineering

**Materials Testing** 

**Building Science** 

Archaeological Services

# Sampling & Analysis Plan

Phase II – Environmental Site Assessment Trail's Edge: Phase 4 (South) Southern Parcel (Commercial Zone) Part of 2284 Mer Bleue Road Ottawa, Ontario

**Prepared For** 

**Richcraft Group of Companies** 

### 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 September 21, 2020

Report: PE4999-1-SAP

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# 1.0 SAMPLING PROGRAM

Paterson Group Inc. (Paterson) was commissioned by Richcraft Group of Companies to conduct a Phase II – Environmental Site Assessment (Phase II ESA) for a portion of land within the proposed Trail's Edge: Phase 4 (South) subdivision development, in the City of Ottawa, Ontario.

This Phase II ESA was completed in conjunction with the Phase II ESA work carried out on another portion of the Trail's Edge: Phase 4 (South) development.

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

Borehole/ Test Pit	Location & Rationale	Proposed Depth & Rationale
BH4-20	Northern portion of Phase II study area; to assess for potential impacts resulting from a former metal workshop as well as the presence of fill material of unknown quality.	4-7 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH5-20	South-central portion of Phase II study area; to assess for potential impacts resulting from a former metal workshop as well as the presence of fill material of unknown quality.	4-7 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
BH6-20	Central portion of Phase II study area; to assess for potential impacts resulting from a former metal workshop as well as the presence of fill material of unknown quality.	4-7 m; to intercept the groundwater table for the purpose of installing a groundwater monitoring well.
TP10-TP14	Throughout the Phase II study area; to assess for potential impacts resulting from the presence of fill material of unknown quality.	0-0.5 m; for general coverage purposes.

Borehole and grab sample locations are shown on Drawing PE4999-3 – 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 all boreholes (BH4-20 to BH6-20) 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 PROCEDURES

### 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 F₁, 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.

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### 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.

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# 6.0 PHYSICAL IMPEDIMENTS

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
- **O** Other site-specific impediments

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

## SOIL PROFILE AND TEST DATA

**Phase II - Environmental Site Assessment** Trail's Edge: Phase 4 (South) Ottawa Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

						lawa, Oi	itano					
DATUM Geodetic									FILE NO.	PE4999	9	
REMARKS				_	(	O a va ta va la l		00	HOLE NO.	BH 4-2	20	
BORINGS BY Track-Mount Power Auge			SAN			Septembe	er 29, 20	Photo Ionization Detector				
SOIL DESCRIPTION	PLOT					DEPTH (m)	ELEV. (m)	Volatile Organic Rdg. (ppm)				
	STRATA	ТҮРЕ	NUMBER	% RECOVERY	VALUE r ROD			○ Lowe	r Explosive	Limit %	Monitoring Well Construction	
GROUND SURFACE	ST	H	ŊŊ	REC	N N		00.04	20	40 60	80	Į≥õ	
FILL: Crushed stone with organics		XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX	1				-88.24					
- with brown sand by 0.8m depth			1									
<u>0.84</u>		F 17										
		ss	2	62	7	1-	-87.24	•			<u>11111</u>	
Stiff, brown SILTY CLAY		ss	3	100	8	2-	-86.24					
Sun, brown SILTY CLAY												
<ul> <li>firm to soft and grey by 2.3m depth</li> </ul>		ss	4	100	3			•				
		$\square$										
		$\overline{\mathbf{N}}$				3-	-85.24					
		ss	5	100	W							
		ss	6	100	w	4-	-84.24					
		17										
		ss	7	100	w	5-	-83.24	•				
							00.24					
		$\sqrt{2}$		100								
5.94		ss	8	100	W					· · · · · · · · · · · · · · · · · · ·		
End of Borehole												
(GWL @ 1.04m - Nov. 12, 2020)												
									200 300 Eagle Rdg. (	(ppm)	00	
								▲ Full G	as Resp. 🛆 M	ethane Elim.		

## SOIL PROFILE AND TEST DATA

FILE NO.

**PE4999** 

Phase II - Environmental Site Assessment Trail's Edge: Phase 4 (South) Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

### REMARKS

DATUM

BORINGS BY Track-Me	ount Power	Auge
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Geodetic

BORINGS BY Track-Mount Power Aug	ər			п		Septemb	er 29-20	20	HOLE NO.	BH 5-2	20
SOIL DESCRIPTION	Р. СОТ	SAMPLI				DEPTH	ELEV.	Photo	Ionization I Atile Organic F	Detector	g Well Stion
	STRATA F	ЭДХТ	NUMBER	% RECOVERY	VALUE Pr ROD	(m)	(m)		er Explosiv		Monitoring Well Construction
GROUND SURFACE	S		N	RE	N 0 H		07.04	20	40 60	80	ž
FILL: Crushed stone with sand 0.23		AU	1			0-	-87.81	•			
		ss	2	92	9	1-	-86.81	•			որիներին որիներին ուներուներին ուներուներում։ 
Stiff, brown SILTY CLAY		ss	3	100	6	2-	-85.81	•			արությունը Անդեսներությունը
- firm to soft and grey by 2.3m depth		ss	4	100	2	3-	-84.81	•			
		ss	5	100	w			•			
		ss	6	100	W	4-	-83.81	•			
		ss	7	100	W	5-	-82.81	•			
5.94		ss	8	100	w						
(GWL @ 0.37m - Nov. 12, 2020)											
									200 300 Eagle Rdg. as Resp. △ N		⊣ ;00

### SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment Trail's Edge: Phase 4 (South) Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

### REMARKS

FILE NO.	PE4999
HOLE NO.	

#### HO **BH 6-20** BORINGS BY CME-55 Low Clearance Drill DATE November 6, 2020 SAMPLE **Photo Ionization Detector** Monitoring Well Construction STRATA PLOT DEPTH ELEV. SOIL DESCRIPTION • Volatile Organic Rdg. (ppm) (m) (m) N VALUE or RQD RECOVERY NUMBER TYPE o/0 O Lower Explosive Limit % **GROUND SURFACE** 80 20 40 60 0 + 87.81FILL: Crushed stone 0.30 AU 1 FILL: Brown silty sand with gravel 1.07 1 + 86.81SS 2 50 10 SS 3 100 6 Stiff, brown SILTY CLAY, some 2+85.81 sand - firm to soft and grey by 2.0m depth SS 4 100 2 3+84.81 SS 5 100 W 4+83.81 SS 6 W 100 4.42 End of Borehole (GWL @ 0.46m - Nov. 12, 2020) 100 200 300 400 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

## SOIL PROFILE AND TEST DATA

FILE NO.

Phase II - Environmental Site Assessment Trail's Edge: Phase 4 (South) Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

Geodetic

											L 110.	PE	4999	9
REMARKS										но	LE NO.	ТГ	P10	
BORINGS BY Backhoe	DATE October 19, 2020													T
SOIL DESCRIPTION	PLOT		SAN			DEPTH (m)	ELEV. (m)				<b>ation</b> rganic F			g Wel
	STRATA	ТҮРЕ	NUMBER	° ≈	N VALUE or RQD		(,	0	Lowe	er Ex	plosiv	e Lim	it %	Monitoring Well Construction
GROUND SURFACE	ū	•	Ĩ	REC	zö	0	-87.55		20	40	60	80	)	ž
FILL: Crushed stone0.10						- 0-	-87.55							
FILL: Brown silty sand, some crushed stone, trace clay, topsoil, organics		G 	1					•						
FILL: Brown silty clay, trace sand and gravel		G	2											
Brown <b>SILTY CLAY</b>		-				1-	-86.55							
End of Test Pit									100 RKI Full G	200 Eagle ias Re	300 ∋ Rdg.	(ppm	)	00

# SOIL PROFILE AND TEST DATA

FILE NO.

Phase II - Environmental Site Assessment Trail's Edge: Phase 4 (South) Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

Geodetic

										PE4999	9
REMARKS					HOLE NO.	<b>TP</b> 11					
BORINGS BY Backhoe				D			1611				
SOIL DESCRIPTION	PLOT		SAN	<b>IPLE</b>		DEPTH (m)	ELEV. (m)		onization tile Organic		Monitoring Well Construction
	STRATA	ТҮРЕ	NUMBER	% RECOVERY	N VALUE or RQD			○ Lowe	r Explosiv	ve Limit %	nitorin onstru
GROUND SURFACE	LS	H	<b>NN</b>	REC	Z O			20	40 60		₹O
FILL: Brown silty clay, some sand, trace gravel, cobbles and boulders		G	1			- 0-	-87.73	•			
0.64 TOPSOIL 0.91		G	2					•			
Brown <b>SILTY CLAY</b>		G	3			1-	-86.73				
								100 RKI E ▲ Full Ga	200 300 Eagle Rdg	0 400 50 . <b>(ppm)</b> Methane Elim.	00

# SOIL PROFILE AND TEST DATA

FILE NO.

Phase II - Environmental Site Assessment Trail's Edge: Phase 4 (South) Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

Geodetic

										PE499	9	
REMARKS						HOLE NO. TO 10						
BORINGS BY Backhoe				D	1	TP12						
SOIL DESCRIPTION	РГОТ		SAN	IPLE		DEPTH		Photo Ionization Detector     Volatile Organic Rdg. (ppm)				
			Ř	RY	Be	(m)	(m)			lio ridg. (ppin)	tring	
	STRATA	ТҮРЕ	NUMBER	% RECOVERY	N VALUE or RQD			O Lowe	r <b>Explo</b> a	sive Limit %	Monitoring Well Construction	
GROUND SURFACE	N		Z	RE	z <sup>o</sup>	0.	-87.35	20	40	60 80	ΣŬ	
FILL: Brown silty sand, some clay and crushed stone		G	1				- 67.33	•				
<u>0.38</u>		G	2					•				
Light reddish brown SILTY CLAY												
- brown by 0.8m depth						1.	-86.35					
1.44		G	3			1-	-80.30	•				
End of Test Pit		-							agle R	300 400 5 dg. (ppm) △ Methane Elim.		

# SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment Trail's Edge: Phase 4 (South) Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

DATUM Geodetic											F	ILE I	NO.		PI	E49	99	
REMARKS BORINGS BY Backhoe					ATE (	October 1	10 2020				н	IOLE	NO	).	Т	P13	3	
Bonings BT Dackride	PLOT		SAN	IPLE					Pho	oto	lon	izat	ion	De				
SOIL DESCRIPTION						DEPTH (m)	ELEV. (m)					Orga					W 20	Monitoring well Construction
	STRATA	ТҮРЕ	NUMBER	° ≈ © ©	VALUE r ROD			m)				xol	osi	ve	Lim	nit %		nstru
GROUND SURFACE	ST	Ĥ	ION	REC	N OL N					20		Explosive Lin				80		<u>50</u>
FILL: Brown silty clay, some sand and gravel, trace cobbles, topsoil		G	1			- 0-	-87.64	•										
<u>0.41</u>		-																
									· · · · · · · · · · · · · · · · · · ·		· · · · · · · · · · · · · · · · · · ·							
Light reddish brown <b>SILTY CLAY</b> - brown by 0.8m depth																		
						1-	-86.64				· · · · · · · · · · · · · · · · · · ·							
		_									· · · · · · · · · · · · · · · · · · ·							
1.81		G	2					•			· · · · · · · · · · · · · · · · · · ·							
End of Test Pit		-							• • • •		· · · · · · · · · · · · · · · · · · ·							
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# SOIL PROFILE AND TEST DATA

FILE NO.

Phase II - Environmental Site Assessment Trail's Edge: Phase 4 (South) Ottawa, Ontario

154 Colonnade Road South, Ottawa, Ontario K2E 7J5

Geodetic

DATUM Geodelic									FILE NO.	PE4999	9	
REMARKS BORINGS BY Backhoe				D	ATE	October 1	9. 2020		HOLE NO.	<b>TP14</b>		
	П						Photo I	o Ionization Detector 🧧				
SOIL DESCRIPTION	A PLOT		А			DEPTH (m)	ELEV. (m)	Vola	tile Organic F	Rdg. (ppm)	Monitoring Well	
	STRATA	ТҮРЕ	NUMBER	° ≈ © © © ©	N VALUE or RQD			○ Lowe	r Explosiv	e Limit %	onitol	
GROUND SURFACE	s S		z	RE	z <sup>0</sup>	0-	-87.44	20	40 60	80	Σ	
FILL: Brown/black silty sand, some clay and gravel, trace cobbles		G	1					•				
Brown SILTY CLAY		G	2			1-	-86.44	•				
1.42 End of Test Pit		-										
								100	200 300	400 5	- 00	
								100 RKI E ▲ Full Ga	200 300 Eagle Rdg. as Resp. △ M	400 50 (ppm) Aethane Elim.	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





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# Certificate of Analysis

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

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Nick Sullivan

Client PO: 30902 Project: PE4999 Custody: 128222

Report Date: 5-Oct-2020 Order Date: 30-Sep-2020

Order #: 2040420

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

Paracel ID	Client ID
2040420-01	BH2-20-AU1
2040420-02	BH2-20-SS5
2040420-03	BH4-20-SS5
2040420-04	BH5-20-SS4

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.



Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 30902

#### **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	1-Oct-20	1-Oct-20
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	1-Oct-20	3-Oct-20
Mercury by CVAA	EPA 7471B - CVAA, digestion	5-Oct-20	5-Oct-20
PHC F1	CWS Tier 1 - P&T GC-FID	1-Oct-20	1-Oct-20
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	1-Oct-20	5-Oct-20
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	2-Oct-20	2-Oct-20
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	1-Oct-20	4-Oct-20
Solids, %	Gravimetric, calculation	1-Oct-20	2-Oct-20

Order #: 2040420

Report Date: 05-Oct-2020 Order Date: 30-Sep-2020 Project Description: PE4999



#### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 30902

Order #: 2040420

Report Date: 05-Oct-2020

Order Date: 30-Sep-2020

Project Description: PE4999

	Client ID: Sample Date: Sample ID: MDL/Units	BH2-20-AU1 29-Sep-20 09:00 2040420-01 Soil	BH2-20-SS5 29-Sep-20 09:00 2040420-02 Soil	BH4-20-SS5 29-Sep-20 12:00 2040420-03 Soil	BH5-20-SS4 29-Sep-20 12:00 2040420-04 Soil		
Physical Characteristics	MDE/Offics	001			0011		
% Solids	0.1 % by Wt.	94.6	59.7	59.9	60.9		
Metals		0.10		ļ			
Antimony	1.0 ug/g dry	<1.0	-	-	-		
Arsenic	1.0 ug/g dry	4.2	-	-	-		
Barium	1.0 ug/g dry	101	-	-	-		
Beryllium	0.5 ug/g dry	<0.5	-	-	-		
Boron	5.0 ug/g dry	5.3	-	-	-		
Cadmium	0.5 ug/g dry	<0.5	-	-	-		
Chromium	5.0 ug/g dry	27.6	-	-	-		
Chromium (VI)	0.2 ug/g dry	<0.2	-	-	_		
Cobalt	1.0 ug/g dry	6.7	_	-	_		
Copper	5.0 ug/g dry	17.9	_	-	-		
Lead	1.0 ug/g dry	38.2	_	-	_		
Mercury	0.1 ug/g dry	0.1	-	-	-		
Molybdenum	1.0 ug/g dry	<1.0	_	-	_		
Nickel	5.0 ug/g dry	17.4	-	-			
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	31.3	_	-			
Zinc	20.0 ug/g dry	72.2	-	-			
Volatiles		12.2	-	ļ	-		
Benzene	0.02 ug/g dry	_	<0.02	<0.02	<0.02		
Ethylbenzene	0.05 ug/g dry	-	<0.05	<0.05	<0.02		
Toluene	0.05 ug/g dry	-	<0.05	<0.05	<0.05		
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		
Toluene-d8	Surrogate	-	114%	114%	114%		
Hydrocarbons					·		
F1 PHCs (C6-C10)	7 ug/g dry	-	<7	<7	<7		
F2 PHCs (C10-C16)	4 ug/g dry	-	<4	<4	<4		
F3 PHCs (C16-C34)	8 ug/g dry	-	<8	<8	<8		
F4 PHCs (C34-C50)	6 ug/g dry	-	<6	<6	<6		



#### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 30902

Report Date: 05-Oct-2020 Order Date: 30-Sep-2020

Project Description: PE4999

	Client ID:	BH2-20-AU1	BH2-20-SS5	BH4-20-SS5	BH5-20-SS4	
	Sample Date:	29-Sep-20 09:00	29-Sep-20 09:00	29-Sep-20 12:00	29-Sep-20 12:00	
	Sample ID:	2040420-01	2040420-02	2040420-03	2040420-04	
	MDL/Units	Soil	Soil	Soil	Soil	
Semi-Volatiles						
Acenaphthene	0.02 ug/g dry	0.08	-	-	-	
Acenaphthylene	0.02 ug/g dry	0.15	-	-	-	
Anthracene	0.02 ug/g dry	0.27	-	-	-	
Benzo [a] anthracene	0.02 ug/g dry	0.72	-	-	-	
Benzo [a] pyrene	0.02 ug/g dry	0.92	-	-	-	
Benzo [b] fluoranthene	0.02 ug/g dry	0.59	-	-	-	
Benzo [g,h,i] perylene	0.02 ug/g dry	0.62	-	-	-	
Benzo [k] fluoranthene	0.02 ug/g dry	0.57	-			
Chrysene	0.02 ug/g dry	0.68	-			
Dibenzo [a,h] anthracene	0.02 ug/g dry	0.16			-	
Fluoranthene	0.02 ug/g dry	1.37			-	
Fluorene	0.02 ug/g dry	0.10	-			
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	0.57	-	-	-	
1-Methylnaphthalene	0.02 ug/g dry	0.03	-	-	-	
2-Methylnaphthalene	0.02 ug/g dry	0.05	-	-	-	
Methylnaphthalene (1&2)	0.04 ug/g dry	0.08	-	-	-	
Naphthalene	0.01 ug/g dry	0.11	-	-	-	
Phenanthrene	0.02 ug/g dry	0.94	-	-	-	
Pyrene	0.02 ug/g dry	1.26	-	-	-	
2-Fluorobiphenyl	Surrogate	112%	-	-	-	
Terphenyl-d14	Surrogate	86.5%	-	-	-	



Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 30902

# Order #: 2040420

Report Date: 05-Oct-2020

Order Date: 30-Sep-2020

Project Description: PE4999

#### 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						
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 Molybdenum	ND ND	0.1 1.0	ug/g						
Nickel	ND	5.0	ug/g						
Selenium	ND	1.0	ug/g ug/g						
Silver	ND	0.3	ug/g ug/g						
Thallium	ND	1.0	ug/g ug/g						
Uranium	ND	1.0	ug/g						
Vanadium	ND	10.0	ug/g						
Zinc	ND	20.0	ug/g						
Semi-Volatiles			-9.9						
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 0.04	ug/g						
Methylnaphthalene (1&2) Naphthalene	ND ND	0.04	ug/g						
Phenanthrene	ND	0.01	ug/g						
Pyrene	ND	0.02	ug/g						
Surrogate: 2-Fluorobiphenyl	1.18	0.02	ug/g <i>ug/g</i>		88.1	50-140			
Surrogate: Terphenyl-d14	1.16		ug/g ug/g		87.3	50-140			
Volatiles	1.10		uy/y		07.5	50-140			
Benzene		0.02							
Ethylbenzene	ND ND	0.02	ug/g						
Toluene	ND	0.05	ug/g						
m,p-Xylenes	ND	0.05	ug/g ug/g						
o-Xylene	ND	0.05							
Xylenes, total	ND	0.05	ug/g ug/g						
Surrogate: Toluene-d8	9.09	0.00	ug/g ug/g		114	50-140			
Canogato. Totacho ao	3.03		ug/g			00 170			



Certificate of Analysis Client: Paterson Group Consulting Engineers

Client PO: 30902

Report Date: 05-Oct-2020

Order Date: 30-Sep-2020

Project Description: PE4999

#### Method Quality Control: Duplicate

Analyte	Pocult	Reporting Limit	1.16.24	Source	0/ 050	%REC		RPD	Notaa
лицую	Result		Units	Result	%REC	Limit	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	ND	1.0	ug/g dry	ND			NC	30	
Arsenic	1.9	1.0	ug/g dry	2.1			9.2	30	
Barium	45.7	1.0	ug/g dry	46.7			2.3	30	
Beryllium	ND	0.5	ug/g dry	ND			NC	30	
Boron	6.0	5.0	ug/g dry	5.7			5.4	30	
Cadmium	ND	0.5	ug/g dry	ND			NC	30	
Chromium (VI)	ND	0.2	ug/g dry	ND			NC	35	
Chromium	16.4	5.0	ug/g dry	16.4			0.3	30	
Cobalt	4.9	1.0	ug/g dry	4.7			3.9	30	
Copper	9.2	5.0	ug/g dry	9.4			1.5	30	
Lead	4.3	1.0	ug/g dry	4.5			2.8	30	
Mercury	ND	0.1	ug/g dry	ND			NC	30	
Molybdenum	ND	1.0	ug/g dry	ND			NC	30	
Nickel	9.3	5.0	ug/g dry	9.3			0.2	30	
Selenium	ND	1.0	ug/g dry	ND			NC	30	
Silver	ND	0.3	ug/g dry	ND			NC	30	
Thallium	ND	1.0	ug/g dry	ND			NC	30	
Uranium	ND	1.0	ug/g dry	ND			NC	30	
Vanadium	26.4	10.0	ug/g dry	26.4			0.1	30	
Zinc	24.1	20.0	ug/g dry	23.8			1.6	30	
Physical Characteristics		20.0		20.0					
% Solids	96.9	0.1	% by Wt.	95.9			1.0	25	
Semi-Volatiles	00.0	0.1	<i>10 by</i> 110	00.0			1.0	20	
Acenaphthene	0.094	0.02	ug/g dry	0.083			12.8	40	
Acenaphthylene	0.124	0.02	ug/g dry	0.146			17.0	40	
Anthracene	0.296	0.02	ug/g dry	0.273			7.9	40	
Benzo [a] anthracene	0.663	0.02	ug/g dry ug/g dry	0.720			8.2	40	
Benzo [a] pyrene	0.835	0.02	ug/g dry ug/g dry	0.921			9.8	40	
Benzo [b] fluoranthene	0.951	0.02	ug/g dry ug/g dry	0.589			47.1	40	QR-04
Benzo [g,h,i] perylene	0.562	0.02	ug/g dry ug/g dry	0.623			10.2	40	QIT 04
Benzo [k] fluoranthene	0.502	0.02	ug/g dry ug/g dry	0.023			11.5	40	
Chrysene	0.637	0.02	ug/g dry ug/g dry	0.679			6.4	40	
Dibenzo [a,h] anthracene	0.037	0.02		0.079			10.5	40	
Fluoranthene	1.36	0.02	ug/g dry	1.37			0.9	40	
Fluorene	0.112	0.02	ug/g dry	0.098			0.9 12.6	40 40	
	0.112	0.02	ug/g dry	0.098			12.0	40 40	
Indeno [1,2,3-cd] pyrene	0.036	0.02	ug/g dry	0.072				40 40	
1-Methylnaphthalene			ug/g dry				12.1 13.7		
2-Methylnaphthalene	0.054	0.02	ug/g dry	0.047			13.7	40	
Naphthalene	0.128	0.01	ug/g dry	0.107				40	
Phenanthrene	1.04	0.02	ug/g dry	0.943			9.7 3.5	40 40	
Pyrene	1.22	0.02	ug/g dry	1.26	100	50-140	3.5	40	
Surrogate: 2-Fluorobiphenyl	1.53		ug/g dry		109				
Surrogate: Terphenyl-d14	1.17		ug/g dry		82.8	50-140			
Volatiles									
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	15.4		ug/g dry		115	50-140			



# Method Quality Control: Spike

Report Date: 05-Oct-2020

Order Date: 30-Sep-2020

Project Description: PE4999

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit Notes
Hydrocarbons								
F1 PHCs (C6-C10)	181	7	ug/g	ND	90.7	80-120		
F2 PHCs (C10-C16)	144	4	ug/g	ND	107	60-140		
F3 PHCs (C16-C34)	359	8	ug/g	ND	109	60-140		
F4 PHCs (C34-C50)	229	6	ug/g	ND	111	60-140		
Metals								
Antimony	44.8	1.0	ug/g	ND	89.2	70-130		
Arsenic	53.1	1.0	ug/g	ND	105	70-130		
Barium	68.5	1.0	ug/g	18.7	99.7	70-130		
Beryllium	48.7	0.5	ug/g	ND	97.1	70-130		
Boron	43.9	5.0	ug/g	ND	83.3	70-130		
Cadmium	49.3	0.5	ug/g	ND	98.6	70-130		
Chromium (VI)	0.1	0.2	ug/g	ND	60.5	70-130		QM-05
Chromium	58.7	5.0	ug/g	6.6	104	70-130		
Cobalt	52.0	1.0	ug/g	1.9	100	70-130		
Copper	52.2	5.0	ug/g	ND	96.9	70-130		
Lead	50.4	1.0	ug/g	1.8	97.3	70-130		
Mercury	1.53	0.1	ug/g	ND	102	70-130		
Molybdenum	49.5	1.0	ug/g	ND	98.8	70-130		
Nickel	52.5	5.0	ug/g	ND	97.6	70-130		
Selenium	48.4	1.0	ug/g	ND	96.7	70-130		
Silver	42.6	0.3	ug/g	ND	85.3	70-130		
Thallium	49.3	1.0	ug/g	ND	98.5	70-130		
Uranium	52.0	1.0	ug/g	ND	104	70-130		
Vanadium	62.2	10.0	ug/g	10.6	103	70-130		
Zinc	55.8	20.0	ug/g	ND	92.6	70-130		
Semi-Volatiles								
Acenaphthene	0.123	0.02	ug/g	ND	74.0	50-140		
Acenaphthylene	0.107	0.02	ug/g	ND	64.3	50-140		
Anthracene	0.116	0.02	ug/g	ND	69.7	50-140		
Benzo [a] anthracene	0.095	0.02	ug/g	ND	57.3	50-140		
Benzo [a] pyrene	0.093	0.02	ug/g	ND	56.0	50-140		
Benzo [b] fluoranthene	0.144	0.02	ug/g	ND	86.2	50-140		
Benzo [g,h,i] perylene	0.112	0.02	ug/g	ND	67.0	50-140		
Benzo [k] fluoranthene	0.130	0.02	ug/g	ND	78.2	50-140		
Chrysene	0.124	0.02	ug/g	ND	74.6	50-140		
Dibenzo [a,h] anthracene	0.108	0.02	ug/g	ND	64.6	50-140		
Fluoranthene	0.113	0.02	ug/g	ND	67.8	50-140		
Fluorene	0.113	0.02	ug/g	ND	67.6	50-140		
Indeno [1,2,3-cd] pyrene	0.111	0.02	ug/g	ND	66.3	50-140		
1-Methylnaphthalene	0.135	0.02	ug/g	ND	81.0	50-140		
2-Methylnaphthalene	0.149	0.02	ug/g	ND	89.5	50-140		
Naphthalene	0.129	0.01	ug/g	ND	77.5	50-140		
Phenanthrene	0.121	0.02	ug/g	ND	72.4	50-140		
Pyrene	0.114	0.02	ug/g	ND	68.2	50-140		
Surrogate: 2-Fluorobiphenyl Surrogate: Terphenyl-d14	0.960 0.962		ug/g ug/g		72.0 72.2	50-140 50-140		
Volatiles								
Benzene	3.93	0.02	ug/g	ND	98.3	60-130		



Report Date: 05-Oct-2020 Order Date: 30-Sep-2020

51der Date: 30-8ep-2020

Project Description: PE4999

## Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Ethylbenzene	3.97	0.05	ug/g	ND	99.2	60-130			
Toluene	3.88	0.05	ug/g	ND	97.0	60-130			
m,p-Xylenes	8.28	0.05	ug/g	ND	104	60-130			
o-Xylene	4.06	0.05	ug/g	ND	101	60-130			
Surrogate: Toluene-d8	7.89		ug/g		98.6	50-140			



#### **Qualifier Notes:**

#### QC Qualifiers :

QM-05: The spike recovery was outside acceptance limits for the matrix spike due to matrix interference.

QR-04 : Duplicate results exceeds RPD limits due to non-homogeneous matrix.

#### 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.

<b>PARAC</b> LABORATORIES	Parace					ent Blvd. (1G 4J8 ) ellabs.com .com		(Lab	Use	Num Only)				(Li	ab Use	Custody Only) 28222	
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# Certificate of Analysis

# **Paterson Group Consulting Engineers**

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

Client PO: 31209 Project: PE4999 Custody: 52670

Report Date: 13-Nov-2020 Order Date: 10-Nov-2020

Order #: 2046234

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

Paracel ID 2046234-01

**Client ID** BH6-20-SS3

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.



Report Date: 13-Nov-2020 Order Date: 10-Nov-2020

Order #: 2046234

Project Description: PE4999

## **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	12-Nov-20	13-Nov-20
PHC F1	CWS Tier 1 - P&T GC-FID	12-Nov-20	13-Nov-20
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	11-Nov-20	13-Nov-20
Solids, %	Gravimetric, calculation	11-Nov-20	12-Nov-20



#### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 31209

Report Date: 13-Nov-2020

Order Date: 10-Nov-2020

Project Description: PE4999

	_				
	Client ID:	BH6-20-SS3	-	-	-
	Sample Date:	06-Nov-20 09:00	-	-	-
	Sample ID:	2046234-01	-	-	-
	MDL/Units	Soil	-	-	-
Physical Characteristics					
% Solids	0.1 % by Wt.	67.7	-	-	-
Volatiles	•		•		
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	119%	-	-	-
Hydrocarbons					
F1 PHCs (C6-C10)	7 ug/g dry	<7	-	-	-
F2 PHCs (C10-C16)	4 ug/g dry	<4	-	-	-
F3 PHCs (C16-C34)	8 ug/g dry	<8	-	-	-
F4 PHCs (C34-C50)	6 ug/g dry	<6	-	-	-



Order #: 2046234

Report Date: 13-Nov-2020

Order Date: 10-Nov-2020

Project Description: PE4999

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						
Volatiles									
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	9.47		ug/g		118	50-140			



Client PO: 31209

Report Date: 13-Nov-2020

Order Date: 10-Nov-2020

Project Description: PE4999

## 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	
Physical Characteristics									
% Solids	91.5	0.1	% by Wt.	91.2			0.3	25	
Volatiles									
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	10.9		ug/g dry		119	50-140			



Report Date: 13-Nov-2020

Order Date: 10-Nov-2020

Project Description: PE4999

## Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	163	7	ug/g	ND	81.5	80-120			
F2 PHCs (C10-C16)	93	4	ug/g	ND	91.3	60-140			
F3 PHCs (C16-C34)	226	8	ug/g	ND	90.6	60-140			
F4 PHCs (C34-C50)	140	6	ug/g	ND	88.9	60-140			
Volatiles									
Benzene	3.55	0.02	ug/g	ND	88.7	60-130			
Ethylbenzene	3.97	0.05	ug/g	ND	99.3	60-130			
Toluene	4.21	0.05	ug/g	ND	105	60-130			
m,p-Xylenes	8.32	0.05	ug/g	ND	104	60-130			
o-Xylene	3.89	0.05	ug/g	ND	97.2	60-130			
Surrogate: Toluene-d8	8.32		ug/g		104	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. 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.

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Revision 3.0



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# Certificate of Analysis

# **Paterson Group Consulting Engineers**

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Nick Sullivan

Client PO: 31420 Project: PE4999 Custody: 55010/11

Report Date: 26-Oct-2020 Order Date: 20-Oct-2020

Order #: 2043350

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

Client ID
Client ID
TP1-G2
TP2-G1
TP4-G2
TP5-G2
TP7-G1
TP8-G2
TP10-G1
TP11-G1
TP12-G1
TP13-G1
TP14-G1

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	21-Oct-20	23-Oct-20
Mercury by CVAA	EPA 7471B - CVAA, digestion	23-Oct-20	23-Oct-20
PHC F4G (gravimetric)	CWS Tier 1 - Extraction Gravimetric	26-Oct-20	26-Oct-20
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	21-Oct-20	22-Oct-20
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	23-Oct-20	23-Oct-20
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	21-Oct-20	23-Oct-20
Solids, %	Gravimetric, calculation	21-Oct-20	22-Oct-20

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

Report Date: 26-Oct-2020 Order Date: 20-Oct-2020

Project Description: PE4999



#### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 31420

Report Date: 26-Oct-2020

Order Date: 20-Oct-2020

Project Description: PE4999

	Client ID: Sample Date: Sample ID: MDL/Units	TP1-G2 19-Oct-20 09:00 2043350-01 Soil	TP2-G1 19-Oct-20 09:00 2043350-02 Soil	TP4-G2 19-Oct-20 09:00 2043350-03 Soil	TP5-G2 19-Oct-20 09:00 2043350-04 Soil
Physical Characteristics	MDE/Onits				
% Solids	0.1 % by Wt.	77.8	85.9	69.1	75.6
Metals		-			
Antimony	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Arsenic	1.0 ug/g dry	3.5	2.6	4.1	3.9
Barium	1.0 ug/g dry	152	45.2	300	149
Beryllium	0.5 ug/g dry	0.6	<0.5	0.8	0.6
Boron	5.0 ug/g dry	<5.0	<5.0	5.9	<5.0
Cadmium	0.5 ug/g dry	<0.5	<0.5	<0.5	<0.5
Chromium	5.0 ug/g dry	70.0	19.5	118	66.8
Chromium (VI)	0.2 ug/g dry	<0.2	<0.2	<0.2	<0.2
Cobalt	1.0 ug/g dry	13.3	5.1	22.3	11.9
Copper	5.0 ug/g dry	23.5	11.3	50.5	22.3
Lead	1.0 ug/g dry	10.8	8.2	7.7	13.9
Mercury	0.1 ug/g dry	<0.1	<0.1	<0.1	<0.1
Molybdenum	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Nickel	5.0 ug/g dry	34.2	12.7	62.6	31.7
Selenium	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Silver	0.3 ug/g dry	<0.3	<0.3	<0.3	<0.3
Thallium	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Uranium	1.0 ug/g dry	1.7	<1.0	<1.0	2.2
Vanadium	10.0 ug/g dry	59.9	23.2	104	57.2
Zinc	20.0 ug/g dry	77.4	33.3	118	84.2
Hydrocarbons			•		
F2 PHCs (C10-C16)	4 ug/g dry	<4	<4	<4	<4
F3 PHCs (C16-C34)	8 ug/g dry	<8	<8	<8	<8
F4 PHCs (C34-C50)	6 ug/g dry	<6	<6	<6	<6
Semi-Volatiles					
Acenaphthene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Acenaphthylene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Anthracene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Benzo [a] anthracene	0.02 ug/g dry	<0.02	0.05	<0.02	<0.02
Benzo [a] pyrene	0.02 ug/g dry	<0.02	0.07	<0.02	<0.02
Benzo [b] fluoranthene	0.02 ug/g dry	<0.02	0.09	<0.02	<0.02
Benzo [g,h,i] perylene	0.02 ug/g dry	<0.02	0.05	<0.02	<0.02
Benzo [k] fluoranthene	0.02 ug/g dry	<0.02	0.04	<0.02	<0.02



Order #: 2043350

Report Date: 26-Oct-2020 Order Date: 20-Oct-2020

Project Description: PE4999

	Client ID: Sample Date:	TP1-G2 19-Oct-20 09:00	TP2-G1 19-Oct-20 09:00	TP4-G2 19-Oct-20 09:00	TP5-G2 19-Oct-20 09:00
	Sample ID: MDL/Units	2043350-01 Soil	2043350-02 Soil	2043350-03 Soil	2043350-04 Soil
Chrysene	0.02 ug/g dry	<0.02	0.07	<0.02	<0.02
Dibenzo [a,h] anthracene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Fluoranthene	0.02 ug/g dry	<0.02	0.13	<0.02	0.03
Fluorene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	<0.02	0.05	<0.02	<0.02
1-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
2-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Methylnaphthalene (1&2)	0.04 ug/g dry	<0.04	<0.04	<0.04	<0.04
Naphthalene	0.01 ug/g dry	<0.01	<0.01	<0.01	<0.01
Phenanthrene	0.02 ug/g dry	<0.02	0.04	<0.02	<0.02
Pyrene	0.02 ug/g dry	<0.02	0.11	<0.02	0.02
2-Fluorobiphenyl	Surrogate	70.4%	84.9%	68.2%	72.3%
Terphenyl-d14	Surrogate	104%	132%	101%	110%



#### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 31420

Report Date: 26-Oct-2020 Order Date: 20-Oct-2020

Project Description: PE4999

	Client ID: Sample Date: Sample ID: MDL/Units	Sample Date: 19-Oct-20 09:00 19-Oct-2		TP10-G1 19-Oct-20 09:00 2043350-07 Soil	TP11-G1 19-Oct-20 09:00 2043350-08 Soil
Physical Characteristics	INDEPOINTS				
% Solids	0.1 % by Wt.	72.6	78.8	93.2	76.5
Metals					
Antimony	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Arsenic	1.0 ug/g dry	3.2	2.9	1.8	3.3
Barium	1.0 ug/g dry	209	106	55.3	167
Beryllium	0.5 ug/g dry	0.7	<0.5	<0.5	0.6
Boron	5.0 ug/g dry	5.6	<5.0	<5.0	<5.0
Cadmium	0.5 ug/g dry	<0.5	<0.5	<0.5	<0.5
Chromium	5.0 ug/g dry	102	41.8	15.4	69.2
Chromium (VI)	0.2 ug/g dry	<0.2	<0.2	<0.2	<0.2
Cobalt	1.0 ug/g dry	18.0	8.6	3.6	13.7
Copper	5.0 ug/g dry	31.5	16.5	5.4	26.8
Lead	1.0 ug/g dry	13.7	23.9	4.0	7.8
Mercury	0.1 ug/g dry	<0.1	<0.1	<0.1	<0.1
Molybdenum	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Nickel	5.0 ug/g dry	50.5	21.0	9.3	38.5
Selenium	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Silver	0.3 ug/g dry	<0.3	<0.3	<0.3	<0.3
Thallium	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Uranium	1.0 ug/g dry	1.8	1.2	<1.0	<1.0
Vanadium	10.0 ug/g dry	78.6	39.0	15.9	54.5
Zinc	20.0 ug/g dry	98.2	78.3	<20.0	65.9
Hydrocarbons	· · ·			-	
F2 PHCs (C10-C16)	4 ug/g dry	<4	<4	<4	<4
F3 PHCs (C16-C34)	8 ug/g dry	<8	<8	<8	<8
F4 PHCs (C34-C50)	6 ug/g dry	<6	<6	<6	<6
Semi-Volatiles					
Acenaphthene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Acenaphthylene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Anthracene	0.02 ug/g dry	<0.02	0.02	<0.02	<0.02
Benzo [a] anthracene	0.02 ug/g dry	<0.02	0.06	<0.02	<0.02
Benzo [a] pyrene	0.02 ug/g dry	<0.02	2 0.08 <0.02		<0.02
Benzo [b] fluoranthene	0.02 ug/g dry	<0.02	2 0.10 <0.02		<0.02
Benzo [g,h,i] perylene	0.02 ug/g dry	<0.02	0.06 <0.02		<0.02
Benzo [k] fluoranthene	0.02 ug/g dry	<0.02	0.05	<0.02	<0.02



Order #: 2043350

Report Date: 26-Oct-2020 Order Date: 20-Oct-2020

Project Description: PE4999

	F		•		-
	Client ID:	TP7-G1	TP8-G2	TP10-G1	TP11-G1
	Sample Date:	19-Oct-20 09:00	19-Oct-20 09:00	19-Oct-20 09:00	19-Oct-20 09:00
	Sample ID:	2043350-05	2043350-06	2043350-07	2043350-08
	MDL/Units	Soil	Soil	Soil	Soil
Chrysene	0.02 ug/g dry	<0.02	0.08	<0.02	<0.02
Dibenzo [a,h] anthracene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Fluoranthene	0.02 ug/g dry	<0.02	0.14	<0.02	<0.02
Fluorene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	<0.02	0.05	<0.02	<0.02
1-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
2-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.02	<0.02
Methylnaphthalene (1&2)	0.04 ug/g dry	<0.04	<0.04	<0.04	<0.04
Naphthalene	0.01 ug/g dry	<0.01	<0.01	<0.01	<0.01
Phenanthrene	0.02 ug/g dry	<0.02	0.05	<0.02	<0.02
Pyrene	0.02 ug/g dry	<0.02	0.12	<0.02	<0.02
2-Fluorobiphenyl	Surrogate	61.8%	75.2%	76.5%	66.0%
Terphenyl-d14	Surrogate	83.5%	118%	102%	70.6%



#### Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 31420

Order #: 2043350

Report Date: 26-Oct-2020

Order Date: 20-Oct-2020

Project Description: PE4999

	Client ID: Sample Date: Sample ID: MDL/Units	TP12-G1 19-Oct-20 09:00 2043350-09 Soil	TP13-G1 19-Oct-20 09:00 2043350-10 Soil	TP14-G1 19-Oct-20 09:00 2043350-11 Soil	- - -
Physical Characteristics	WDL/OTITS	001			
% Solids	0.1 % by Wt.	81.0	80.8	87.2	-
Metals					
Antimony	1.0 ug/g dry	<1.0	<1.0	<1.0	-
Arsenic	1.0 ug/g dry	4.4	3.0	3.1	-
Barium	1.0 ug/g dry	157	116	138	-
Beryllium	0.5 ug/g dry	0.6	<0.5	<0.5	-
Boron	5.0 ug/g dry	6.2	6.2	5.4	-
Cadmium	0.5 ug/g dry	<0.5	<0.5	<0.5	-
Chromium	5.0 ug/g dry	69.1	48.2	53.5	-
Chromium (VI)	0.2 ug/g dry	<0.2	<0.2	<0.2	-
Cobalt	1.0 ug/g dry	12.0	9.0	10.9	-
Copper	5.0 ug/g dry	22.8	17.2	23.3	-
Lead	1.0 ug/g dry	12.8	11.6	15.1	-
Mercury	0.1 ug/g dry	<0.1	<0.1	<0.1	-
Molybdenum	1.0 ug/g dry	1.2	<1.0	<1.0	-
Nickel	5.0 ug/g dry	35.5	25.0	31.1	-
Selenium	1.0 ug/g dry	<1.0	<1.0	<1.0	-
Silver	0.3 ug/g dry	<0.3	<0.3	<0.3	-
Thallium	1.0 ug/g dry	<1.0	<1.0	<1.0	-
Uranium	1.0 ug/g dry	2.4	1.3	<1.0	-
Vanadium	10.0 ug/g dry	56.1	39.6	51.3	-
Zinc	20.0 ug/g dry	84.2	53.7	65.1	-
Hydrocarbons					
F2 PHCs (C10-C16)	4 ug/g dry	<4	<4	<4	-
F3 PHCs (C16-C34)	8 ug/g dry	40	<8	60	-
F4 PHCs (C34-C50)	6 ug/g dry	122	<6	188 [1]	-
F4G PHCs (gravimetric)	50 ug/g dry	-	-	722	-
Semi-Volatiles					
Acenaphthene	0.02 ug/g dry	<0.02	<0.02	<0.02	-
Acenaphthylene	0.02 ug/g dry	<0.02	0.02	0.04	-
Anthracene	0.02 ug/g dry	<0.02	0.07	0.09	-
Benzo [a] anthracene	0.02 ug/g dry	0.05	0.16	0.36	-
Benzo [a] pyrene	0.02 ug/g dry	0.18	0.16	0.41	-
Benzo [b] fluoranthene	0.02 ug/g dry	0.50	0.18 0.79		-
Benzo [g,h,i] perylene	0.02 ug/g dry	0.22	0.10	0.34	-



Order #: 2043350

Report Date: 26-Oct-2020 Order Date: 20-Oct-2020

Project Description: PE4999

	-				
	Client ID:	TP12-G1	TP13-G1	TP14-G1	-
	Sample Date:	19-Oct-20 09:00	19-Oct-20 09:00	19-Oct-20 09:00	-
	Sample ID:	2043350-09	2043350-10	2043350-11	-
	MDL/Units	Soil	Soil	Soil	-
Benzo [k] fluoranthene	0.02 ug/g dry	0.14	0.09	0.44	-
Chrysene	0.02 ug/g dry	0.06	0.16	0.42	-
Dibenzo [a,h] anthracene	0.02 ug/g dry	0.04	0.03	0.08	-
Fluoranthene	0.02 ug/g dry	0.11	0.35	0.84	-
Fluorene	0.02 ug/g dry	<0.02	<0.02	<0.02	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	0.13	0.09	0.29	-
1-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.02	-
2-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.02	-
Methylnaphthalene (1&2)	0.04 ug/g dry	<0.04	<0.04	<0.04	-
Naphthalene	0.01 ug/g dry	<0.01	<0.01	<0.01	-
Phenanthrene	0.02 ug/g dry	0.06	0.17	0.29	-
Pyrene	0.02 ug/g dry	0.16	0.27	0.67	-
2-Fluorobiphenyl	Surrogate	103%	71.1%	107%	-
Terphenyl-d14	Surrogate	107%	112%	110%	-



## Method Quality Control: Blank

Report Date: 26-Oct-2020

Order Date: 20-Oct-2020

Project Description: PE4999

		Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Hydrocarbons									
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			-3,3						
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			-3.3						
Acenaphthene	ND	0.02	ug/g						
Acenaphthylene	ND	0.02	ug/g ug/g						
Anthracene	ND	0.02	ug/g ug/g						
Benzo [a] anthracene	ND	0.02	ug/g ug/g						
Benzo [a] pyrene	ND	0.02	ug/g ug/g						
Benzo [b] fluoranthene	ND	0.02	ug/g ug/g						
Benzo [g,h,i] perylene	ND	0.02	ug/g 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 ug/g						
Indeno [1,2,3-cd] pyrene	ND	0.02							
1-Methylnaphthalene	ND	0.02	ug/g						
2-Methylnaphthalene	ND	0.02	ug/g						
	ND	0.02	ug/g						
Methylnaphthalene (1&2) Naphthalene	ND ND	0.04 0.01	ug/g						
Phenanthrene	ND	0.01	ug/g						
	ND ND	0.02	ug/g						
Pyrene Surrogate: 2 Eluorobiohenyl	0.869	0.02	ug/g		65.1	50-140			
Surrogate: 2-Fluorobiphenyl			ug/g						
Surrogate: Terphenyl-d14	1.59		ug/g		119	50-140			



Client PO: 31420

Report Date: 26-Oct-2020

Order Date: 20-Oct-2020

Project Description: PE4999

## Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
-			Gillo	rtesuit	, or CEO	Laint		Lurint	
Hydrocarbons									
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	ND	1.0	ug/g dry	ND			NC	30	
Arsenic	3.8	1.0	ug/g dry	3.5			6.2	30	
Barium	170	1.0	ug/g dry	152			11.0	30	
Beryllium	0.7	0.5	ug/g dry	0.6			18.8	30	
Boron	6.1	5.0	ug/g dry	ND			NC	30	
Cadmium	ND	0.5	ug/g dry	ND			NC	30	
Chromium (VI)	ND	0.2	ug/g dry	ND			NC	35	
Chromium	77.6	5.0	ug/g dry	70.0			10.3	30	
Cobalt	14.6	1.0	ug/g dry	13.3			8.7	30	
Copper	26.2	5.0	ug/g dry	23.5			10.8	30	
Lead	12.2	1.0	ug/g dry	10.8			11.6	30	
Mercury	ND	0.1	ug/g dry	ND			NC	30	
Molybdenum	ND	1.0	ug/g dry	ND			NC	30	
Nickel	37.7	5.0	ug/g dry	34.2			9.9	30	
Selenium	ND	1.0	ug/g dry	ND			NC	30	
Silver	ND	0.3	ug/g dry	ND			NC	30	
Thallium	ND	1.0	ug/g dry	ND			NC	30	
Uranium	2.0	1.0	ug/g dry	1.7			16.0	30	
Vanadium	66.5	10.0	ug/g dry	59.9			10.5	30	
Zinc	86.0	20.0	ug/g dry	77.4			10.4	30	
Physical Characteristics									
% Solids	84.7	0.1	% by Wt.	84.6			0.0	25	
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g dry	ND			NC	40	
Acenaphthylene	ND	0.02	ug/g dry	ND			NC	40	
Anthracene	0.023	0.02	ug/g dry	ND			NC	40	
Benzo [a] anthracene	0.036	0.02	ug/g dry	0.028			22.8	40	
Benzo [a] pyrene	0.037	0.02	ug/g dry	0.028			27.5	40	
Benzo [b] fluoranthene	0.046	0.02	ug/g dry	0.036			25.6	40	
Benzo [g,h,i] perylene	0.028	0.02	ug/g dry	ND			NC	40	
Benzo [k] fluoranthene	0.020	0.02	ug/g dry	ND			NC	40	
Chrysene	0.044	0.02	ug/g dry	0.030			37.0	40	
Dibenzo [a,h] anthracene	ND	0.02	ug/g dry	ND			NC	40	
Fluoranthene	0.097	0.02	ug/g dry	0.079			21.3	40	
Fluorene	ND	0.02	ug/g dry	ND			NC	40	
Indeno [1,2,3-cd] pyrene	0.023	0.02	ug/g dry	ND			NC	40	
1-Methylnaphthalene	ND	0.02	ug/g dry	ND			NC	40	
2-Methylnaphthalene	ND	0.02	ug/g dry	ND			NC	40	
Naphthalene	0.021	0.01	ug/g dry	0.013			NC	40	
Phenanthrene	0.091	0.02	ug/g dry	0.059			NC	40	
Pyrene	0.078	0.02	ug/g dry	0.062			22.4	40	
Surrogate: 2-Fluorobiphenyl	0.852		ug/g dry		55.6	50-140			
Surrogate: Terphenyl-d14	1.47		ug/g dry		95.7	50-140			



# Method Quality Control: Spike

Report Date: 26-Oct-2020

Order Date: 20-Oct-2020

Project Description: PE4999

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F2 PHCs (C10-C16)	110	4	ug/g	ND	107	60-140			
F3 PHCs (C16-C34)	280	8	ug/g	ND	111	60-140			
F4 PHCs (C34-C50)	175	6	ug/g	ND	109	60-140			
F4G PHCs (gravimetric)	980	50	ug/g	ND	98.0	80-120			
Metals									
Antimony	40.3	1.0	ug/g	ND	80.3	70-130			
Arsenic	47.6	1.0	ug/g	1.4	92.3	70-130			
Barium	109	1.0	ug/g	60.9	95.3	70-130			
Beryllium	43.0	0.5	ug/g	ND	85.5	70-130			
Boron	38.0	5.0	ug/g	ND	72.1	70-130			
Cadmium	42.6	0.5	ug/g	ND	85.0	70-130			
Chromium (VI)	4.1	0.2	ug/g	ND	81.5	70-130			
Chromium	75.7	5.0	ug/g	28.0	95.4	70-130			
Cobalt	50.3	1.0	ug/g	5.3	89.9	70-130			
Copper	53.2	5.0	ug/g	9.4	87.6	70-130			
Lead	46.0	1.0	ug/g	4.3	83.3	70-130			
Mercury	1.66	0.1	ug/g	ND	111	70-130			
Molybdenum	42.4	1.0	ug/g	ND	84.1	70-130			
Nickel	57.6	5.0	ug/g	13.7	87.9	70-130			
Selenium	44.1	1.0	ug/g	ND	87.8	70-130			
Silver	36.5	0.3	ug/g	ND	73.0	70-130			
Thallium	44.7	1.0	ug/g	ND	89.3	70-130			
Uranium	46.0	1.0	ug/g	ND	90.6	70-130			
Vanadium	71.3	10.0	ug/g	23.9	94.7	70-130			
Zinc	74.7	20.0	ug/g	31.0	87.5	70-130			
Semi-Volatiles									
Acenaphthene	0.201	0.02	ug/g	ND	105	50-140			
Acenaphthylene	0.154	0.02	ug/g	ND	80.6	50-140			
Anthracene	0.265	0.02	ug/g	ND	138	50-140			
Benzo [a] anthracene	0.275	0.02	ug/g	0.028	128	50-140			
Benzo [a] pyrene	0.294	0.02	ug/g	0.028	139	50-140			
Benzo [b] fluoranthene	0.294	0.02	ug/g	0.036	135	50-140			
Benzo [g,h,i] perylene	0.218	0.02	ug/g	ND	114	50-140			
Benzo [k] fluoranthene	0.265	0.02	ug/g	ND	139	50-140			
Chrysene	0.292	0.02	ug/g	0.030	136	50-140			
Dibenzo [a,h] anthracene	0.175	0.02	ug/g	ND	91.3	50-140			
Fluoranthene	0.112	0.02	ug/g	ND	67.1	50-140			
Fluorene	0.213	0.02	ug/g	ND	111	50-140			
Indeno [1,2,3-cd] pyrene	0.220	0.02	ug/g	ND	115	50-140			
1-Methylnaphthalene	0.120	0.02	ug/g	ND	62.7	50-140			
2-Methylnaphthalene	0.136	0.02	ug/g	ND	71.2	50-140			
Naphthalene	0.181	0.01	ug/g	0.013	87.6	50-140			
Phenanthrene	0.109	0.02	ug/g	ND	65.2	50-140			
Pyrene	0.108	0.02	ug/g	ND	65.0	50-140			
Surrogate: 2-Fluorobiphenyl	0.829		ug/g		54.1	50-140			
Surrogate: Terphenyl-d14	1.66		ug/g		108	50-140			



Sample Qualifiers :

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

#### QC Qualifiers :

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.

GPARACEL				2043350	m		(Lab	Order I Use O		r		(L	n Of C ab Use 550		y
Client Name: Paterson Group		Projec	t Ref:	2E11999							,		age _	of Z	
Contact Name: Mick Sullivan		Quote										Turi	naroun	d Time	
Address:		PO #:	31	120							□ 1 d	lay			3 day
154 Colonnade Rd. S.		E-mail									□ 2 d	lay			Regular
Telephone: 613-226 + 7381			ns	ullvanes	polycon	<u>مرمر</u>	Q. (	a			Date Re	quired	:		
Regulation 153/04 Other Regulation						1	1								
Table 1 Res/Park Med/Fine REG 558 PWQO				(Soil/Sed.) GW (Gr /ater) SS (Storm/Sar						Re	quired Ar	nalysis			
Table 2 Ind/Comm Coarse CCME MISA		(		aint) A (Air) O (Oth			1	13	3	5		Τ	1		
Table 3 Agri/Other USU-Sani USU-Sto	rm	1	s.			4		1	2	+ 2					
Table Mun:		e	Containers	Sample	Taken	L.	Ľ,			0.0					
For RSC: Yes No Other:	- xi	Air Volume	Cont				2		West in	hremiu.					
Sample ID/Location Name	Matrix	Air	# of	Date	Time	2		,5	Ś	5					
1 TP1-G2	5		1	oct. 19.20	AM	×	X	4	V			1			
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3 TP4-GZ			1												
4 195-62			1			$\uparrow \uparrow$		$\square$	Ħ				1		
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7 TP10-G1	+		1			++	$\ddagger$					+-	1		
8 TR 11 - GI			1			++		Ħ				+	+		<b>—</b> ,
9 TP12-G1	++		J			++	$\square$	+	+			+			-
10 TP13-G1				V	V	V	V		V	V		+	+		
Comments:							V	V	<u> </u>	Metho	d of Delive	ry: CEL	. 4	NAR	ž
Relinquished By (Sign): N- Surfinvenn Received B	y Driver/D	epot:	D	0 3 25	Refeived at Lab;	ovn		Doh	maj	Verifie	d By:	St	AM		
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Date/Time: Oct. ZO. 20 Temperatu	re:	1		°C PH.	Temperature:	10	4 °(	C			ified: 🗌		γ:		

Chain of Custody (Blank) xlsx

Revision 3.0

C PARACEL III				2043350			acel On (Lab U	se Onl	y)		C Nº	Stor Bar	of C buse (	Only)	dy	
Client Name: Paterson Group		Project	Ref:	PE4999								Pa	nge 🔼	of Z	,	
Contact Name: Nick Sulliver		Quote	#:									Turna	aroun	d Tim	e	
Address:		PO #: •	311	120							🗆 1 da	зу			🗆 3 da	ıy
154 Colongade Rd. S.		E-mail:								_	🗆 2 da	ay .			Reg	ula
154 Colonnade Rd. S. Telephone: 613-226-7381		0	Sul	Ivan @ patersor	201		C.a.			D	ate Req	uired:				
Regulation 153/04 Other Regulation					0	T			2							10
Table 1 Res/Park Med/Fine REG 558 PWQ0				(Soil/Sed.) GW (Ground Wa /ater) SS (Storm/Sanitary Sew						Requ	ired An	alysis				
Table 2 Ind/Comm Coarse CCME MISA				aint) A (Air) O (Other)				1		<		Τ				
Table 3 🗆 Agri/Other			S			E.		TC	1	2						
Table Mun:		e	taine	Sample Taken		1,0	AHS	$\sim$	5	in m						
For RSC: Yes No Other:	,iz	Air Volume	of Containers			5	×	Mehals	Nereve	chremium			1.1			
Sample ID/Location Name	Matrix	Air	to #	Date Tim	ne	PHCs	<u></u>	Ne	Z	C.P.						
1 JP14 - GI	S		1	004.19.20 A	И	X	X	r	6	$\mathbf{x}$						
2																
3							-									
4	$\square$	-													1	
5								1				1				
6									+	-	-				-	
7						$\square$		-		+		+			-	
8	$\left  \right $							-	-	+	+-				-	
9										+	+	1				
10								+								
Comments:						-			M	lethod o	f Deliver	I.	4	1 auc	EC	
Relinquished By (Sign):	1	/	Tu.	Received a	eph	rM	(	bla	mai	erified B	γ:	36	2m			
Relinquished By (Print): Nich Sufficien Date/Time:	20/1	10	12	325 Dater Jime	V . d	HPO		R	1. ]	ate/Tim	00		1,2020	0	13:30	0
Date/Time: Oct. 20, 20 Temperature		-		°C PH. Temperatu	ure: 1	10	/ ) °C		pł	H Verifie	2d: 🛄	By:				

Chain of Custody (Blank) xlsx

Revision 3.0



RELIABLE.

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# Certificate of Analysis

# **Paterson Group Consulting Engineers**

154 Colonnade Road South Nepean, ON K2E 7J5 Attn: Nick Sullivan

Client PO: 31397 Project: PE4999 Custody: 116611

Report Date: 8-Oct-2020 Order Date: 5-Oct-2020

Order #: 2041094

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

Paracel ID	Client ID
2041094-01	BH2-GW1
2041094-02	BH4-GW1
2041094-03	BH5-GW1
2041094-04	DUP1

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: 08-Oct-2020 Order Date: 5-Oct-2020

Order #: 2041094

Project Description: PE4999

## **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 624 - P&T GC-MS	6-Oct-20	6-Oct-20
PHC F1	CWS Tier 1 - P&T GC-FID	6-Oct-20	6-Oct-20
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	7-Oct-20	7-Oct-20
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	6-Oct-20	6-Oct-20



Client PO: 31397

Report Date: 08-Oct-2020 Order Date: 5-Oct-2020

Project Description: PE4999

ſ	Client ID: Sample Date: Sample ID: MDL/Units	BH2-GW1 02-Oct-20 12:00 2041094-01 Water	BH4-GW1 02-Oct-20 12:00 2041094-02 Water	BH5-GW1 02-Oct-20 12:00 2041094-03 Water	DUP1 02-Oct-20 12:00 2041094-04 Water
Volatiles	i		i		
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 (dibromoethane, 1,2-)	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



Report Date: 08-Oct-2020 Order Date: 5-Oct-2020

Project Description: PE4999

	Client ID: Sample Date:	BH2-GW1 02-Oct-20 12:00 2041094-01	BH4-GW1 02-Oct-20 12:00 2041094-02	BH5-GW1 02-Oct-20 12:00 2041094-03	DUP1 02-Oct-20 12:00 2041094-04
	Sample ID: MDL/Units	Water	204 1094-02 Water	2041094-03 Water	204 1094-04 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	-	97.0%	97.0%	96.3%
Dibromofluoromethane	Surrogate	-	98.1%	98.2%	79.0%
Toluene-d8	Surrogate	-	104%	105%	106%
Benzene	0.5 ug/L	<0.5	-	-	-
Ethylbenzene	0.5 ug/L	<0.5	-	-	-
Toluene	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	-	-	-
Toluene-d8	Surrogate	105%	-	-	-
Hydrocarbons					•
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	-



Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 31397

## Method Quality Control: Blank

Report Date: 08-Oct-2020

Order Date: 5-Oct-2020

Project Description: PE4999

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	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 Methyl Ethyl Ketone (2-Butanone)	ND ND	1.0 5.0	ug/L						
Methyl Isobutyl Ketone	ND	5.0	ug/L						
Methyl tert-butyl ether	ND	2.0	ug/L ug/L						
Methylene Chloride	ND	5.0	ug/L						
Styrene	ND	0.5	ug/L						
1,1,1,2-Tetrachloroethane	ND	0.5	-						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L 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	77.9	0.5	ug/L		97.4	50-140			
Surrogate: Dibromofluoromethane	76.5		ug/L		95.6	50-140 50-140			
-			-						
Surrogate: Toluene-d8	84.3	0.5	ug/L		105	50-140			
Benzene	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
		0.5							
Xylenes, total <i>Surrogate: Toluene-d8</i>	ND 84.3	0.5	ug/L ug/L		105	50-140			



# Method Quality Control: Duplicate

Report Date: 08-Oct-2020 Order Date: 5-Oct-2020

Project Description: PE4999

Analyté         Result         Limit         Units         Result         S,REC         Limit         RPD         Limit         Notes           Hydrocarbons         F1 FHGs (CS-01)         ND         25         ug/L         ND         NC         30           F1 FHGs (CS-01)         ND         5.0         ug/L         ND         NC         30           Adefone         ND         0.5         ug/L         ND         NC         30           Branzen         ND         0.5         ug/L         ND         NC         30           Gromonethane         ND         0.5         ug/L         ND         NC         30           Gromonethane         ND         0.5         ug/L         ND         NC         30           Carbon Ferachioride         ND         0.5         ug/L         ND         NC         30           Chiorobarzene         ND         0.5         ug/L         ND         NC         30           Diskinordifusormethane         ND         0.5         ug/L         ND         NC         30           Labelintorbarzene         ND         0.5         ug/L         ND         NC         30           Labe			Reporting		Source		%REC		RPD	
r F1P(b2) (Cb-C10)ND25ug/LNDNC30VolticeValueBarszneND5.0ug/LNDNC30BarszneND0.5ug/LNDNC30BromadicioronthaneND0.5ug/LNDNC30BromadicioronthaneND0.5ug/LNDNC30BromadicioronthaneND0.5ug/LNDNC30BromadicioronthaneND0.5ug/LNDNC30ChioroformND0.5ug/LNDNC30DibrondicioronthaneND0.5ug/LNDNC30DibrondicioronthaneND0.5ug/LNDNC301/2.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3.DichioroberzeneND0.5ug/LNDNC301/3	Analyte	Result		Units		%REC		RPD		Notes
r F1P(b2) (Cb-C10)ND25ug/LNDNC30VolumeND5.0ug/LNDNC30BarnzmeND0.5ug/LNDNC30Barnzme/InternationND0.5ug/LNDNC30Bromodic/incomblancND0.5ug/LNDNC30Bromodic/incomblancND0.5ug/LNDNC30Bromodic/incomblancND0.5ug/LNDNC30Bromone/InternationND0.5ug/LNDNC30ChickotombrazeneND0.5ug/LNDNC30ChickotombrazeneND0.5ug/LNDNC301/2.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5ug/LNDNC301/3.DickitordenzeneND0.5 <td>Hydrocarbons</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	Hydrocarbons									
Volatilies         ND         S.0         ugL         ND         NC         30           Bernarde         ND         0.5         ugL         ND         NC         30           Bromadium         ND         0.5         ugL         ND         NC         30           Bromadium         ND         0.5         ugL         ND         NC         30           Bromadium         ND         0.5         ugL         ND         NC         30           Chirotherazone         ND         0.5         ugL         ND         NC         30           Dichtorotherane         ND         0.5         ugL         ND         NC         30           Dichtorotherane         ND         0.5         ugL         ND         NC         30           1.3-Dichtorotherane         ND         0.5         ugL         ND         NC         30           1.3-Dichtorotherane         ND         0.5         ugL         ND         NC         30           1.3-Dichtorothylene         ND         0.5         ugL         ND         NC         30           1.3-Dichtorothylene         ND         0.5         ugL         ND         NC	•	ND	25	ua/l	ND			NC	30	
ActoneND5.0ug/LNDNDNC30BenenderkhonenhaneND0.5ug/LNDNC30BronenderkhonenhaneND0.5ug/LNDNC30BronenderkhonenhaneND0.5ug/LNDNC30BronenderkhonenhaneND0.5ug/LNDNC30Cathon TetrobardenND0.5ug/LNDNC30Cathon TetrobardenND0.5ug/LNDNC30DibronenhanenenhaneND0.5ug/LNDNC301/2 OlchonebrazeneND0.5ug/LNDNC301/3 OlchonebrazeneND <td></td> <td>ne -</td> <td>20</td> <td>ug/L</td> <td>ne -</td> <td></td> <td></td> <td>110</td> <td>00</td> <td></td>		ne -	20	ug/L	ne -			110	00	
BenzeneND0.5up/LNDNDNC30BromodifinomethaneND0.5up/LNDNC30BromodifinomethaneND0.5up/LNDNC30Carbon TetrachlorideND0.5up/LNDNC30ChioroberzeneND0.5up/LNDNC30ChioroberzeneND0.5up/LNDNC30DichrondfloromethaneND0.5up/LNDNC30J.2-OchricoberzeneND0.5up/LNDNC30J.3-DichrondfloromethaneND0.5up/LNDNC30J.4-DichronderzeneND0.5up/LNDNC30J.2-DichronderzeneND0.5up/LNDNC30J.1-DichronderzeneND0.5up/LNDNC30J.1-DichronderyeneND0.5up/LNDNC30J.1-DichronderyeneND0.5up/LNDNC30J.1-DichronderyeneND0.5up/LNDNC30J.1-DichronderyeneND0.5up/LNDNC30L-1-DichronderyeneND0.5up/LNDNC30L-1-DichronderyeneND0.5up/LNDNC30L-1-DichronderyeneND0.5up/LNDNC30L-1-DichronderyeneND<			FO					NC	20	
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Carbon TetrachlorideND0.2ug/LNDNC30ChlorobernzeneND0.5ug/LNDNC30DichorochloromethaneND0.5ug/LNDNC30DichorochloromethaneND0.5ug/LNDNC301.3-DichlorobenzeneND0.5ug/LNDNC301.3-DichlorobenzeneND0.5ug/LNDNC301.4-DichlorobenzeneND0.5ug/LNDNC301.4-DichlorobenzeneND0.5ug/LNDNC301.4-DichlorobenzeneND0.5ug/LNDNC301.4-DichlorobethaneND0.5ug/LNDNC301.4-DichlorobethaneND0.5ug/LNDNC301.4-DichlorobethyleneND0.5ug/LNDNC301.4-DichlorobethyleneND0.5ug/LNDNC301.4-DichlorobethyleneND0.5ug/LNDNC301.4-DichlorophyleneND0.5ug/LNDNC301.4-DichlorobethyleneND0.5ug/LNDNC301.4-DichlorobethyleneND0.5ug/LNDNC301.4-DichlorobethyleneND0.5ug/LNDNC301.4-DichlorobethyleneND0.5ug/LNDNC301.4-										
ChicolenzaneND0.5ug/LNDNC30OhrondermachiaromethaneND0.5ug/LNDNC30DicharonationomethaneND0.5ug/LNDNC301.2-DicharobenzeneND0.5ug/LNDNC301.3-DicharobenzeneND0.5ug/LNDNC301.4-DicharobenzeneND0.5ug/LNDNC301.1-DicharonethaneND0.5ug/LNDNC301.1-DicharonethaneND0.5ug/LNDNC301.1-DicharonethaneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND0.5ug/LNDNC301.1-DicharonethyleneND5.0ug/LNDNC301.1-Dicharone				-						
ChlorodrinomethaneND0.5ug/LNDNC30DichonochtomethaneND0.5ug/LNDNC301.2-DichorobenzeneND0.5ug/LNDNC301.4-DichorobenzeneND0.5ug/LNDNC301.4-DichorobenzeneND0.5ug/LNDNC301.4-DichorobenzeneND0.5ug/LNDNC301.2-DichorobenzeneND0.5ug/LNDNC301.2-DichoroethaneND0.5ug/LNDNC301.2-DichoroethyleneND0.5ug/LNDNC301.2-DichoroethyleneND0.5ug/LNDNC301.2-DichoroethyleneND0.5ug/LNDNC301.2-DichoroptyleneND0.5ug/LNDNC301.2-DichoroptyleneND0.5ug/LNDNC301.2-DichoroptyleneND0.5ug/LNDNC301.2-DichoroptyleneND0.5ug/LNDNC301.2-DichoroptyleneND0.5ug/LNDNC301.2-DichoroptyleneND0.5ug/LNDNC301.2-DichoroptyleneND0.5ug/LNDNC301.2-DichoroptyleneND5.0ug/LNDNC301.2-DichoroptyleneND <td></td> <td></td> <td></td> <td>-</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>				-						
DibronchloromethaneND0.5ug/LNDNC301.2-DichlorobenzeneND0.5ug/LNDNC301.3-DichlorobenzeneND0.5ug/LNDNC301.3-DichlorobenzeneND0.5ug/LNDNC301.4-DichlorobenzeneND0.5ug/LNDNC301.1-DichlorobenzeneND0.5ug/LNDNC301.2-DichlorobenzeneND0.5ug/LNDNC301.2-DichlorobethyleneND0.5ug/LNDNC301.1-DichlorobethyleneND0.5ug/LNDNC301.2-DichlorobethyleneND0.5ug/LNDNC301.2-DichlorophyleneND0.5ug/LNDNC301.2-DichlorophyleneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30ItarasitionoblaneND0.5ug/L<				-						
DichlorodifluorometaneND1.0ug/LNDNC3.01.2-DichlorobenzeneND0.5ug/LNDNC301.3-DichlorobenzeneND0.5ug/LNDNC301.4-DichlorobenzeneND0.5ug/LNDNC301.2-DichloroethaneND0.5ug/LNDNC301.2-DichloroethaneND0.5ug/LNDNC301.2-DichloroethyleneND0.5ug/LNDNC301.2-DichloroethyleneND0.5ug/LNDNC301.2-DichloroethyleneND0.5ug/LNDNC301.2-DichloroethyleneND0.5ug/LNDNC301.2-DichloropropyleneND0.5ug/LNDNC301.2-DichloropropyleneND0.5ug/LNDNC301.2-DichloropropyleneND0.5ug/LNDNC30Ethylene dibromide (dibromoethane, 1,2ND0.5ug/LNDNC30Methyl (stone (2-Bulanone)ND5.0ug/LNDNC30Methyl (stone (2-Bulanone)ND0.5ug/LNDNC30Methyl (stone (2-Bulanone)ND0.5ug/LNDNC301.1.2.2-TietachioroethaneND0.5ug/LNDNC301.1.1.2.2-TietachioroethaneND0.5ug/L<				-						
1.2-DichloroberzeneND0.5ug/LNDNDNC301.3-DichloroberzeneND0.5ug/LNDNC301.4-DichloroberzeneND0.5ug/LNDNC301.1-DichloroberzeneND0.5ug/LNDNC301.1-DichloroberzeneND0.5ug/LNDNC301.1-DichloroberzeneND0.5ug/LNDNC301.1-DichloroberzeneND0.5ug/LNDNC301.1-DichloroberzeneND0.5ug/LNDNC301.1-DichloroberzeneND0.5ug/LNDNC301.2-DichloropopaneND0.5ug/LNDNC30Ethylene dibromole (dibromoethane, 1.2ND0.5ug/LNDNC30Ethylene dibromide (dibromoethane, 1.2ND0.5ug/LNDNC30HexaneND0.5ug/LNDNC30Hethyle Ethyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl terbulyl etherND0.5ug/LNDNC30StyreneND0.5ug/LNDNC301,1,1.2-TickhorobethaneND0.5ug/LNDNC301,1,2-TickhorobethaneND0.5ug/LNDNC301,1,2-TickhorobethaneND0.5ug/LND				-						
1,3-DichlorobenzeneND0.5ug/LNDNDNC301,4-DichlorobenzeneND0.5ug/LNDNC301,2-DichloroethaneND0.5ug/LNDNC301,2-DichloroethyleneND0.5ug/LNDNC301,2-DichloroethyleneND0.5ug/LNDNC301,2-DichloroptropeneND0.5ug/LNDNC301,2-DichloroptropeneND0.5ug/LNDNC301,2-DichloroptropeneND0.5ug/LNDNC301,2-DichloroptropeneND0.5ug/LNDNC301,3-DichloroptropeneND0.5ug/LNDNC301,3-DichloroptropeneND0.5ug/LNDNC30Ethylene dibromide (dibromoethane, 1,2ND0.5ug/LNDNC30HexaneND5.0ug/LNDNC30Hethyl Ethyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl Ethyl ketoneND5.0ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/L<				-						
1.4-DichlorobenzeneND0.5ug/LNDNDNC301,1-DichloroethaneND0.5ug/LNDNC301,2-DichloroethyleneND0.5ug/LNDNC30(5:1.2-DichloroethyleneND0.5ug/LNDNC30(5:1.2-DichloroethyleneND0.5ug/LNDNC30(5:1.2-DichloroethyleneND0.5ug/LNDNC30(1.2-DichloroethyleneND0.5ug/LNDNC30(1.2-DichloroethyleneND0.5ug/LNDNC30(1.2-DichloroethyleneND0.5ug/LNDNC30(1.2-Dichloroethane, 1.2ND0.5ug/LNDNC30Ethylene dibromide (dibromoethane, 1.2ND0.5ug/LNDNC30Ethylene dibromide (dibromoethane, 1.2ND0.5ug/LNDNC30Methyl Ethyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl Ethyl KetoneND5.0ug/LNDNC30Methyl Ethyl KetoneND5.0ug/LNDNC301,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1-TrichoroethaneND0.5ug/LNDNC301,1,1-TrichoroethaneND0.5ug/LNDNC301,1,1-TrichoroethaneND	,			-						
1.1-DichloroethaneND0.5ug/LNDNDNC301.2-DichloroethyleneND0.5ug/LNDNC30cis 1.2-DichloroethyleneND0.5ug/LNDNC30cis 1.2-DichloroethyleneND0.5ug/LNDNC30trans-1.2-DichloroethyleneND0.5ug/LNDNC301.2-DichloropropaneND0.5ug/LNDNC30cis 1.3-DichloropropyleneND0.5ug/LNDNC30trans-1.2-DichloropropyleneND0.5ug/LNDNC30EthylencaND0.5ug/LNDNC30HexaneND0.5ug/LNDNC30HexaneND1.0ug/LNDNC30Hethyl felotoutyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl fel-thyl felotoutyl KetoneND5.0ug/LNDNC30Methyl fel-thyl felotoutyl KetoneND5.0ug/LNDNC301,1.2-TetrachloroethaneND0.5ug/LNDNC301,1.1.2-TetrachloroethaneND0.5ug/LNDNC301,1.1.2-TetrachloroethaneND0.5ug/LNDNC301,1.1.2-TetrachloroethaneND0.5ug/LNDNC301,1.1.2-TetrachloroethaneND0.5				-						
1.2-DichlorosethaneND0.5ug/LNDNC301.1-DichlorosethyleneND0.5ug/LNDNC30cis.1.2-DichlorosethyleneND0.5ug/LNDNC30cis.1.2-DichlorosethyleneND0.5ug/LNDNC30cis.1.3-DichlorosethyleneND0.5ug/LNDNC30cis.1.3-DichlorosethyleneND0.5ug/LNDNC30cis.1.3-Dichlorosethane,1.2ND0.5ug/LNDNC30Ethylene dibronide (dibromoethane, 1.2ND0.5ug/LNDNC30Ethylene dibronide (dibromoethane, 1.2ND0.5ug/LNDNC30Methyl Ethyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl Isbulyl KetoneND5.0ug/LNDNC30Methyl Isbulyl KetoneND5.0ug/LNDNC30StyreneND0.5ug/LNDNC30Methyl Isbulyl KetoneND5.0ug/LNDNC30StyreneND0.5ug/LNDNC30Methyl Isbulyl KetoneND5.0ug/LNDNC301.1,1.2-TetrachlorosethaneND0.5ug/LNDNC301.1,1.2-TetrachlorosethaneND0.5ug/LNDNC301.1,1.2-TetrachlorosethaneND </td <td></td> <td></td> <td></td> <td>-</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>				-						
1.1 DicklarosethyleneND0.5ug/LNDNC30cis-1.2-DicklarosethyleneND0.5ug/LNDNC301.2-DicklarosethyleneND0.5ug/LNDNC301.2-DicklarospropaneND0.5ug/LNDNC301:1-3-DicklarospropaneND0.5ug/LNDNC30tis-1.3-DicklarospropyleneND0.5ug/LNDNC30Ethylene diromoethane, 1.2ND0.5ug/LNDNC30HexaneND0.5ug/LNDNC30Hethyl Kotone (2-Butanone)ND5.0ug/LNDNC30Methyl Ethyl kotone (2-Butanone)ND5.0ug/LNDNC30Methyl Isobutyl KotoneND5.0ug/LNDNC30Methyl Isobutyl KotoneND5.0ug/LNDNC30Methyl Isobutyl KotoneND5.0ug/LNDNC30StyreneND0.5ug/LNDNC301,1,1.2-TeitachlorosethaneND0.5ug/LNDNC301,1,1.2-TeitachlorosethaneND0.5ug/LNDNC301,1,1.2-TeitachlorosethaneND0.5ug/LNDNC301,1,1.2-TeitachlorosethaneND0.5ug/LNDNC301,1,1.2-TeitachlorosethaneND0.5ug/L	,			-						
cis-12-DichloroethyleneND0.5ug/LNDNDNC30trans-12-DichloroethyleneND0.5ug/LNDNC3012-DichloropropaneND0.5ug/LNDNC30cis-13-DichloropropyleneND0.5ug/LNDNC30trans-13-DichloropropyleneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30EthylbenzeneND0.5ug/LNDNC30HexaneND1.0ug/LNDNC30Methyl Ethyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl Istrbulyl KetoneND5.0ug/LNDNC30Methyl Istrbulyl KetoneND5.0ug/LNDNC30Methyl Istrbulyl KetoneND5.0ug/LNDNC30SyreneND0.5ug/LNDNC301,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,1,2-TetrachloroethaneND0.5ug/LND <t< td=""><td></td><td></td><td></td><td>-</td><td></td><td></td><td></td><td></td><td></td><td></td></t<>				-						
trans.12-DichloredhyleneND0.5ug/LNDNDNC301,2-DichloropropaneND0.5ug/LNDNC30cis1.3-DichloropropyleneND0.5ug/LNDNC30trans.13-DichloropropyleneND0.5ug/LNDNC30EthylenzeneND0.5ug/LNDNC30EthylenzeneND0.5ug/LNDNC30Methyl Ethyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl Isobutyl KetoneND5.0ug/LNDNC30Methyl Isobutyl KetoneND5.0ug/LNDNC30Methyl Isobutyl KetoneND5.0ug/LNDNC30Methyl Isobutyl KetoneND5.0ug/LNDNC30StyreneND0.5ug/LNDNC301,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,2-TetrachloroethaneND0.5ug/LNDNC301,1,2-TichloroethaneND0.5ug/LNDNC301,1,1-TichloroethaneND0.5ug/LNDNC301,1,1-TichloroethaneND0.5ug/LNDNC301,1,1-TichloroethaneND0.5ug/LNDNC301,1,1-TichloroethaneND0.5ug/LNDNC <t< td=""><td></td><td></td><td></td><td>-</td><td></td><td></td><td></td><td></td><td></td><td></td></t<>				-						
1.2-Dichloropropane       ND       0.5       ug/L       ND       <				-						
cis-1,3-Dichloropropylene         ND         0.5         ug/L         ND         NC         30           trans-1,3-Dichloropropylene         ND         0.5         ug/L         ND         NC         30           Ethylenzene         ND         0.5         ug/L         ND         NC         30           Ethylenzene         ND         0.2         ug/L         ND         NC         30           Methyl Ethyl Ketone (2-Butanone)         ND         5.0         ug/L         ND         NC         30           Methyl Ethyl Ketone (2-Butanone)         ND         5.0         ug/L         ND         NC         30           Methyl Isrbutyl ether         ND         5.0         ug/L         ND         NC         30           Methyl Isrbutyl ether         ND         5.0         ug/L         ND         NC         30           Styrene         ND         0.5         ug/L         ND         NC         30           1,1,2-Tetrachoroethane         ND         0.5         ug/L         ND         NC         30           1,1,2-Tetrachoroethane         ND         0.5         ug/L         ND         NC         30           1,1,2-Tetrachoroethane				-						
trans-1.3-Dichloropropylene         ND         0.5         ug/L         ND         NC         30           Ethylpenzene         ND         0.5         ug/L         ND         NC         30           Ethylpenzene         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 terbulyl Ketone         ND         5.0         ug/L         ND         NC         30           Methyl terbulyl				-						
Ethylbenzene         ND         0.5         ug/L         ND         NC         30           Ethylbene 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 Ketone         ND         5.0         ug/L         ND         NC         30           Methyl Ketone         ND         5.0         ug/L         ND         NC         30           Methyl Ketone         ND         5.0         ug/L         ND         NC         30           Methyl Ethyl Ketone         ND         5.0         ug/L         ND         NC         30           Methyl Ethyl Ketone         ND         0.5         ug/L         ND         NC         30           Styrene         ND         0.5         ug/L         ND         NC         30           1,1,2.7 Ethtachloroethane         ND         0.5         ug/L         ND         NC         30           1,1,2.7 Ethtachloroethane         ND         0.				•						
Ethylene dibromide (dibromoethane, 1,2:         ND         0.2         ug/L         ND         ND<				-						
HexaneND1.0ug/LNDNDNC30Methyl Ethyl Ketone (2-Butanone)ND5.0ug/LNDNC30Methyl Bobulyl KetoneND5.0ug/LNDNC30Methyl Letr-butyl etherND2.0ug/LNDNC30Methyl Letr-butyl etherND5.0ug/LNDNC30Methyl Letr-butyl etherND0.5ug/LNDNC30StyreneND0.5ug/LNDNC301,1,2.7etrachloroethaneND0.5ug/LNDNC301,1,2.7etrachloroethaneND0.5ug/LNDNC301,1,2.7etrachloroethaneND0.5ug/LNDNC301,1,2.7etrachloroethaneND0.5ug/LNDNC301,1,1.7etrichloroethaneND0.5ug/LNDNC301,1,1.7etrichloroethaneND0.5ug/LNDNC301,1,2.7etrichloroethaneND0.5ug/LNDNC301,1,2.7etrichloroethaneND0.5ug/LNDNC301,1.2.7etrichloroethaneND0.5ug/LNDNC301,1.2.7etrichloroethaneND0.5ug/LNDNC301,1.2.7etrichloroethaneND0.5ug/LNDNC301,1.2.7etrichloroethaneND0.5ug/L	•			-						
Methyl Ethyl Ketone (2-Butanone)         ND         5.0         ug/L         ND         ND         S0         ug/L         ND         NC         30           Methyl Isobutyl Ketone         ND         5.0         ug/L         ND         NC         30           Methyl Ierbutyl ether         ND         5.0         ug/L         ND         NC         30           Methyl Isobutyl Ketone         ND         5.0         ug/L         ND         NC         30           Methylene Chloride         ND         0.5         ug/L         ND         NC         30           1,1,2,2-Tetrachloroethane         ND         0.5         ug/L         ND         NC         30           1,1,1,2-Tetrachloroethane         ND         0.5         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           1,1,1-Trichloroethane         ND         0.5         ug/L         ND         NC         30           Vinyl chloride         ND         0.5         ug/L         ND         NC				-						
Methyl lsobutyl Ketone         ND         5.0         ug/L         ND         NC         30           Methyl terl-butyl ether         ND         2.0         ug/L         ND         NC         30           Methyl terl-butyl ether         ND         5.0         ug/L         ND         NC         30           Methylene Chloride         ND         5.0         ug/L         ND         NC         30           Styrene         ND         0.5         ug/L         ND         NC         30           1,1,2Tetrachloroethane         ND         0.5         ug/L         ND         NC         30           Tetrachloroethane         ND         0.5         ug/L         ND         NC         30           Toluene         ND         0.5         ug/L         ND         NC         30           1,1,2-Trichloroethane         ND         0.5         ug/L         ND         NC         30           1,1,2-Trichloroethane         ND         0.5         ug/L         ND         NC         30           1,1,2-Trichloroethane         ND         0.5         ug/L         ND         NC         30           Trichloroethane         ND         0.5 <td>Methyl Ethyl Ketone (2-Butanone)</td> <td>ND</td> <td>5.0</td> <td>-</td> <td>ND</td> <td></td> <td></td> <td>NC</td> <td>30</td> <td></td>	Methyl Ethyl Ketone (2-Butanone)	ND	5.0	-	ND			NC	30	
Methylene         ND         2.0         ug/L         ND		ND		-					30	
Methylene Chloride         ND         5.0         ug/L         ND         ND         30           Styrene         ND         0.5         ug/L         ND         NC         30           1,1,2-Tetrachloroethane         ND         0.5         ug/L         ND         NC         30           1,1,2-Tetrachloroethane         ND         0.5         ug/L         ND         NC         30           Tetrachloroethane         ND         0.5         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-Tichloroethane         ND         0.5         ug/L         ND         NC         30           1,1,2-Tichloroethane         ND         0.5         ug/L         ND         NC         30           Trichloroffuoromethane         ND         0.5         ug/L         ND         NC         30           Vinyl chloride         ND         0.5         ug/L         ND         NC         30           o-Xylene         ND         0.5         ug/L		ND		-	ND			NC	30	
1,1,2-Tetrachloroethane       ND       0.5       ug/L       ND       ND       NC       30         1,1,2,2-Tetrachloroethane       ND       0.5       ug/L       ND       NC       30         Tetrachloroethylene       ND       0.5       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         1,1,2-Trichloroethane       ND       0.5       ug/L       ND       NC       30         Trichloroethylen       ND       0.5       ug/L       ND       NC       30         Trichlorofluoromethane       ND       0.5       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         Surrogate: Toluene-d8       83.6       ug/L       ND       NC       30         Surrogate: Toluene-d8       83.6	Methylene Chloride	ND	5.0	ug/L	ND			NC	30	
1,1,2,2-Tetrachloroethane       ND       0.5       ug/L       ND       NC       30         Tetrachloroethylene       ND       0.5       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         1,1,2-Trichloroethylene       ND       0.5       ug/L       ND       NC       30         Trichloroethylene       ND       0.5       ug/L       ND       NC       30         Trichlorofluoromethane       ND       0.5       ug/L       ND       NC       30         Vinyl chloride       ND       0.5       ug/L       ND       NC       30         Surrogate: 4-Bromofluorobenzene       ND       0.5       ug/L       ND       NC       30         Surrogate: Toluene-d8       83.6       ug/L       ND       50-140       50-140       50-140       50-140       50-140       50-140       50-140       50-140       50-140       50-140       50-140       50-140	Styrene	ND	0.5	ug/L	ND			NC	30	
TetrachloroethyleneND $0.5$ $ug/L$ NDND $NC$ $30$ TolueneND $0.5$ $ug/L$ NDNC $30$ 1,1,1-TrichloroethaneND $0.5$ $ug/L$ NDNC $30$ 1,1,2-TrichloroethaneND $0.5$ $ug/L$ NDNC $30$ 1,1,2-TrichloroethyleneND $0.5$ $ug/L$ NDNC $30$ TrichloroethyleneND $0.5$ $ug/L$ NDNC $30$ TrichloroethyleneND $0.5$ $ug/L$ NDNC $30$ Vinyl chlorideND $0.5$ $ug/L$ NDNC $30$ Vinyl chlorideND $0.5$ $ug/L$ NDNC $30$ o-XyleneND $0.5$ $ug/L$ NDNC $30$ o-XyleneND $0.5$ $ug/L$ NDNC $30$ Surrogate: A-Bromofluorobenzene $77.6$ $ug/L$ $97.7$ $50-140$ $50-140$ Surrogate: Toluene-d8 $83.6$ $ug/L$ ND $50-140$ $104$ $50-140$ BenzeneND $0.5$ $ug/L$ ND $NC$ $30$ EthylbenzeneND $0.5$ $ug/L$ ND $NC$ $30$ TolueneND $0.5$ $ug/L$ <	1,1,1,2-Tetrachloroethane	ND	0.5	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           1,1,2-Trichloroethane         ND         0.5         ug/L         ND         NC         30           Trichloroethylene         ND         0.5         ug/L         ND         NC         30           Trichloroethylene         ND         0.5         ug/L         ND         NC         30           Vinyl chloride         ND         0.5         ug/L         ND         NC         30           o-Xylene         ND         0.5         ug/L         ND         NC         30           Surrogate: Dibromofluorobenzene         78.2         ug/L         ND         NC         30           Surrogate: Toluene-d8         83.6         ug/L         97.7         50-140         50-140           Benzene         ND         0.5         ug/L         ND         104         50-140           Benzene         ND         0.5         ug/L         ND	1,1,2,2-Tetrachloroethane	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       78.2       ug/L       ND       NC       30         Surrogate: Toluene-d8       83.6       ug/L       97.7       50-140       -         Surrogate: Toluene-d8       83.6       ug/L       104       50-140       -         Benzene       ND       0.5       ug/L       ND       -       -         Benzene       ND       0.5       ug/L       ND       NC       30         Toluene       ND       0.5       ug/L       ND       NC	Tetrachloroethylene	ND	0.5	ug/L	ND			NC	30	
$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	Toluene	ND	0.5	ug/L	ND			NC	30	
$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	1,1,1-Trichloroethane	ND	0.5	ug/L	ND			NC	30	
$ \begin{array}{c c c c c c c c c c c c c c c c c c c $	1,1,2-Trichloroethane	ND	0.5	ug/L	ND			NC	30	
Vinyl chloride         ND         0.5         ug/L         ND	Trichloroethylene	ND	0.5	ug/L	ND			NC	30	
m,p-Xylenes         ND         0.5         ug/L         ND	Trichlorofluoromethane	ND	1.0	ug/L	ND			NC	30	
o-Xylene       ND       0.5       ug/L       ND	Vinyl chloride	ND	0.5	ug/L	ND			NC		
Surrogate: 4-Bromofluorobenzene       78.2       ug/L       97.7 $50-140$ Surrogate: Dibromofluoromethane       77.6       ug/L       97.0 $50-140$ Surrogate: Toluene-d8       83.6       ug/L       104 $50-140$ Benzene       ND       0.5       ug/L       ND $50-140$ Ethylbenzene       ND       0.5       ug/L       ND $50-140$ Toluene       ND       0.5       ug/L       ND $50-140$ m,p-Xylenes       ND       0.5       ug/L       ND $50-140$ m,p-Xylene       ND       0.5       ug/L       ND $50-140$ m,p-Xylenes       ND       0.5       ug/L       ND $NC$ $30$ o-Xylene       ND       0.5       ug/L       ND       NC $30$	m,p-Xylenes	ND	0.5		ND			NC	30	
Surrogate: Dibromofluoromethane         77.6         ug/L         97.0         50-140           Surrogate: Toluene-d8         83.6         ug/L         104         50-140           Benzene         ND         0.5         ug/L         ND         50-140           Ethylbenzene         ND         0.5         ug/L         ND         NC         30           Toluene         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	o-Xylene	ND	0.5	ug/L	ND			NC	30	
Surrogate: Toluene-d8         83.6         ug/L         104         50-140           Benzene         ND         0.5         ug/L         ND         NC         30           Ethylbenzene         ND         0.5         ug/L         ND         NC         30           Toluene         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			ug/L						
Benzene         ND         0.5         ug/L         ND         NC         30           Ethylbenzene         ND         0.5         ug/L         ND         NC         30           Toluene         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: Dibromofluoromethane			ug/L		97.0	50-140			
Ethylbenzene         ND         0.5         ug/L         ND         NC         30           Toluene         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: Toluene-d8	83.6		ug/L		104	50-140			
Toluene         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	Benzene	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	Ethylbenzene	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		ND	0.5	ug/L	ND			NC	30	
	m,p-Xylenes	ND			ND			NC	30	
Surrogate: Toluene-d8 83.6 ug/l 104 50-140			0.5	ug/L	ND			NC	30	
	Surrogate: Toluene-d8	83.6		ug/L		104	50-140			



Surrogate: Toluene-d8

Benzene

Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 31397

# Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1740	25	ug/L	ND	86.8	68-117			
F2 PHCs (C10-C16)	1840	100	ug/L	ND	115	60-140			
F3 PHCs (C16-C34)	4590	100	ug/L	ND	117	60-140			
F4 PHCs (C34-C50)	2840	100	ug/L	ND	115	60-140			
Volatiles			9						
Acetone	96.6	5.0	ug/L	ND	96.6	50-140			
Benzene	40.3	0.5	ug/L	ND	101	60-130			
Bromodichloromethane	41.9	0.5	ug/L	ND	105	60-130			
Bromoform	42.4	0.5	ug/L	ND	106	60-130			
Bromomethane	47.4	0.5	ug/L	ND	119	50-140			
Carbon Tetrachloride	41.7	0.2	ug/L	ND	104	60-130			
Chlorobenzene	41.3	0.5	ug/L	ND	103	60-130			
Chloroform	41.6	0.5	ug/L	ND	104	60-130			
Dibromochloromethane	43.5	0.5	ug/L	ND	109	60-130			
Dichlorodifluoromethane	44.2	1.0	ug/L	ND	110	50-140			
1,2-Dichlorobenzene	41.9	0.5	ug/L	ND	105	60-130			
1,3-Dichlorobenzene	42.4	0.5	ug/L	ND	106	60-130			
1,4-Dichlorobenzene	41.9	0.5	ug/L	ND	105	60-130			
1,1-Dichloroethane	42.4	0.5	ug/L	ND	106	60-130			
1,2-Dichloroethane	40.4	0.5	ug/L	ND	101	60-130			
1,1-Dichloroethylene	40.7	0.5	ug/L	ND	102	60-130			
cis-1,2-Dichloroethylene	41.5	0.5	ug/L	ND	104	60-130			
trans-1,2-Dichloroethylene	42.4	0.5	ug/L	ND	106	60-130			
1,2-Dichloropropane	41.7	0.5	ug/L	ND	104	60-130			
cis-1,3-Dichloropropylene	39.5	0.5	ug/L	ND	98.8	60-130			
trans-1,3-Dichloropropylene	39.2	0.5	ug/L	ND	98.0	60-130			
Ethylbenzene	40.2	0.5	ug/L	ND	101	60-130			
Ethylene dibromide (dibromoethane, 1,2	39.0	0.2	ug/L	ND	97.5	60-130			
Methyl Ethyl Ketone (2-Butanone)	86.2	5.0	ug/L	ND	86.2	50-140			
Methyl Isobutyl Ketone	89.7	5.0	ug/L	ND	89.7	50-140			
Methyl tert-butyl ether	87.2	2.0	ug/L	ND	87.2	50-140			
Methylene Chloride	37.0	5.0	ug/L	ND	92.6	60-130			
Styrene	38.6	0.5	ug/L	ND	96.6	60-130			
1,1,1,2-Tetrachloroethane	40.2	0.5	ug/L	ND	101	60-130			
1,1,2,2-Tetrachloroethane	39.8	0.5	ug/L	ND	99.4	60-130			
Tetrachloroethylene	40.6	0.5	ug/L	ND	102	60-130			
Toluene	40.8	0.5	ug/L	ND	102	60-130			
1,1,1-Trichloroethane	41.7	0.5	ug/L	ND	104	60-130			
1,1,2-Trichloroethane	40.2	0.5	ug/L	ND	100	60-130			
Trichloroethylene	42.7	0.5	ug/L	ND	107	60-130			
Trichlorofluoromethane	43.3	1.0	ug/L	ND	108	60-130			
Vinyl chloride	46.8	0.5	ug/L	ND	117	50-140			
m,p-Xylenes	81.0	0.5	ug/L	ND	101	60-130			
o-Xylene	40.0	0.5	ug/L	ND	100	60-130			
Surrogate: 4-Bromofluorobenzene	83.9		ug/L		105	50-140			
Surrogate: Dibromofluoromethane	84.2		ug/L		105	50-140			
	01.0		· 3· =		100	50 1 10			

Report Date: 08-Oct-2020 Order Date: 5-Oct-2020

Project Description: PE4999

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

ug/L

ug/L

102

101

ND

50-140

60-130

81.2

40.3

0.5



Report Date: 08-Oct-2020 Order Date: 5-Oct-2020

Project Description: PE4999

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Ethylbenzene	40.2	0.5	ug/L	ND	101	60-130			
Toluene	40.8	0.5	ug/L	ND	102	60-130			
m,p-Xylenes	81.0	0.5	ug/L	ND	101	60-130			
o-Xylene	40.0	0.5	ug/L	ND	100	60-130			
Surrogate: Toluene-d8	81.2		ug/L		102	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. 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.

GPARACEL LABORATORIES LTI									3( O p:	ttawa 1-80	19 St. , Onta 0-749	Laurent ario K1G -1947 baracella	6 4J8		.Nº	Lab Use	6611	
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Nick Sullivan				Quote #										011	Day		□ 3 I	Day
154 Colonnade Pd. 5 Telephone: 613-226-7381				Email Address:	an@pa	ters	×0.	30	000	D.C	a			Date	Day Requi	red:	Re	gular
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2041094	Matrix	Air Volume	of Containers	Sample	Taken	PHCs F1-F4+BTEX	cs	Hs	Metals by ICP		B (HWS)	PHCs	5					
Sample ID/Location Name		Ϋ́	壮	Date	Time	PH	vocs	PAHs	Š	Εğ		6L	Ĩ					
1 BHZ-GWI	GW		3	Oct. 2.20	PM	Х	_		_									
2 BH4-GW1			3				X					X						
3 BH5-GWI			3				X					X						
4 DUPI	V		Ζ	V	V		Х											
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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: 31218 Project: PE4999 Custody: 54893

Report Date: 18-Nov-2020 Order Date: 12-Nov-2020

Order #: 2046443

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

Paracel ID 2046443-01

**Client ID** BH6-20-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: 18-Nov-2020 Order Date: 12-Nov-2020

Order #: 2046443

Project Description: PE4999

## **Analysis Summary Table**

Analysis	Method Reference/Description	Extraction Date	Analysis Date
PHC F1	CWS Tier 1 - P&T GC-FID	13-Nov-20	14-Nov-20
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	17-Nov-20	18-Nov-20
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	13-Nov-20	14-Nov-20



Report Date: 18-Nov-2020

Order Date: 12-Nov-2020

Project Description: PE4999

	Client ID: Sample Date: Sample ID:	BH6-20-GW1 12-Nov-20 09:00 2046443-01		- - -	
	MDL/Units	Water	-	-	-
Volatiles					
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 (dibromoethane, 1,2-)	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	-	-	-



Report Date: 18-Nov-2020

Order Date: 12-Nov-2020

Project Description: PE4999

	F				
	Client ID:	BH6-20-GW1	-	-	-
	Sample Date:	12-Nov-20 09:00	-	-	-
	Sample ID:	2046443-01	-	-	-
	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	116%	-	-	-
Dibromofluoromethane	Surrogate	102%	-	-	-
Toluene-d8	Surrogate	115%	-	-	-
Hydrocarbons					
F1 PHCs (C6-C10)	25 ug/L	<25	-	-	-
F2 PHCs (C10-C16)	100 ug/L	<100	-	-	-
F3 PHCs (C16-C34)	100 ug/L	<100	-	-	-
F4 PHCs (C34-C50)	100 ug/L	<100	-	_	-



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Certificate of Analysis Client: Paterson Group Consulting Engineers Client PO: 31218

# Method Quality Control: Blank

Report Date: 18-Nov-2020

Order Date: 12-Nov-2020

Project Description: PE4999

		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		440	50 4 40			
Surrogate: 4-Bromofluorobenzene	90.2		ug/L		113	50-140			
Surrogate: Dibromofluoromethane	78.1		ug/L		97.7	50-140			
Surrogate: Toluene-d8	95.2		ug/L		119	50-140			



# Method Quality Control: Duplicate

Order #: 2046443
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Report Date: 18-Nov-2020 Order Date: 12-Nov-2020

Project Description: PE4999

		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			0						
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	3.03	0.5	ug/L	3.07			1.3	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	6.90	0.5	ug/L	6.81			1.3	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	ND	0.5	ug/L	ND			NC	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	
Styrene	ND	0.5	ug/L	ND			NC	30	
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L	ND			NC	30	
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L	ND			NC	30	
Tetrachloroethylene	ND	0.5	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	94.7		ug/L		118	50-140			
Surrogate: Dibromofluoromethane	80.2		ug/L		100	50-140			
Surrogate: Toluene-d8	93.6		ug/L		117	50-140			



# Method Quality Control: Spike

Report Date: 18-Nov-2020

Order Date: 12-Nov-2020

Project Description: PE4999

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	2120	25	ug/L	ND	106	68-117			
F2 PHCs (C10-C16)	1800	100	ug/L	ND	112	60-140			
F3 PHCs (C16-C34)	4220	100	ug/L	ND	108	60-140			
F4 PHCs (C34-C50)	2940	100	ug/L	ND	119	60-140			
Volatiles									
Acetone	110	5.0	ug/L	ND	110	50-140			
Benzene	36.6	0.5	ug/L	ND	91.4	60-130			
Bromodichloromethane	34.1	0.5	ug/L	ND	85.2	60-130			
Bromoform	37.8	0.5	ug/L	ND	94.4	60-130			
Bromomethane	40.9	0.5	ug/L	ND	102	50-140			
Carbon Tetrachloride	33.8	0.2	ug/L	ND	84.4	60-130			
Chlorobenzene	41.8	0.5	ug/L	ND	105	60-130			
Chloroform	36.4	0.5	ug/L	ND	91.1	60-130			
Dibromochloromethane	40.2	0.5	ug/L	ND	100	60-130			
Dichlorodifluoromethane	37.2	1.0	ug/L	ND	93.1	50-140			
1,2-Dichlorobenzene	46.5	0.5	ug/L	ND	116	60-130			
1,3-Dichlorobenzene	46.0	0.5	ug/L	ND	115	60-130			
1,4-Dichlorobenzene	43.9	0.5	ug/L	ND	110	60-130			
1,1-Dichloroethane	36.3	0.5	ug/L	ND	90.8	60-130			
1,2-Dichloroethane	31.0	0.5	ug/L	ND	77.4	60-130			
1,1-Dichloroethylene	36.1	0.5	ug/L	ND	90.3	60-130			
cis-1,2-Dichloroethylene	38.3	0.5	ug/L	ND	95.8	60-130			
trans-1,2-Dichloroethylene	38.4	0.5	ug/L	ND	96.0	60-130			
1,2-Dichloropropane	37.2	0.5	ug/L	ND	93.0	60-130			
cis-1,3-Dichloropropylene	38.0	0.5	ug/L	ND	95.0	60-130			
trans-1,3-Dichloropropylene	34.8	0.5	ug/L	ND	86.9	60-130			
Ethylbenzene	40.6	0.5	ug/L	ND	102	60-130			
Ethylene dibromide (dibromoethane, 1,2	40.6	0.2	ug/L	ND	102	60-130			
Hexane	50.0	1.0	ug/L	ND	125	60-130			
Methyl Ethyl Ketone (2-Butanone)	97.5	5.0	ug/L	ND	97.5	50-140			
Methyl Isobutyl Ketone	89.3	5.0	ug/L	ND	89.3	50-140			
Methyl tert-butyl ether	92.7	2.0	ug/L	ND	92.7	50-140			
Methylene Chloride	37.3	5.0	ug/L	ND	93.4	60-130			
Styrene	37.6	0.5	ug/L	ND	94.0	60-130			
1,1,1,2-Tetrachloroethane	39.4	0.5	ug/L	ND	98.5	60-130			
1,1,2,2-Tetrachloroethane	44.9	0.5	ug/L	ND	112	60-130			
Tetrachloroethylene	44.5	0.5	ug/L	ND	111	60-130			
Toluene	42.6	0.5	ug/L	ND	106	60-130			
1,1,1-Trichloroethane	35.3	0.5	ug/L	ND	88.4	60-130			
1,1,2-Trichloroethane	37.9	0.5	ug/L	ND	94.7	60-130			
Trichloroethylene	37.4	0.5	ug/L	ND	93.6	60-130			
Trichlorofluoromethane	35.6	1.0	ug/L	ND	89.0	60-130			
Vinyl chloride	33.0	0.5	ug/L	ND	82.6	50-140			
m,p-Xylenes	85.0	0.5	ug/L	ND	106	60-130			
o-Xylene	39.8	0.5	ug/L	ND	99.6	60-130			
Surrogate: 4-Bromofluorobenzene	92.5		ug/L		116	50-140			
Surrogate: Dibromofluoromethane	73.5		ug/L		91.9	50-140			
Surrogate: Toluene-d8	83.3		ug/L		104	50-140			



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. 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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