

Phase II – Environmental Site Investigation

137-141 George Street and 110-116 York Ottawa, Ontario

Prepared for Claridge Homes

Report: PE2709-5



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EXECUTIVE SUMMARY

Assessment

A Phase II ESA was conducted for the property addressed 137-141 George Street and 110-116 York Street, in the City of Ottawa, Ontario. The purpose of the Phase IIESA was to address potentially contaminating activities (PCAs) that were identified during the Phase I ESA and considered to result in areas of potential environmental concern (APECs) on the Phase II Property.

The Phase II ESA was carried out in conjunction with a Geotechnical Investigation and consisted of three drilling programs were carried out on the following dates: February 24-28, 2023, August 14-16, 2023, and May 8-May 9. Together, the field programs consisted of drilling 16 boreholes to address the APECS identified in the Phase I ESA. Twelve of the boreholes were instrumented with groundwater monitoring wells to assess the groundwater beneath the Phase II Property. The monitoring well installed in BH1-23 was installed at a greater depth than the remainder of the wells, for geotechnical purposes.

The borehole profiles generally consist of a surficial layer of asphaltic concrete (ranging from 0.05-0.06 m in thickness), followed by fill material consisting of brown silty sand with gravel, crushed stone, cobbles and/or trace topsoil (with occasional pieces of building debris). The fill layer extended to a maximum depth of approximately 3.8 m and was underlain by glacial till consisting of a silty sand to sandy silt matrix with gravel, cobbles, and boulders. Grey limestone bedrock was encountered at a maximum depth of approximately 5.9 m below the existing ground surface.

<u>Soil</u>

A total of 40 samples and 4 duplicate samples were submitted for analysis of metals (including As, Sb, Se, Hg and CrVI), benzene, toluene, ethylbenzene xylenes (BTEX), petroleum hydrocarbons (PHCs, F1-F4), polycyclic aromatic hydrocarbons (PAHs), volatile organic compounds (VOCs) and/or pH.

Metal, Mercury (Hg), PAH and/or PHC impacts were identified in the fill material across most of the Phase II Property. PHC impacts were also identified in the native glacial till in the southern portion of the Phase II Property.

Groundwater

Groundwater samples from monitoring wells installed in BH2-23, BH3-23, BH5-23 and BH6-23 were collected during the March 8, 2023, sampling event and submitted for laboratory analysis of BTEX, PHCs (F1-F4), and/or VOCs. The majority of the analyzed

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PHC parameters were non-detect apart from the PHC F1 concentration in BH1-GW1, which complied with the applicable MECP Table 3 standards. The concentrations of PHC fractions F1 and F2 identified in BH5-23-GW1 exceeded the applicable MECP Table 3 standards, resulting in the completion of a second round of sampling.

All the analyzed VOC parameters were non-detect and therefore comply with the applicable MECP Table 3 standards, apart from the identified chloroform concentrations in BH3-23-GW1 and its duplicate sample (DUP1-23-GW1).

A second round of groundwater sampling was completed on March 23, 2023, and involved the analytical testing of PHCs and/or VOCs from groundwater obtained within BH3-23 and BH5-23. All the analyzed PHC and VOC parameters complied with the applicable MECP Table 3 standards.

A third round of groundwater sampling was completed on June 27, 2023, to obtain a second clean groundwater result from BH5-23, to comply with industry standards. The groundwater sample was submitted for BTEX and PHCs. Based on the analytical test results the concentrations of PHC fractions F1 and F2 once again exceeded the applicable MECP Table 3 standards.

Groundwater samples recovered from BH7-23 through BH9-23 were sampled on August 23, 2023, and submitted for analytical testing of BTEX or VOCs and PHCs. The monitoring well installed in BH10-23 was dry at the time of the sampling event. Apart from the chloroform concentration identified in BH9-23, all parameters complied with the MECP Table 3 standards. It should be noted that the PHC detection limits for Sample BH9-23 were elevated above the standards due to low sample volume.

Groundwater samples recovered from BH1-24, BH3-24 and BH4-24 on May 22, 2024, were submitted for analytical testing of Metals, Mercury, BTEX, VOCs and/or PHCs. Based on the analytical test results, all identified concentrations complied with the MECP Table 3 standards.

Recommendations

A remediation program was recommended in conjunction with site redevelopment, to remove all impacted soil and groundwater from the Phase II Property. Based on the findings of the Phase II ESA, it is expected that most of the fill will require of-site disposal at a registered landfill site. Deeper impacts within the native material are contained primarily to the southwestern corner of the site where groundwater impacts were also observed. It is expected that impacted groundwater, identified within the upper bedrock, near the soil-bedrock interface, will be remediation through the removal of impacted soil and underlying bedrock.

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The remediation program is underway, with most of the soil having been removed from the site for disposal at a licenced landfill site or for beneficial reuse at a Class 1 Management Site or Reuse Site, in accordance with O.Reg. 406/19.

This Phase II ESA report will be updated with the findings of the remediation program once post-remediation groundwater monitoring has been completed, in accordance with O.Reg. 153/04, to support the filing of a Record of Site Condition.

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1.0 INTRODUCTION

At the request of Claridge Homes, Paterson Group (Paterson) conducted a Phase II Environmental Site Assessment for the property addressed 141 George Street, which includes the civic addresses 137 George Street, 141 George Street, 110 York Street and 116 York Street, in the City of Ottawa, Ontario.

The purpose of this Phase II ESA has been to further address areas of potential environmental concern (APECs) identified on the Phase II Property and to delineated previously identified soil and groundwater impacts. The report also includes findings of previous field investigations.

1.1 Site Description

Address: 137-141 George Street and 110-116 York Street,

Ottawa, Ontario

Location: The Phase I Property is located between York Street

and George Street, approximately 20m east of Dalhousie Street, in the City of Ottawa, Ontario. Refer to Figure 1 - Key Plan in the Figures section following

the text.

Latitude and Longitude: 45° 25' 43.2" N, 75° 41' 24.9" W

Site Description:

Configuration: Rectangular (approximate)

Site Area: 0.41 ha (approximate)

Zoning: MD2 – Mixed Downtown Use (with a Mature

Neighbourhoods Overlay on the York Street parcels)

1.2 Property Ownership

Paterson was engaged to conduct this Phase I-ESA by Mr. Stephen Poon with Claridge Homes, the current property owner. The Claridge Homes head office is located at 210 Gladstone Avenue, Suite 2001, Ottawa, Ontario K2P and 0Z9. Mr. Poon can by reached by telephone at (613) 233-6030.

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1.3 Current and Proposed Future Uses

The northern portion of the Phase II Property addressed 110 York Street was most recently occupied by a vacant, two-storey commercial building fronting onto York Street. The building was demolished in April of 2024.

The remainder of the Phase II Property was occupied by commercial parking.

A multi-storey mixed used development is proposed. An underground parking lot associated with the proposed multi-storey residential building will occupy the entirety of the Phase II Property, while above-grade levels will be shared between the residential building and commercial hotel.

1.4 Applicable Site Condition Standard

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

J	Coarse-grained soil conditions
J	Full depth generic site conditions
J	Non-potable groundwater conditions
J	Residential land use

Section 35 of O.Reg. 153/04 applies to the Phase II Property as the Phase II Property and neighbouring properties are all serviced by the municipality.

Section 41 of O.Reg. 153/04 does not apply to the Phase II Property, as the property is not within 30m of an environmentally sensitive area and the pH of the surface soil is between 5 and 9, while the pH of the subsurface soil is between 5 and 11.

Section 43.1 of O.Reg. 153/04 does not apply to the Phase II Property in that the property is not a Shallow Soil property and the property is not within 30m of a water body.

Coarse-grained soil standards were chosen as a conservative approach. Grain size analysis was not completed. The most sensitive intended use of the Phase II Property is residential; therefore, the Residential Standards have been selected for the purpose of this Phase II ESA.

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2.0 BACKGROUND INFORMATION

2.1 Physical Setting

The Phase II Property is located between York Street and George Street, approximately 20m east of Dalhousie Street in the City of Ottawa, Ontario. According to the City of Ottawa website, the Phase II Property is situated in a mixed used downtown zone with surrounding properties consisting of commercial and residential land use.

At the time of the most recent site visits, the Phase II Property was primarily occupied by a paved asphaltic parking lot that occupies the central and southern portions of the property. The northern portion of the Phase II Property addressed 110 York Street was occupied by the foundation and building rubble associated with the recently demolished commercial building.

The Phase II Property is relatively flat and at grade with York and Geroge Streets. The regional topography slopes gently downward to the north, towards the Ottawa River. Site drainage occurs primarily through sheet flow to catch basins located in the paved asphaltic parking lot on the Phase II Property as well as along York Street and George Street.

2.2 Past Investigations

As discussed in the Phase I-ESA, the following reports were reviewed prior to conducting this assessment:

'Phase I Environmental Site Assessment, 325 Dalhousie Street, 110 York Street and 137 George Street – Ottawa, Ontario, prepared by Pinchin Environmental Ltd., dated May 2012.

Based on the findings of the Phase I ESA, a Phase II ESA was recommended to address the former RFOs on the properties addressed 321 to 325 Dalhousie Street (previously 327 Dalhousie Street, and 351 Dalhousie Street, situated at the northeast corner of the Dalhousie Street and George Street intersection, immediately west/southwest of the Phase I Property.

Phase I and Limited Phase II Environmental Site Assessment, Commercial Property, 325 Dalhousie Street, 110 York Street and 137-141 George Street, Ottawa, Ontario", prepared by Paterson Group Inc., dated June 8, 2012.

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Based on the findings of the 2013 Phase I - ESA, environmental concerns were identified with regards to the historical uses of the Phase I Property and adjacent/neighbouring properties.

Based on the findings of the 2012 Phase I - ESA, APECs were identified with regards to the historical uses of the Phase I Property and adjacent/neighbouring properties.

Historical uses of concern included former gasoline service stations, automotive service garages, printers, dry cleaners, and a roofing manufacturer.

A limited subsurface investigation was conducted at 141 George Street on May 11, 2012, and consisted of the placement of one borehole, equipped with a groundwater monitoring well, that was installed at a depth of 5.6 m below the existing ground surface. One water sample was collected from the installed monitoring well and was submitted for analytical testing of volatile organic compounds (VOC) and petroleum hydrocarbon (PHC) parameters. No detectable concentrations of these parameters were identified, and the groundwater was therefore considered to comply with the 2011 MECP standards.

Further investigative work was recommended to assess the potential for impact from remaining potential environmental concerns that were not addressed as part of the Limited Phase II ESA.

'Phase II Environmental Site Assessment, 110 York Street, 325 Dalhousie Street, 137 and 141 George Street, Ottawa, Ontario, prepared by Paterson Group Inc., dated July 27, 2012.

Four boreholes were advanced on the Phase I Property in July of 2012. The borehole locations were selected to assess the potential for impacts resulting from the identification of historical on-or off-site PCAs. Three of the boreholes were cored into bedrock and instrumented with groundwater monitoring wells.

Based on the results of the Phase II-ESA, petroleum hydrocarbon impacted soil exceeding the applicable MECP Table 3 standards was present on the southern portion of the Phase I Property, in the vicinity of BH1. The hydrocarbon impacts were considered to have resulted from the former RFO at 353 Dalhousie Street. Given that no additional soil impacts were identified at nearby borehole locations, the PHC-impacted material was considered limited in extent.

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Demolition debris including pieces of concrete and brick, as well as lead concentrations in exceedance of the applicable MECP Table 3 standards were identified in the fill material underlying the asphaltic pavement in the southern portion of the Phase I Property. The fill layer was determined to vary in thickness from approximately 1.2 to 3 m.

Based on the analytical test results, the PHC and lead impacts were considered limited in extent and did not present an immediate risk to the use of the Phase I Property.

Based on the findings of the Phase II–ESA, it was recommended that the impacted soil be remediated at the time of site redevelopment.

'Supplemental Phase II Environmental Site Assessment, 110 York Street, 321 Dalhousie Street, 137 and 141 George Street, Ottawa, Ontario", prepared by Paterson Group Inc., dated August 10, 2012.

Five additional boreholes were advanced on the Phase I Property on August 9, 2012. The borehole locations were selected to further delineate the PHC impacted soil identified in BH1 during the previous investigation. No monitoring wells were installed at the time of the supplemental investigation.

Based on the results of the Phase II-ESA, PHC-impacted soil exceeding the applicable MECP Table 3 standards was identified in the vicinity of BH1 and BH5. Like the findings presented in the previous subsurface investigation, the soil impacts were considered to have been from the former use of the property at 351/353 Dalhousie Street, as an RFO. Given that the test results met the MECP Table 3 standards at nearby borehole locations, the contaminated soil was considered limited in extent.

'Phase I Environmental Site Assessment, Proposed Mixed Use Development, 321-325 Dalhousie Street and 137-141 George Street – Ottawa, Ontario, prepared by Paterson dated April 2013.

According to historical research conducted as part of the 2013 Phase I ESA, the Phase I Property was initially developed with residential dwellings in the late 1800s. The portion of the Phase I Property addressed 137 George Street (formerly addressed 125 to 127 George Street) was occupied by a printing establishment from 1950 until 1965. The portion of the Phase I Property addressed 141 George Street (formerly addressed 137, 139 and 141 George Street and was occupied by three (3) units of an 11-unit row house in the 1950s. The neighbouring properties primarily consisted of residential dwellings, a school, office buildings and commercial establishments.

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Five PCAs that resulted in APECs on the Phase I Property were identified with respect to the historical use of the Phase I Property and surrounding lands. The identified PCA's are as follows:

Retail fuel outlet on the property addressed as 321 Dalhousie Street.
Former retail fuel outlet on the property addressed 351 Dalhousie Street
(currently addressed as 353 Dalhousie Street), immediately to the west of 137
George Street.
Printer formerly located at 125-127 George Street (currently the parking lot
addressed as 137 George Street).
Dry-cleaning business at 343 Dalhousie Street, immediately to the west of
137 George Street.
Fill material of unknown quality across the Phase I Property.

Past subsurface environmental investigations conducted for the subject property identified soil and groundwater impacted with petroleum hydrocarbons (PHCs) above the 2011 MECP Table 3 standards, on the southwestern portion of the site. A lead concentration exceeding the 2011 standard was also identified in the fill layer on the southwestern portion of the site. Based on the heterogeneous nature of the fill material and the presence of fill material across the subject property, pockets of metal-impacted fill were expected to be encountered across the site.

It was recommended that groundwater conditions be reassessed, prior to conducting a soil and groundwater remediation program.

□ 'Updated Phase II-Environmental Site Assessment, Commercial Property, 325 Dalhousie Street, 110 York Street and 137-141 George Street, Ottawa, Ontario, prepared by Paterson dated November 29, 2013.

The Phase II-ESA Updated consisted of third round of sampling at the subject property, carried out on November 12, 2013. The purpose of the sampling event was to confirm the chloroform concentrations previous identified in the groundwater at the Phase II Property, had decreased, and to reassess the PHC F2 concentrations previously identified in BH1.

Due to insufficient sample volume at BH1, the analysis of the F2 parameter could not be completed, however samples from BH1, BH2 and BH4 were submitted for analytical testing of VOCs. Chloroform was not identified in any of the samples analysed; BTEX parameters as well and/or acetone were identified at concentrations well below the Table 3 Standards.

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A remedial action plan consisting of a full-depth approach whereby all PHC and metal impacted soil would be removed from the boundaries of the subject property was recommended.

'Phase I-Environmental Site Assessment, 137-141 George Street and 110-116 York Street, Ottawa, Ontario, prepared by Paterson dated September 4, 2024

Based on the findings of the recent Phase I-ESA in combination with a review of previous reports, no new APECs were identified on the subject parcels of land addressed 137-141 George Street and 110 York Street. Fill material of unknown quality and the use of road salt for deicing purposes during winter conditions were identified as APECs on the newly included parcel of land addressed 116 York. A Phase II-ESA was recommended and carried out.

3.0 SCOPE OF INVESTIGATION

3.1 Overview of Site Investigation

The Phase II ESA, carried out in conjunction with a Geotechnical Investigation, included three separate drilling programs carried out during the interim of February 24, 2023, and May 2024.

The February 2023 investigation consisted of drilling six (6) boreholes, four (4) of which (BH2-23, BH3-23, BH5-23, and BH6-23) were instrumented with groundwater monitoring wells. These boreholes were placed to further address APECs identified during past Phase I-ESAs and provide coverage of the site for Geotechnical purposes.

The following August 2023 field program consists of the placement of four (4) boreholes, all instrumented with groundwater monitoring wells (BH7-23 through BH10-23). The purpose of the boreholes was to laterally and vertically delineate groundwater impacts identified during the February 2023 investigation.

Finally, the May 2024 field program consisted of drilling five (5) boreholes across the parcel addressed 116 York Street; this parcel had not been included in the previous investigations. Three (3) of the boreholes were instrumented with groundwater monitoring wells to access the groundwater table, if required.

The boreholes were drilled to a maximum depth of 15.3 m below the ground surface (mbgs).

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3.2 Media Investigated

During the subsurface investigation, soil samples 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 during the Phase I ESA.

The contaminants of potential concern for the soil and/or groundwater on the Phase II Property include the following:

Benzene, Toluene, Ethylbenzene, Xylenes (BTEX)
Petroleum Hydrocarbons (PHCs)
Metals (including arsenic (As), antimony (Sb), selenium (Se), mercury (Hg
and hexavalent chromium (CrVI)
Polycyclic Aromatic Hydrocarbons (PAHs)
Volatile Organic Compounds

In accordance with Section 49.1 of O.Reg.153/04, as amended, electrical conductivity (EC) and sodium adsorption ratio (SAR) are not considered to be CPCs.

3.3 Phase I Conceptual Site Model

Geological and Hydrogeological Setting

The Geological Survey of Canada website on the Urban Geology of the National Capital Area was consulted as part of this assessment. Based on this information, the bedrock in the area of the Phase I Property consists of interbedded limestone and shale of the Verulam Formation. Overburden soils are shown as glacial till, with a drift thickness on the order of 3 to 5 m. The findings of the previously completed surface investigations confirm the reported depths.

Groundwater is anticipated to flow in a northwesterly direction, towards the Ottawa River.

Existing Buildings and Structures

The northern portion of the Phase I Property is occupied by a vacant two-storey commercial building.

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Water Bodies and Areas of Natural and Scientific Interest

No water bodies are present on the Phase I Property. The closest water body is the Ottawa River, located approximately 720m west of the Phase I Property.

No areas of natural scientific interest were identified within the Phase I Study Area.

Water Well Records

A search of the MECP 's web site for all drilled well records within 250 m of the Phase I Property was conducted on February 13, 2023. Two well records were documented for the Phase I Property and pertain to monitoring wells that were installed in 2012. The monitoring wells were installed in conjunction with the previously mentioned Phase II – ESA that was completed following the identification of APECs on the Phase I Property.

In addition to the documented records, three more monitoring wells were installed on the Phase I Property in conjunction with previously completed subsurface investigations. Based on the observations made during the geotechnical and environmental assessments, the water table was intercepted at depths ranging from 3.7 to 4.5 m below the existing ground surface. Interbedded limestone and shale bedrock was encountered at a maximum depth of 5 m below the existing ground surface.

Two monitoring well records were documented for the adjacent property to the west addressed 325 Dalhousie Street. An additional three monitoring well records were documented for properties within the Phase I Study Area. The monitoring wells were drilled to depths ranging from 1.5 to 4.57 m below grade.

The identified off-site monitoring wells are not considered to be indicative of the potential for environmental impacts on the Phase I Property.

Bedrock was encountered at a maximum depth of 5.41 m below the existing ground surface, and the groundwater table was intercepted at an average depth of 4.2 m. A copy of the well records has been appended to this report.

Neighbouring Land Use

The neighbouring lands within the Phase I Study Area is primarily residential with some commercial land use. Current land use is shown on Drawing PE2709-2R – Surrounding Land Use Plan, in the Figures section of this report.

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Potentially Contaminating Activities and Areas of Potential Environmental Concern

Based on the findings of this Phase I ESA Update, four on-site and five off-site potentially contaminating activities (PCAs), were deemed to result in areas of potential environmental concern (APECs) with respect to the Phase I Property. These APECs and associated PCAs and CPCs are presented in the table below.

Table 1: Areas of Potential Environmental Concern										
Area of Potential Environmental Concern	Location of Area of Potential Environmental Concern with respect to Phase I Property	Potentially Contaminating Activity	Location of PCA (on-site or off-site)	Contaminants of Potential Concern	Media Potentially Impacted (Groundwater, Soil, and/or Sediment)					
APEC 1 Importation of Fill Material of Unknown Quality	Entire Phase I Property	"Item 30 – Importation of Fill Material of Unknown Quality"	On-site	Metals Hg CrVI PAHs BTEX PHCs (F1-F4)	Soil					
APEC 2 Former Aboveground Storage Tank (AST)	Southwestern portion of the Phase I Property	"Item 28 – Gasoline and Associated Products Storage in Fixed Tanks"	On-site	BTEX PHCs (F ₁ –F ₄)	Soil Groundwater					
APEC 3 Former Printer	Southern/sout hwestern portion of the Phase I Property	"Item 31 – Ink Manufacturing, Processing and Bulk Storage"	On-site	VOCs	Soil Groundwater					
APEC 4 Former Dry Cleaner and Machine Shop	Western portion of the Phase I Property	"Item 37 – Operation of Dry-Cleaning Equipment (where chemicals are used)"and other	Adjacent property to the west	Metals PHCs BTEX VOCs	Soil and/or Groundwater					
APEC 5 Former Retail Fuel Outlet	Southern portion of the Phase I Property	"Item 28 – Gasoline and Associated Products Storage in Fixed Tanks"	Adjacent property to the southwest	BTEX PHCs (F ₁ –F ₄)	Soil Groundwater					
APEC 6 Former Refined Petroleum Industry and Roofing Manufacturer	Southern portion of the Phase I Property	"Item 41 – Petroleum- derived Gas Refining, Manufacturing, Processing and Bulk Storage' Other	Further South of the Phase I Property, across George Street	BTEX PHCs (F ₁ –F ₄) PAHs VOCs	Soil and/or Groundwater					

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Table 1: Areas of Potential Environmental Concern										
Area of Potential Environmental Concern	Location of Area of Potential Environmental Concern with respect to Phase I Property	Potentially Contaminating Activity	Location of PCA (on-site or off-site)	Contaminants of Potential Concern	Media Potentially Impacted (Groundwater, Soil, and/or Sediment)					
APEC 7 Former Automotive Service Garage	Southern portion of the Phase I Property	"Item 52 – Storage, Maintenance, Fueling and Repair of Equipment Vehicles, and Material Used to Maintain Transportation Systems"	Further South of the Phase I Property, across George Street	BTEX PHCs (F ₁ –F ₄) PAHs VOCs	Soil Groundwater					
APEC 8 Former Printer and Dry- Cleaner	Southern portion of the Phase I Property	"Item 31 – Ink Manufacturing, Processing and Bulk Storage" "Item 37 – Operation of Dry- Cleaning Equipment (where chemicals are used)"	Further South of the Phase I Property, across George Street	Metals VOCs	Soil Groundwater					
APEC 9 Application of road salt for snow removal and de-icing purposes	Central, southern, and eastern portions of the Phase I Property	N/A	On-site	EC/SAR	Soil and Groundwater					

^{1 –} In accordance with Section 49.1 of O.Reg. 153/04 standards are deemed to be met if an applicable site condition standard is exceeded at a property solely because the qualified person has determined that a substance has been applied to surfaces for the safety of vehicular or pedestrian traffic under conditions of snow or ice or both. The exemption outlined in Section 49.1 is being relied up with respect to the RSC property.

Based on the significant coverage of asphaltic concrete on the Phase I Property, the use of salt during conditions of snow and ice is highly probable. As such, an additional APEC was included to account for the use of salt on the property. O.Reg. 153/04 is being relied upon.

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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 historical on-site and off-site PCAs that have resulted in APECs on the Phase I Property.

Additional off-site PCAs identified within the study area are not considered to represent APECs on the Phase I Property based on their separation distances and/or orientations relative to the subject land.

3.4 Deviations from Sampling and Analysis Plan

The Sampling and Analysis Plan for this project is included in Appendix 1 of this report. No deviations from the sampling and analysis plan were identified during the Phase II ESA.

3.5 Impediments

Physical impediments encountered during the Phase II ESA program include underground utilities, the former building foundation, and parked vehicles which limited the location of certain boreholes.

4.0 INVESTIGATION METHOD

4.1 Subsurface Investigation

The subsurface investigation was conducted in stages during the interim of February 24, 2023, through May 9, 2024, in conjunction with a Geotechnical Investigation.

A total of 15 boreholes were drilled to depths ranging from approximately 3.6m to 15.3m below ground surface (bgs). Eleven of the boreholes were instrumented with groundwater monitoring wells upon their completion.

The boreholes were drilled with a low clearance drill rig operated by George Downing Estate Drilling of Hawkesbury, Ontario, under full-time supervision of Paterson personnel. The borehole locations are indicated on the attached Drawing PE2709-11- Test Hole Location Plan.

Borehole and monitoring well locations from previous investigations are also presented on Drawing PE2709-11.

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4.2 Soil Sampling

A total of 83 soil samples were obtained from the boreholes by means of grab sampling from auger flights/auger samples and split spoon sampling.

Split spoon samples were taken at approximate 0.76 m intervals. Rock core samples were collected with the use of coring equipment.

The depths at which split spoon, auger flight and rock core samples were obtained from the boreholes are shown as "SS", "AU" and "RC" respectively on the Soil Profile and Test Data Sheets provided in Appendix 2. Note that borehole logs from previous investigations are also included.

The borehole profiles generally consist of a surficial layer of asphaltic concrete (ranging from 0.05-0.06 m in thickness), followed by fill material consisting of brown silty sand with gravel, crushed stone, cobbles and/or trace topsoil. Building debris fragments, including concrete, brick, metal and/or wood, were identified in the fill layer at BH2-23, BH7-23, BH8-23, BH1-24 and BH3-24.

The fill layer extended to a maximum depth of 3.58 m in BH6-23 and was underlain by glacial till consisting of a silty sand to sandy silt matrix with gravel, cobbles, and boulders. The fill material within BH7-23 was underlain by a concrete slab which had previously served as a crane base during the development of the adjacent property to the west. Grey limestone bedrock was encountered at a maximum depth of 5.64 m below the existing ground surface.

Petroleum hydrocarbon odours were noted in soil samples recovered from BH5-23, BH7-23, BH8-23 and BH10-23.

Borehole locations are shown on Drawing PE2709-11 – Test Hole Location Plan.

4.3 Field Screening Measurements

Soil samples recovered at the time of sampling 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. Allowing the samples to stabilize to room temperature ensures consistency of readings between samples.

To measure the soil vapours, the analyser probe is inserted into the nominal headspace above the soil sample. A photo ionization detector (PID) or Gastech was used to measure the vapour concentrations.

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The sample is agitated/manipulated gently as the measurement is taken. The peak reading registered within the first 15 seconds is recorded as the vapour measurement.

The maximum vapour readings measured were obtained from Samples BH5-23-SS6 and BH8-23-SS6 and were recorded as 341.5 ppm and 1,500 ppm. Otherwise, the vapour readings were generally less than 50ppm. The high vapour reading is indicative of potential significant contamination from lighter fraction petroleum hydrocarbons. The vapour screening can not be relied upon to identify heavier petroleum products. Vapour readings are noted on the Soil Profile and Test Data Sheets in Appendix 1.

4.4 Groundwater Monitoring Well Installation

Four environmental groundwater monitoring wells and one deep geotechnical monitoring well were installed on the Phase II Property as part of the subsurface investigation.

The monitoring wells consisted of 32 mm Schedule 40 threaded PVC risers and screens. Monitoring well construction details are listed in Table 2 and are also presented on the Soil Profile and Test Data Sheets provided in Appendix 1.

Borehole locations and elevations were surveyed geodetically by Paterson personnel.

TABLE 2 - Monitoring Well Construction Details											
Well ID	Ground Surface Elevation	Total Depth (m BGS)	Screened Interval (m BGS)	Sand Pack (m BGS)	Bentonite Seal (m BGS)	Casing Type					
BH1-23 (Geo)	61.52	15.32	12.3-15.3	11.5-15.3	3.6-11.5	Flushmount					
BH2-23	61.53	7.6	4.6-7.6	4.2-7.6	2.4-4.2	Flushmount					
BH3-23	61.87	7.6	4.5-7.6	3.9-7.6	2.4-4	Flushmount					
BH5-23	62.08	6.8	3.7-6.8	3.3-6.8	1.8-3.3	Flushmount					
BH6-23	62.09	7.1	4.1-7.1	4.1-7.1	2.4-3.4	Flushmount					
BH1-24	60.78	10.03	7.03-10.03	6.67-10.03	0.00-7.03	Flushmount					
BH3-24	60.92	10.26	7.26-10.26	6.67-10.26	0.00-7.26	Flushmount					
BH4-24	61.33	10.34	7.34-10.34	6.12-10.34	0.00-7.34	Flushmount					

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4.5 Field Measurement of Water Quality Parameters

The first round of groundwater sampling was conducted on March 8, 2023. Water quality parameters were measured in the field using a multi-parameter analyzer. Parameters measured in the field included temperature, pH, and electrical conductivity. Secon and third rounds of sampling were carried out in March 23, 2023 and June 27, 2023 for BH3-23 and/or BH5-23, however the results of the initial sampling event are presented below.

Field parameters were measured after each well volume purged. Wells were purged prior to sampling until at least three well volumes had been removed, the field parameters were relatively stable, or the well was dry. Stabilized field parameter values are summarized in Table 3.

Table 3: Groundwater Quality Parameters										
Well ID	Temperature (ºC)	Conductivity (µs)	рН							
March 8, 202	3									
BH2-23	9.2	3989	7.96							
BH3-23	9.7	2778	8.19							
BH5-23	11.3	2453	7.58							
BH6-23	12.3	3595	12.4							

4.6 Groundwater Sampling

Groundwater sampling protocols were followed using the MECP document entitled "Guidance on Sampling and Analytical Methods for Use at Contaminated Sites in Ontario", dated May 1996. Groundwater samples were obtained from each monitoring well, using dedicated sampling equipment.

Standing water was purged from each well prior to sampling. Samples were stored in coolers to reduce analyte volatilization during transportation.

Details of our standard operating procedure for groundwater sampling are provided in the Sampling and Analysis Plan in Appendix 1.

4.7 Analytical Testing

Based on the guidelines outlined in the Sampling and Analysis Plan appended to this report, the following soil and groundwater samples, as well as analyzed parameters are presented in Tables 4 and 5, respectively.

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TABLE	TABLE 4 – Analyzed Parameters for Submitted Soil Samples											
			Pa	rameter								
Sample ID	Sample Depth & Stratigraphic Unit	Metals¹	втех	PHCs F ₁ -F ₄	PAHs	VOCs	Rationale					
BH1-23- SS2	0.76 – 1.37 m Silty Sand (Fill Material)	Х	х	Х	х		Assess fill material of unknown quality. Sample depth determined by CPC, location of APEC and associated medium.					
BH1-23- SS3	1.60 – 2.20 m Silty Sand (Fill Material)	х	х	х	x		Assess fill material of unknown quality. Sample depth determined by CPC, location of APEC and associated medium.					
BH1-23- SS4	2.30 – 2.90 m Silty Sand (Fill Material)			Х		х	Assess potential soil impacts resulting from various off-site industries (eastern portion of Phase II Property). Sample depth determined by CPC, location of APEC and associated medium.					
BH1-23- SS5	3.2 – 3.4 m Glacial Till (Native)		x	х			Assess potential impacts resulting from various off-site industries (eastern portion of Phase II Property). Sample depth determined by CPC, location of APEC and associated medium.					
BH2-23- SS2	0.8 – 1.4 m Silty Sand (Fill Material)	Х	х	Х	х		Assess fill material of unknown quality. Sample depth determined by CPC, location of APEC and associated medium.					
BH2-23- SS4	2.4-3.0 m Silty Sand (Fill Material)	Х	Х	Х			Assess fill material of unknown quality. Sample depth determined by CPC, location of APEC and associated medium.					
BH2-23- SS5	3.10 – 3.70 m Glacial till (Native)			х		х	Assess potential soil impacts resulting from various off-site industries (northeastern portion of Phase II Property) and general coverage. Sample depth determined by CPC, location of APEC and associated medium.					
BH3-23- AU2	0.25 –0.61 m Silty Sand (Fill)	х	х	Х	х		Assess fill material of unknown quality. Sample depth determined by CPC, location of APEC and associated medium.					
BH3-23- SS4	1.6 - 2.2 m Silty Sand (Fill)	х	x	x	х	Х	Assess fill material of unknown quality and general coverage. Sample depth determined by CPC, location of APEC and associated medium.					
BH3-23- SS5	3.2 – 3.6 m Glacial Till (Native)	Х	х	Х		х	Assess potential soil impacts resulting from the former dry cleaner and machine shop previously located on the adjacent property to the west (western portion of Phase II Property) and general coverage. Sample depth determined by CPC, location of APEC and associated medium.					
DUP1-23 (duplicate of BH3- 23-SS5)	3.2 – 3.6 m Glacial Till (Native)					х	Assess potential soil impacts resulting from the former dry cleaner and machine shop previously located on the adjacent property to the west (western portion of Phase II Property) and general coverage. Sample depth determined by CPC, location of APEC and associated medium.					
BH3-23- SS6	4.6 – 5.2 m Glacial Till (Native)	х				х	Assess potential soil impacts resulting from the former dry cleaner and machine shop previously located on the adjacent property to the west (western portion of Phase II Property) and general coverage. Sample depth determined by CPC, location of APEC and associated medium.					
BH4-23- SS3	0.77 - 1.37 m Silty Sand (Fill)	Х	х	Х	х		Assess fill material of unknown quality and general coverage. Sample depth determined by CPC, location of APEC and associated medium.					
BH4-23- SS4	1.6 - 2.2 m Glacial Till (Native)	Х	x	x	х		Assess fill material of unknown quality and general coverage. Sample depth determined by CPC, location of APEC and associated medium.					

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TABLE	TABLE 4 – Analyzed Parameters for Submitted Soil Samples										
Sample	Sample Depth		Pa	ırameter							
ID	& Stratigraphic Unit	Metals¹	втех	PHCs F ₁ -F ₄	PAHs	VOCs	Rationale				
BH5-23- AU1	0.06 – 0.25 m Silty Sand (Fill)	х	х	х	х		Assess fill material of unknown quality and potential for soil impacts resulting from the former AST and printer previously located on the Phase II Property (southwestern portion of the Phase II Property). Sample depth determined by CPC, location of APEC and associated medium.				
BH5-23- SS3	1.6 – 2.2 m Silty Sand (Fill)	Х	х	х	х		Assess fill material of unknown quality and potential for soil impacts resulting from the former AST and printer previously located on the Phase II Property (southwestern portion of the Phase II Property). Sample depth determined by CPC, location of APEC and associated medium.				
BH5-23- SS6	4 – 4.6 m Glacial Till (Native)	х		х		х	Assess potential for soil impacts resulting from the former AST and printer previously located on the Phase II Property, as well as former retail fuel outlet on adjacent property to the west (southwestern portion of the Phase II Property). Sample depth determined by CPC, location of APEC and associated medium.				
DUP (duplicate of BH5- 23-SS6	4 – 4.6 m Glacial Till (Native)	х		х		х	Assess potential for soil impacts resulting from the former AST and printer previously located on the Phase II Property, as well as former retail fuel outlet on adjacent property to the west (southwestern portion of the Phase II Property). Sample depth determined by CPC, location of APEC and associated medium.				
BH6-23- SS2	0.77 – 1.37 m Silty Sand (Fill)	х	х	Х	Х		Assess fill material of unknown quality and general coverage. Sample depth determined by CPC, location of APEC and associated medium.				
BH6-23- SS5	3.2 – 3.4 m Silty Sand (Fill)	х	х	Х	Х		Assess fill material of unknown quality and general coverage Sample depth determined by CPC, location of APEC and associated medium.				
BH7-23- SS3	1.5-2.1 m Glacial Till (Native)		Х	Х			Delineation of previously identified PHC impacts. Samples selected based on results of vapour screening in				
BH8-23- SS6	3.8-4.4 m Glacial Till (Native)		Х	Х			combination with depth.				
BH1-24- AU1	0-0.3 m Silty Sand (Fill)	Х	Х	Х	Х		Assess potential impacts in the fill. Sample selected based				
BH1-24- SS3	1.5-2.1 m Silty Sand (Fill)	Х	Х	х	Х		on depth and visual observations.				
BH1-24- SS5	3.0-3.6 m Glacial Till (Native)	х	Х	Х	Х		Sample selected based on vapour screening and depth, to characterize soil for off-site disposal purposes.				
BH2-24- AU1	0.1-0.3 m Silty Sand (Fill)	Х	х	Х	Х		Assess potential impacts in the fill. Sample selected based on depth and visual observations.				
BH2-24- SS3	1.5-2.1 m Glacial Till (Native)	Х	Х	Х	Х		Sample selected based on vapour screening and depth, to characterize soil for off-site disposal purposes.				
BH3-24- AU1	0-0.46 m Silty Sand (Fill)	Х	Х	Х	Х		Assess potential impacts in the fill. Sample selected based on depth and visual observations.				
BH3-24- SS5	3.0-3.6 m Glacial Till (Native)	Х	Х	Х	Х		Sample selected based on vapour screening and depth, to characterize soil for off-site disposal purposes.				
BH4-24- AU1	0-0.46 m Silty Sand (Fill)	Х	Х	Х	х		Assess potential impacts in the fill. Sample selected based				
BH4-24- SS2	0.73-1.37 m Silty Sand (Fill)	х	Х	Х	Х		on depth and visual observations.				

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	Sample		Pa	rameter				
Sample ID	Depth & Stratigraphic Unit	Metals ¹	втех	PHCs F ₁ -	PAHs	VOCs	Rationale	
BH4-24- SS3	1.52-2.13 m Glacial Till (Native)	Х			х		Sample selected based on vapour screening and depth,	
BH4-24- SS6	3.81-4.42 m Glacial Till (Native)	х	х	Х			characterize soil for off-site disposal purposes.	
BH5-24- AU1	0.05-0.46 m Silty Sand (Fill)	х	х	х	Х		Assess potential impacts in the fill. Sample selected based on depth and visual observations.	
BH5-24- SS2 (Bottom)	0.76-1.37 m Glacial Till (Native)	х	х	х			Sample selected based on vapour screening and depth, to characterize soil for off-site disposal purposes.	

The submitted soil samples were selected for analysis based on vapour screening and field observations, in combination with information obtained during previous investigations, to assess APECs and associated CPCs identified in the Phase I – ESA.

As indicated, samples were also analysed to characterize soil for possible off-site beneficial reuse, in accordance with O.Reg. 406/19, as excess soil is expected to be generated during redevelopment. A comparison of the soil results to the Excess Soil Quality Standards, is provided under separate cover.

TABLE 5- Testing Parameters for Submitted Groundwater Samples									
		Parameters Analyzed							
Sample ID	Screened Interval	PHCs F ₁ -F ₄	ВТЕХ	VOCs	Metals ¹	PAHs	Rationale		
B2-23-GW1	4.6-7.6 m Glacial Till into Bedrock	x	x	x			Assess potential groundwater impacts resulting from various off-site industries (northeastern portion of Phase II Property). The screened interval was selected based on the depth at which the highest concentrations would likely be present (i.e., Top of water table, which straddles the soil/bedrock interface)		

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TABLE 5- Testing Parameters for Submitted Groundwater Samples									
			Parameters Analyzed						
Sample ID	Screened Interval	PHCs F ₁ -F ₄	ВТЕХ	NOCs	Metals 1	PAHs	Rationale		
BH3-23-GW1	4.5-7.6 m Glacial Till into Bedrock		X	Х			Assess potential groundwater impacts resulting from the former dry cleaner and machine shop previously located on the adjacent property to the west and for general coverage (western portion of Phase II Property) The screened interval was selected based on the depth at which the highest concentrations would likely be present (i.e., Top of water table, which straddles the soil/bedrock interface)		
DUP1-23- GW1 (duplicate of BH3-23-GW1)	4.5-7.6 Glacial Till into Bedrock		X	X			Assess potential groundwater impacts resulting from the former dry cleaner and machine shop previously located on the adjacent property to the west and for general coverage (western portion of Phase II Property). The screened interval was selected based on the depth at which the highest concentrations would likely be present (i.e., Top of water table, which straddles the soil/bedrock interface)		
BH5-23-GW1	3.7-6.8 Glacial Till into Bedrock	X	X	X			Assess potential groundwater impacts resulting from the former AST and printer on the Phase II Property, as well as the former retail fuel outlet on the adjacent property to the west (southwestern portion of the Phase II Property) The screened interval was selected based on the depth at which the highest concentrations would likely be present (i.e., Top of water table, which straddles the soil/bedrock interface)		
BH6-23-GW1	4.1-7.1 Glacial Till into Bedrock	Х	X	Х			General Coverage		
BH3-23-GW2	4.5-7.6 m Glacial Till into Bedrock	Х	X	Х			Confirm previous chloroform results have dissipated.		
BH5-23-GW2	3.7-6.8 m Glacial Till into Bedrock	Х	X				Confirm previous PHC results.		
BH5-23-GW3	3.7-6.8 m Glacial Till into Bedrock	Х	X				Commin previous Fino results.		
BH7-23-GW1	3.8-6.8 m Glacial Till into Bedrock	X	X				Lateral delineation of impacts at BH5-23.		
BH8-23-GW1	9.1-12.2 m Bedrock	Х	Χ	Χ			Vertical delineation of impacts at BH5-23.		
BH9-23-GW1	6.0-9.0 Bedrock	Χ	Χ				Lateral delineation of impacts at BH5-23.		
BH1-24-GW1	7.0-10.3 m Bedrock	Х	Х		Х	Х			
BH3-24-GW1	7.0-10.3 m Bedrock				Х	Х	General Coverage (no groundwater CPC identified)		
BH4-24-GW1	7.0-10.3 m Bedrock				Х	Х			

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

4.8 Residue Management

All soil cuttings, purge water and fluids from equipment cleaning were retained on-site.

4.9 Elevation Surveying

The ground surface elevations at each borehole location were surveyed by Paterson personnel with a high-precisions GPS unit.

4.10 Quality Assurance and Quality Control Measures

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

5.0 REVIEW AND EVALUATION

5.1 Geology

The borehole profiles generally consist of a surficial layer of asphaltic concrete (ranging from 0.05-0.06 m in thickness), followed by fill material consisting of brown silty sand with gravel, crushed stone, cobbles and/or trace topsoil. Building debris fragments, including concrete, brick, metal and/or wood, were identified in the fill layer at BH2-23, BH7-23, BH8-23, BH1-24 and BH3-24.

The fill layer extended to a maximum depth of 3.58 m in BH3-23 and was underlain by glacial till consisting of a silty sand to silty clay matrix with gravel, cobbles and boulders. The fill material within BH6-23 was underlain by a concrete slab which had previously served as a crane base during the development of the adjacent property to the west. Grey limestone bedrock was encountered at a maximum depth of 5.64 m below the existing ground surface.

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Groundwater was measured at depths ranging from approximately 2.5 to 6.3 mbgs. Based on field observations, the overburden was not considered to be the water bearing unit. Monitoring well BH10-23 installed within the overburden was dry. The groundwater is expected to be present within the bedrock near the soil-bedrock interface.

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

5.2 Groundwater Elevations, Flow Direction, and Hydraulic Gradient

Groundwater levels were measured during the groundwater sampling event on March 8, 2023, using an electronic water level meter. Groundwater levels were recorded from the monitoring wells installed in BH1-23, B2-23, BH3-23, BH5-23 and BH6-23. Groundwater levels are summarized below in Table 6.

TABLE 6 - Groundwater Level Measurements								
Borehole Location	Ground Surface Elevation (m)	Water Level Depth (m below grade)	Water Level Elevation (m ASL)	Date of Measurement				
BH1-23	61.52	4.49	57.03	March 8, 2023				
BH2-23	60.53	4.98	55.55	March 8, 2023				
BH3-23	61.87	3.48	58.39	March 8, 2023				
BH5-23	62.09	2.51	59.58	March 8, 2023				
BH6-23	62.08	2.77	59.31	March 8, 2023				
BH3-23	61.87	4.52	57.35	March 23, 2023				
BH5-23	62.09	3.99	58.1	March 23, 2023				
BH6-23	62.08	4.91	57.17	March 23, 2023				
BH1-24	60.78	3.46	57.32	May 22, 2024				
BH3-24	60.92	6.15	54.78	May 22, 2024				
BH4-24	61.33	6.29	55.04	May 22, 2024				

Based on the groundwater elevations measured during the March 2023 sampling event, groundwater contour mapping was completed. Groundwater contours are shown on Drawing PE2709-11 – Test Hole Location Plan.

Based on the contour mapping, groundwater flow at the Phase II Property is in an easterly direction. The local groundwater flow may have been influenced by the deep excavation on the adjacent property to the east. Regional groundwater flow is expected to be in a northerly direction. It should be noted that groundwater levels are expected to fluctuate throughout the year with seasonal variations.

A horizontal hydraulic gradient of approximately 0.07 m/m was calculated.

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5.3 Fine-Coarse Soil Texture

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

5.4 Soil: Field Screening

Field screening of the soil samples collected during drilling resulted in vapour readings ranging generally from 0.1 less than 50 ppm. Hydrocarbon odours were observed in soil Samples BH5-23-SS6 and BH8-23-SS6 which had the highest vapour readings of approximately 345 ppm and 1,500 ppm.

Building debris fragments, including concrete, brick, metal and/or wood, were identified in the fill layer at BH2-23, BH7-23, BH8-23, BH1-24 and BH3-24. No other unusual/deleterious materials or substances were identified within the fill material across the Phase II Property during the subsurface investigations.

The 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

Based on the findings of the field screening in combination with sample depth and location, 33 soil samples and 4 duplicate samples were submitted for analysis of metals (including As, Sb, Se, Hg and/or CrVI), PAHs, BTEX, PHCs (F1-F4), VOCs and/or pH. The results of all the analytical testing completed on the Phase II Property (including past assessments) are presented in Table A1 appended to this report. The laboratory Certificates of Analysis are also provided in the Appendix.

Metals (including As, Sb, Se), Hg and CrVI

All of the analyzed metal parameters are in compliance with the applicable MECP Table 3 Standards, with the exception of lead, mercury, and zinc in soil Sample BH2-23-SS2 and lead and mercury in soil Samples BH3-23-AU2 and BH4-24-SS2. Additionally, soil Sample BH4-SS3, which was submitted during the 2012 assessment, and Sample BH5-24-AU1, exceeded the MECP Table 3 standard for only lead.

Sample BH4-24-SS3 was submitted for analysis of methyl mercury due to the exceedance of mercury in Sample BH4-24-SS2. Methyl mercury was not detected in the sample analysed.

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The analytical results for metals (including As, Sb, Se, CrVI and Hg) tested in soil are shown on Drawing PE2709-12— Analytical Testing Plan — Soil (Metals).

<u>PAHs</u>

Various PAH parameters identified in soil Samples BH2-23-SS2, BH3-23-AU2, BH1-24-SS3, BH2-24-AU1, BH3-24-AU1, BH4-24-AU1, BH4-24-SS2 and BH5-24-AU1, exceed the MECP Table 3 Standards. All other samples analysed meet the MECP Table 3 Residential Standards.

The analytical results for PAHs tested in soil are shown on PE2709-13 – Analytical Testing Plan – Soil (PAHs).

PHCs (F₁-F₄)

All the detected PHC concentrations in the analysed soil samples comply with the applicable MECP Table 3 Standards, apart from PHC fraction F_3 concentrations in soil Samples BH1-23-SS4, BH2-23-SS2, BH5-23-AU1, BH4-24-AU1, and PHC fractions F_1 and F_2 in Sample BH8-23-SS6. The analytical results for PHCs tested in soil are shown on Drawing PE2709-14 – Analytical Testing Plan – Soil (PHCs).

Additionally, soil Sample BH1-SS7 that was submitted as part of the 2012 assessment, had PHC fraction F_1 and F_2 that exceeded the applicable MECP Table 3 standards.

BTEX

No BTEX concentrations were identified in any of the samples analysed except for Sample BH8-23-SS6 which identified concentrations of BTEX well below the MECP Table 3 standards. All samples analysed comply with the MECP Table 3 standards. The analytical results for BTEX tested in soil are shown on Drawing PE2709-15 – Analytical Testing Plan – Soil (BTEX).

<u>VOCs</u>

No VOC parameters were identified in the samples analysed. As such, the results comply with the MECP Table 3 standards. The analytical results for VOCs tested in soil are shown on Drawing PE2709-16 – Analytical Testing Plan – Soil (VOCs).

The maximum parameter concentrations identified within the soil samples are listed below in Table 7.

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TABLE 7: Maximum Concentrations – Soil							
Parameter	Maximum Concentration (μg/g)	Soil Sample	Depth Interval (m BGS)				
Antimony	2.2	BH2-23-SS2	0.8-1.4				
Arsenic	10.2	BH6-23-SS5	3.2-3.4				
Barium	321	BH2-23-SS2	0.8-1.4				
Beryllium	0.6	BH2-23-SS2	0.8-1.4				
Boron	16.2	BH6-23-SS2	0.76-1.37				
Cadmium	0.6	BH2-23-SS2	0.8-1.4				
Chromium	29.2	BH2-23-SS2	0.8-1.4				
Cobalt	11.8	BH6-23-SS5	3.2-3.4				
Copper	120	BH2-23-SS2	0.8-1.4				
Lead	524	BH4-SS3	0.8-1.2				
Mercury	2.7	BH2-23-SS2	0.8-1.4				
Molybdenum	5.4	BH6-23-SS5	3.2-3.4				
Nickel	20.3	BH6-23-SS5	3.2-3.4				
Selenium	1.3	BH2-23-SS2	0.8-1.4				
Silver	1.8	BH1-AU1	0-0.25				
Vanadium	321	BH1-AU1	0-0.25				
Zinc	429	BH2-23-SS2	0.8-1.4				
Acenaphthene	0.76	BH2-23-SS2	0.8-1.4				
Acenaphthylene	0.19	BH2-23-SS2	0.8-1.4				
Anthracene	1.58	BH2-23-SS2	0.8-1.4				
Benzo[a]anthracene	2.72	BH2-23-SS2	0.8-1.4				
Benzo[a]pyrene	2.37	BH2-23-SS2	0.8-1.4				
Benzo[b]fluoranthene	<u>3</u>	BH2-23-SS2	0.8-1.4				
Benzo[g,h,i]perylene	1.45	BH2-23-SS2	0.8-1.4				
Benzo[k]fluoranthene	1.68	BH2-23-SS2	0.8-1.4				
Chrysene	2.68	BH2-23-SS2	0.8-1.4				
Dibenzo[a,h]anthracene	0.38	BH2-23-SS2	0.8-1.4				
Fluoranthene	6.09	BH2-23-SS2	0.8-1.4				
Fluorene	0.53	BH2-23-SS2	0.8-1.4				
Indeno [1,2,3-cd] pyrene	1.4	BH2-23-SS2	0.8-1.4				
1-Methylnaphthalene	0.13	BH2-23-SS2	0.8-1.4				
2-Methylnaphthalene	0.18	BH2-23-SS2	0.8-1.4				
Methylnaphthalene (1&2)	0.31	BH2-23-SS2	0.8-1.4				
Naphthalene	0.22	BH2-23-SS2	0.8-1.4				
Phenanthrene	4.82	BH2-23-SS2	0.8-1.4				
Pyrene	4.84	BH2-23-SS2	0.8-1.4				
F1 PHCs (C6-C10)	<u>182</u>	BH1-SS7	4.6-4.7				
F2 PHCs (C10-C16)	118	BH1-SS7	4.6-4.7				
F3 PHCs (C16-C34)	699	BH5-23-AU1	0.06-0.25				
F4 PHCs (C34-C50)	1650	BH5-23-AU1	0.06-0.25				
Ethylbenzene	0.1	BH5-23-SS6	4-4.6				
Xylenes, total	0.22	BH5-23-SS6	4-4.6				

Notes:

- * Duplicate of BH3-AU1
- Bold and Underlined Results exceed the selected MECP standards

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5.6 Groundwater Quality

Five groundwater samples (including one duplicate) from monitoring wells installed in BH2-23, BH3-23, BH5-23 and BH6-23 were submitted for laboratory analysis of BTEX, PHCs and VOCs. The groundwater samples were obtained from the screened intervals noted in Table 2.

Six groundwater samples collected during the 2012 investigation, from previously existing monitoring wells in BH1, BH4 and BH1-11, were submitted for analytical testing of BTEX, PHCs and VOCs.

Analytical testing results from both the current program and 2012 investigation are presented in Table A2. The laboratory Certificates of Analysis are provided in Appendix 1.

PHCs (F₁-F₄)

Hydrocarbon parameters were not identified in any of the samples except for samples recovered from BH1 (2012), BH5-23, BH7-23 and BH1-24. The concentrations of PHC F₁ or PHC F₄ identified in BH1 (2012), BH7-23 and BH1-24, complied with the MECP Table 3 standards. Concentrations of PHC F₁ and F₂ identified in groundwater samples BH5-23-GW1 exceeded the MECP Table 3 standards, resulting in the completion of a second round of sampling.

Groundwater from BH5-23 was resampled and submitted for analytical testing of BTEX and PHCs. All the analytical test results from the second groundwater sampling event complied with the applicable MECP Table 3 standards.

A third groundwater sampling event with clean results was required to meet industry standards, however the results of this sampling event identified concentrations of PHC F_1 and F_2 similar to that identified in BH5-23-GW1, at levels above the MECP Table 3 standards.

The analytical results for groundwater tested are shown on Drawing PE2709-19–Analytical Testing Plan – Groundwater (PHCs).

BTEX/VOCs

No VOC parameters were identified in any of the samples analysed, apart from chloroform identified in BH1-GW1, BH4-GW1, BH4-GW2, BH3-23-GW1, DUP1-23-GW1, and BH8-23-GW1. Except for BH8-23-GW1, all wells were resampled, and chloroform was not identified.

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The chloroform in the groundwater samples resulted from the use of municipally treated water for the bedrock coring process required for installation of monitoring wells. All concentrations of chloroform identified were less than 240 μ g/L as lited in Table A of the MECP document entitled "Guidance for Addressing Chloroform at a Record of Site Condition Property" and as such, the exemption in Paragraph 2 of Section 49.1 or O.Reg. 153/04 is being relied upon.

All BTEX parameters identified were in compliance with the MECP Table 3 standards.

The analytical results for BTEX and VOCs in groundwater are shown on Drawing PE2709-17– Analytical Testing Plan – Groundwater (BTEX, VOCs).

5.7 Quality Assurance and Quality Control Results

All samples submitted as part of the March 2023 sampling events were handled in accordance with the Analytical Protocol with respect to preservation method, storage requirement, and container type. As per Subsection 47(3) of O.Reg. 153/04, as amended, under the Environmental Protection Act, a Certificate of Analysis has been received for each sample submitted for analysis and all Certificates of Analysis are appended to this report.

Four duplicate soil samples were collected from BH3-23-SS5 (DUP1-23), BH5-23-SS6 (DUP), BH1-24-SS5 (DUP1) and BH4-24-SS6 (DUP2), and were submitted for metals, PHCs, and/or VOCs. The duplicates were collected with the intent of calculating the relative percent difference (RPD) between duplicate sample values, as a way of assessing the quality of the analytical test results.

The RPD calculations for BH5-23-SS6, BH1-24-SS5 and BH4-24-SS6 and their respective duplicate samples are provided in Tables 8A-8C.

All of the analyzed VOC concentrations in BH3-23-SS5 and its duplicate sample, DUP1-23, were not detected above the laboratory method detection limit; as such these results are not tabulated below.

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Table 8A - QA/QC - Soil - BH5-23-SS6 and DUP								
Parameter	MDL (μg/g)	BH5-23-SS6	DUP	RPD (%)	QA/QC Result			
Arsenic	1.0	4.1	4.6	11.5	Meets Target			
Barium	1.0	129	211	48	Does Not Meet Target			
Boron	5.0	8	8.9	5.3	Meets Target			
Chromium	5.0	13.5	15.5	13.7	Meets Target			
Cobalt	1.0	4.5	4.9	4.2	Meets Target			
Copper	5.0	5	5.5	9.5	Meets Target			
Lead	1.0	5	5.1	2	Meets Target			
Nickel	5.0	11.4	12.5	9.2	Meets Target			
Vanadium	10.0	15.4	18	4.2	Meets Target			
Zinc	20.0	32.4	36.8	12.1	Meets Target			
PHC F ₁	7	7	8	13	Meets Target			
PHC F ₂	4	4	28	150	Does Not Meet Target			
Ethylbenzene	0.05	0.1	0.08	22	Does Not Meet Target			
Xylenes, total	0.05	0.22	0.21	4.6	Meets Target			

Notes:

- MDL Method Detection Limit
- nd not detected above the MDL

The remaining parameter concentrations were not detected in either or both the original sample and duplicate, therefore, the RPD values cannot be calculated.

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Table 8B - QA/QC - Soil - BH1-24-SS5 and DUP1							
Parameter	MDL BH1-24- (μg/g) SS5		DUP1	RPD (%)	QA/QC Result		
Arsenic	1.0	2.1	2.2	4.7	Meets Target		
Barium	1.0	35.8	35.9	0.28	Meets Target		
Boron	5.0	8	5.2	46.7	Does not Meet Target		
Chromium	5.0	11.8	12.2	3.3	Meets Target		
Cobalt	1.0	4.6	4.8	4.3	Meets Target		
Copper	5.0	6.8	7.8	13.7	Meets Target		
Lead	1.0	4	4.2	4.9	Meets Target		
Nickel	5.0	8.3	8.7	4.7	Meets Target		
Vanadium	10.0	18.7	19.1	2.1	Meets Target		
Acenaphthene	0.02	0.02	0.02	0	Meets Target		
Anthracene	0.02	0.06	0.06	0	Meets Target		
Benzo[a]anthracene	0.02	0.05	0.05	0	Meets Target		
Benzo[b]pyrene	0.02	0.04	0.04	0	Meets Target		
Benzo[b]fluoranthene	0.02	0.04	0.03	28.6	Does not Meet Target		
Benzo[g,h,i]perylene	0.02	0.02	0.03	40	Does not Meet Target		
Benzo[k]fluoranthene	0.02	0.03	0.03	0	Meets Target		
Chrysene	0.02	0.05	0.05	0	Meets Target		
Fluoranthene	0.02	0.2	0.18	10.5	Meets Target		
Fluorene	0.02	0.03	0.02	40	Does not Meet Target		
Naphthalene	0.01	0.04	0.02	67	Does not Meet Target		
Phenanthrene	0.02	0.19	0.18	5.4	Meets Target		
Pyrene	0.02	0.14	0.14	0	Meets Target		

Notes:

- MDL Method Detection Limit
- nd not detected above the MDL

Table 8C - QA/QC - Soil - BH4-24-SS6 and DUP2								
Parameter	MDL (µg/g)	BH1-24- SS5	DUP2	RPD (%)	QA/QC Result			
Arsenic	1.0	3	2.7	10.5	Meets Target			
Barium	1.0	36.7	43	15.8	Meets Target			
Boron	5.0	7.4	7.7	4.0	Meets Target			
Chromium	5.0	9.5	10.5	10	Meets Target			
Cobalt	1.0	2.9	3.3	12.9	Meets Target			
Lead	1.0	5.6	5.6	0	Meets Target			
Nickel	5.0	6.5	7.1	8.8	Meets Target			
Vanadium	10.0	13.6	14.9	9.1	Meets Target			

Notes:

- MDL Method Detection Limit
- nd not detected above the MDL

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Typically, RPD values below 20% indicate satisfactory quality. The relative percent difference (RPD) results calculated for three soil parameters (barium, ethylbenzene and PHC F2) identified in Sample BH5-23-SS6 and its duplicate, as well as for boron and several PAH parameters in Sample BH1-24-SS4 and its duplicate, fell outside of the acceptable range of 20%, and thus do not meet the data quality objectives outlined in the Sampling and Analysis Plan, appended to this report.

Despite the exceeded RPD values calculated for the soil samples and the corresponding duplicate, it should be noted that the concentrations of the parameters were well within the applicable MECP Table 3 Standards in both samples.

As a result, it is our opinion that the decision-making usefulness of the samples is not considered to be impaired, and thus the quality of the collected field data is sufficient to meet the overall objectives of this assessment.

A duplicate groundwater sample (DUP1-23-GW1) was obtained from the monitoring well installed in BH3-23 and submitted for laboratory analysis of VOC parameters. The RPD calculations for the original groundwater and duplicate sample are provided in Table 13.

Table 9 - QA/QC Calculations – Groundwater								
Parameter	MDL (µg/L)	BH3-23- GW1	DUP1-23- GW1	RPD (%)	QA/QC Result			
Benzene	0.5	0.6	0.6	0	Meets Target			
Chloroform	0.5	2.5	2.6	4	Meets Target			
Toluene	0.5	3.4	3.5	2.8	Meets Target			
Xylenes, Total	0.5	4.5	4.2	7	Meets Target			

Notes:

- MDL Method Detection Limit
- nd not detected above the MDL

All of the calculated RPD values from BH3-23-GW1 and the duplicate samples BH1-23-GW1 are well below the required 20%.

The quality of the field data collected during the Phase II ESA is considered to be sufficient to meet the overall objectives of the assessment.

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5.8 Phase II Conceptual Site Model

The following section has been prepared in accordance with the requirements of O.Reg. 153/04, as 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

Based on the results of the Phase I ESA completed for the subject property, nine (9) PCAs were considered to result in APECs on the Phase II Property. The identified APECs on the Phase II Property are presented in Table 1 in the Phase I CSM and are as follows:

APEC 1: Resulting from the importation of fill material of unknown quality (PCA #30);
APEC 2: Resulting from the presence of a former on-site aboveground storage tank (AST) (PCA #28);
APEC 3: Resulting from the presence of a former printer previously located on the southwestern portion of the Phase II Property (PCA #31);
APEC 4: Resulting from the presence of a former dry cleaner and machine shop on the adjacent property to the west (PCA #37).
APEC 5: Resulting from the presence of a former retail fuel outlet on the adjacent property to the west (PCA #28).
APEC 6: Resulting from the presence of a former refined petroleum industry and roofing manufacturer previously located to the south, across George Street (PCAs #41 and #28);
APEC 7: Resulting from the presence of a former automotive service garage previously located to the south, across George Street (PCA#52);
APEC 8: Resulting from the presence of a former printer and drycleaner previously located to the south, across George Street (PCA #31 and #37)

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APEC 9: Application of road salt for the removal of snow and de-icing purposes (PCA #NA)

Based on the findings of the Phase I ESA, it is considered likely that road salt was applied to the surface of the parking lot across the Phase II Property for the safety of vehicular and pedestrian traffic under conditions of ice and/or snow. Although not defined as a specific PCA under Column A of Table 2 of O.Reg. 153/04, the use of salt for safety purposes is considered to result in an APEC on the Phase I Property (APEC 4).

According to Section 49.1 of O.Reg. 153/04, if an applicable site condition standard is exceeded at a property solely because of the following reason, the applicable site condition standard is deemed not to be exceeded for the purpose of Part XV.1 of the Act: "The qualified person has determined, based on a phase one environmental site assessment or a phase two environmental site assessment, that a substance has been applied to surfaces for the safety of vehicular or pedestrian traffic under conditions of snow or ice or both."

In accordance with Section 49.1 of O.Reg. 153/04, any EC and SAR concentrations on the RSC Property that exceed the MECP Table 3 Standards for a residential/institutional land use are deemed not to be exceeded for the purpose of Part XV.1 of the Act. This exemption is being relied on for APEC 9.

Contaminants of Potential Concern

The following CPCs are identified with respect to the Phase II Property:

	Metals (including arsenic (As), antimony (Sb), selenium (Se), mercury (Hg) and hexavalent chromium (CrVI)) (Soil);
	Benzene, Toluene, Ethylbenzene, Xylenes (BTEX) (Soil and/or Groundwater);
	Petroleum Hydrocarbons (PHCs) (Soil and/or Groundwater);
	Polycyclic Aromatic Hydrocarbons (PAHs) (Soil); and
	Volatile Organic Compounds (VOCs) (Soil and/or Groundwater)
1	accordance with Ocation 40.1 of O.D. of 150/04 are accorded alcoholical

In accordance with Section 49.1 of O.Reg.153/04, as amended, electrical conductivity (EC) and sodium adsorption ratio (SAR) are not considered to be CPCs.

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Subsurface Structures and Utilities

The Phase I Property is situated in a municipally serviced area. Underground utility services on the subject land during the Phase I site visits include natural gas, electrical, cable, sewer and water services. Services have since been decommissioned.

No potable wells or private sewage systems were observed on the Phase I Property at the time of the site visit. No other subsurface structures were identified at the time of the site visit apart from the foundation associated with the former structure at 110 York Street.

Physical Setting

Site Stratigraphy

The stratigraphy of the Phase II Property generally consists of:

Groundwater was encountered within the overburden at depths ranging from 2.51 to 6.29 mbgs.

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

Asphaltic concrete; encountered at depths ranging from approximately 0.05 to 0.06 mbgs.
Fill material consisting of brown or grey silty sand with gravel, crushed stone, cobbles and/or trace topsoil; extending to a maximum depth of approximately 3.6 mbgs in BH6-23. Building debris fragments, including concrete, brick, metal and/or wood, were identified in the fill layer at BH2-23, BH7-23, BH8-23, BH1-24 and BH3-24.
Glacial Till; consisting of brown to grey silty sand to sandy silt matrix with gravel, cobbles and boulders extending to a maximum depth of 5.64 mbgs. (The fill material within BH7-23 was underlain by a concrete slab which had previously served as a crane base during the development of the adjacent property to the west.)
Grey limestone bedrock was encountered at a maximum depth of 5.64 mbgs.

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.

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Hydrogeological Characteristics

Groundwater was encountered at the soil-bedrock interface at depths ranging from 2.77 to 6.29 mbgs. Based on field observations, the soil is not considered to be a water-bearing unit; it is anticipated that the water table exists within the bedrock.

Based on the March 2023 groundwater monitoring event, groundwater flow was measured in an eastern direction with a hydraulic gradient of 0.07 m/m. It should be noted that groundwater levels are expected to fluctuate throughout the year with seasonal variations. The flow direction in the immediate vicinity of the Phase II Property is expected to have been influenced by the deep excavation on the adjacent property to the east. Regional groundwater flow is expected to be in a northerly direction, toward the Ottawa River. Groundwater contours are shown on Drawing PE2709-11 – Test Hole Location Plan.

Approximate Depth to Bedrock

Grey limestone bedrock was confirmed at each borehole location; depth to bedrock at the Phase II Property ranges from approximately 4.6 to 5.6m below grade.

Approximate Depth to Water Table

The water table was measured at depths ranging from 2.77 to 6.29 mbgs.

Sections 41 and 43.1 of the Regulation

Section 41 of the Regulation does not apply to the Phase II Property, in that the subject property is not within 30m of an environmentally sensitive area and the pH of the surface soil is between 5 and 9 while the pH of the subsurface soil is between 5 and 11.

Section 43.1 of the Regulation does not apply to the Phase II Property given the property is not a shallow soil property and is not within 30m of a body of water.

Areas Where Excess Soil Has Been Finally Placed on the Phase II Property

No excess soil has been finally placed on, in or under the Phase II Property.

Existing Buildings and Structures

The northern portion of the Phase I Property was most recently occupied by a vacant, two-storey commercial building fronting onto York Street (civic address 110 York Street). The former structure was demolished in April 2024.

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Proposed Buildings and Other Structures

The proposed site development for the Phase II Property will consist of a multistorey, mixed-used development comprised of residential dwellings and a commercial hotel. An underground parking lot will occupy the entire Phase II Property below grade.

Environmental Condition

Areas Where Contaminants are Present

Based on the findings of the current Phase II ESA, including the results of past investigations, fill material impacted with metals, mercury, PHCs and PAHs was identified in pockets across the site. Additionally, PHC impacted glacial till was also identified on the southern portion of the Phase II Property, primarily within the southwest corner. A PHC F₃ concentration identified at BH1-23 on the southeastern portion of the site within the native material, is considered to be from a separate source.

Groundwater impacted with PHCs was identified at BH5-23; based on the findings of the Phase II ESA, the impacted groundwater is confined within the southwestern portion of the Phase II Property. The groundwater impacts are expected to be present within the soil-bedrock interface.

Types of Contaminants

Contaminants include metals, mercury, PHC and PAH concentrations in the soil, as well as PHC F₁ and F₂ concentrations in the groundwater. Note that methyl mercury was also analysed in a soil sample recovered from BH4-24, given the mercury exceedance; no methyl mercury was identified in the sample analysed.

Contaminated Media

Based on the findings of this Phase II ESA, the fill material across most of the site and the native glacial till on the southern portion of Phase II Property, has been impacted with PHCs, metals, mercury and/or PAHs.

Groundwater within the southwestern portion of the site has been impacted with PHC F₁ and PHC F₂ concentrations.

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What Is Known About Areas Where Contaminants Are Present

The impacted fill material is considered to have resulted from the importation of fill material of unknown quality, for historical grading purposes following the demolition of previously existing buildings on the Phase II Property.

The native glacial till within the southwestern corner of the site is considered to have resulted from the former operation of a retail fuel outlet (RFO) on the adjacent property to the west.

Similarly, the groundwater on this portion of the site is impacted with residual PHC concentrations considered to have originated from the former RFO.

A PHC F₃ concentration identified at BH1-23 on the southeastern portion of the site within the native material, is expected to be from a separate source, possibly associated with former buildings on this portion of the site.

Distribution and Migration of Contaminants

Based on the findings of the Phase I and Phase II ESA, contaminants of concern identified within the fill material are not considered to have migrated significantly beyond the fil layer.

The native glacial till impacts are also considered to be limited in extent and are isolated primarily to the southern portion of the Phase II Property.

Discharge of Contaminants

Based on the findings of this Phase II ESA, the shallow soil impacts are considered to have resulted from the importation of fill material of unknown quality. The PHC impacted native glacial till on the southwestern portion of the Phase II Property is considered to have resulted from the former operation of an RFO on the adjacent property to the west, while the PHC F₃ concentration identified on the southeastern portion of the site within the native soil, may be associated with former on-site buildings on this portion of the Phase II Property.

Climatic and Meteorological Conditions

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

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Given the site was covered with asphaltic concrete outside of the former building footprints, significant downward leaching of contaminants is expected to have occurred, although some leaching by means of infiltration of precipitation may have occurred through cracks in the asphaltic concrete.

Groundwater levels and/or flow are not considered to have had a significant affect on contaminant distribution at the RSC Property; based on the Phase II ESA, groundwater impacts are confined to the southwestern portion of the property within the upper layers of bedrock near the soil-bedrock interface.

Potential for Vapour Intrusion

Based on the findings of the Phase II ESA, the potential for vapour intrusion is negligible given that soil impacted with volatile compounds (PHCs) was situated outside of the building footprint.

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6.0 CONCLUSIONS

Assessment

A Phase II ESA was conducted for the property addressed 137-141 George Street and 110-116 York Street, in the City of Ottawa, Ontario. The purpose of the Phase IIESA was to address potentially contaminating activities (PCAs) that were identified during the Phase I ESA and considered to result in areas of potential environmental concern (APECs) on the Phase II Property.

The Phase II ESA was carried out in conjunction with a Geotechnical Investigation and consisted of three drilling programs were carried out on the following dates: February 24-28, 2023, August 14-16, 2023, and May 8-May 9. Together, the field programs consisted of drilling 16 boreholes to address the APECS identified in the Phase I ESA. Twelve of the boreholes were instrumented with groundwater monitoring wells to assess the groundwater beneath the Phase II Property. The monitoring well installed in BH1-23 was installed at a greater depth than the remainder of the wells, for geotechnical purposes.

The borehole profiles generally consist of a surficial layer of asphaltic concrete (ranging from 0.05-0.06 m in thickness), followed by fill material consisting of brown silty sand with gravel, crushed stone, cobbles and/or trace topsoil (with occasional pieces of building debris). The fill layer extended to a maximum depth of approximately 3.8 m and was underlain by glacial till consisting of a silty sand to sandy silt matrix with gravel, cobbles, and boulders. Grey limestone bedrock was encountered at a maximum depth of approximately 5.9 m below the existing ground surface.

Soil

A total of 40 samples and 4 duplicate samples were submitted for analysis of metals (including As, Sb, Se, Hg and CrVI), benzene, toluene, ethylbenzene xylenes (BTEX), petroleum hydrocarbons (PHCs, F1-F4), polycyclic aromatic hydrocarbons (PAHs), volatile organic compounds (VOCs) and/or pH.

Metal, Mercury (Hg), PAH and/or PHC impacts were identified in the fill material across most of the Phase II Property. PHC impacts were also identified in the native glacial till in the southern portion of the Phase II Property.

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Groundwater

Groundwater samples from monitoring wells installed in BH2-23, BH3-23, BH5-23 and BH6-23 were collected during the March 8, 2023, sampling event and submitted for laboratory analysis of BTEX, PHCs (F1-F4), and/or VOCs. The majority of the analyzed PHC parameters were non-detect apart from the PHC F1 concentration in BH1-GW1, which complied with the applicable MECP Table 3 standards. The concentrations of PHC fractions F1 and F2 identified in BH5-23-GW1 exceeded the applicable MECP Table 3 standards, resulting in the completion of a second round of sampling.

All the analyzed VOC parameters were non-detect and therefore comply with the applicable MECP Table 3 standards, apart from the identified chloroform concentrations in BH3-23-GW1 and its duplicate sample (DUP1-23-GW1).

A second round of groundwater sampling was completed on March 23, 2023, and involved the analytical testing of PHCs and/or VOCs from groundwater obtained within BH3-23 and BH5-23. All the analyzed PHC and VOC parameters complied with the applicable MECP Table 3 standards.

A third round of groundwater sampling was completed on June 27, 2023, to obtain a second clean groundwater result from BH5-23, to comply with industry standards. The groundwater sample was submitted for BTEX and PHCs. Based on the analytical test results the concentrations of PHC fractions F1 and F2 once again exceeded the applicable MECP Table 3 standards.

Groundwater samples recovered from BH7-23 through BH9-23 were sampled on August 23, 2023, and submitted for analytical testing of BTEX or VOCs and PHCs. The monitoring well installed in BH10-23 was dry at the time of the sampling event. Apart from the chloroform concentration identified in BH9-23, all parameters complied with the MECP Table 3 standards. It should be noted that the PHC detection limits for Sample BH9-23 were elevated above the standards due to low sample volume.

Groundwater samples recovered from BH1-24, BH3-24 and BH4-24 on May 22, 2024, were submitted for analytical testing of Metals, Mercury, BTEX, VOCs and/or PHCs. Based on the analytical test results, all identified concentrations complied with the MECP Table 3 standards.

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Recommendations

A remediation program was recommended in conjunction with site redevelopment, to remove all impacted soil and groundwater from the Phase II Property. Based on the findings of the Phase II ESA, it is expected that most of the fill will require of-site disposal at a registered landfill site. Deeper impacts within the native material are contained primarily to the southwestern corner of the site where groundwater impacts were also observed. It is expected that impacted groundwater, identified within the upper bedrock, near the soil-bedrock interface, will be remediation through the removal of impacted soil and underlying bedrock.

The remediation program is underway, with most of the soil having been removed from the site for disposal at a licenced landfill site or for beneficial reuse at a Class 1 Management Site or Reuse Site, in accordance with O.Reg. 406/19.

This Phase II ESA report will be updated with the findings of the remediation program once post-remediation groundwater monitoring has been completed, in accordance with O.Reg. 153/04, to support the filing of a Record of Site Condition.

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7.0 STATEMENT OF LIMITATIONS

This Phase II - Environmental Site Assessment report has been prepared under the supervision of a Qualified Person, in general accordance with O. Reg 153/04. 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 Claridge Homes. Notification from Claridge Homes and Paterson Group will be required to release this report to any other party.

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Paterson Group Inc.

Mohammed Ramadan, B.Sc.

Karyn Munch, P.Eng., Q.P.ESA

Claridge HomesPaterson Group

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FIGURES

DRAWING PE2709-11 - TEST HOLE LOCATION PLAN

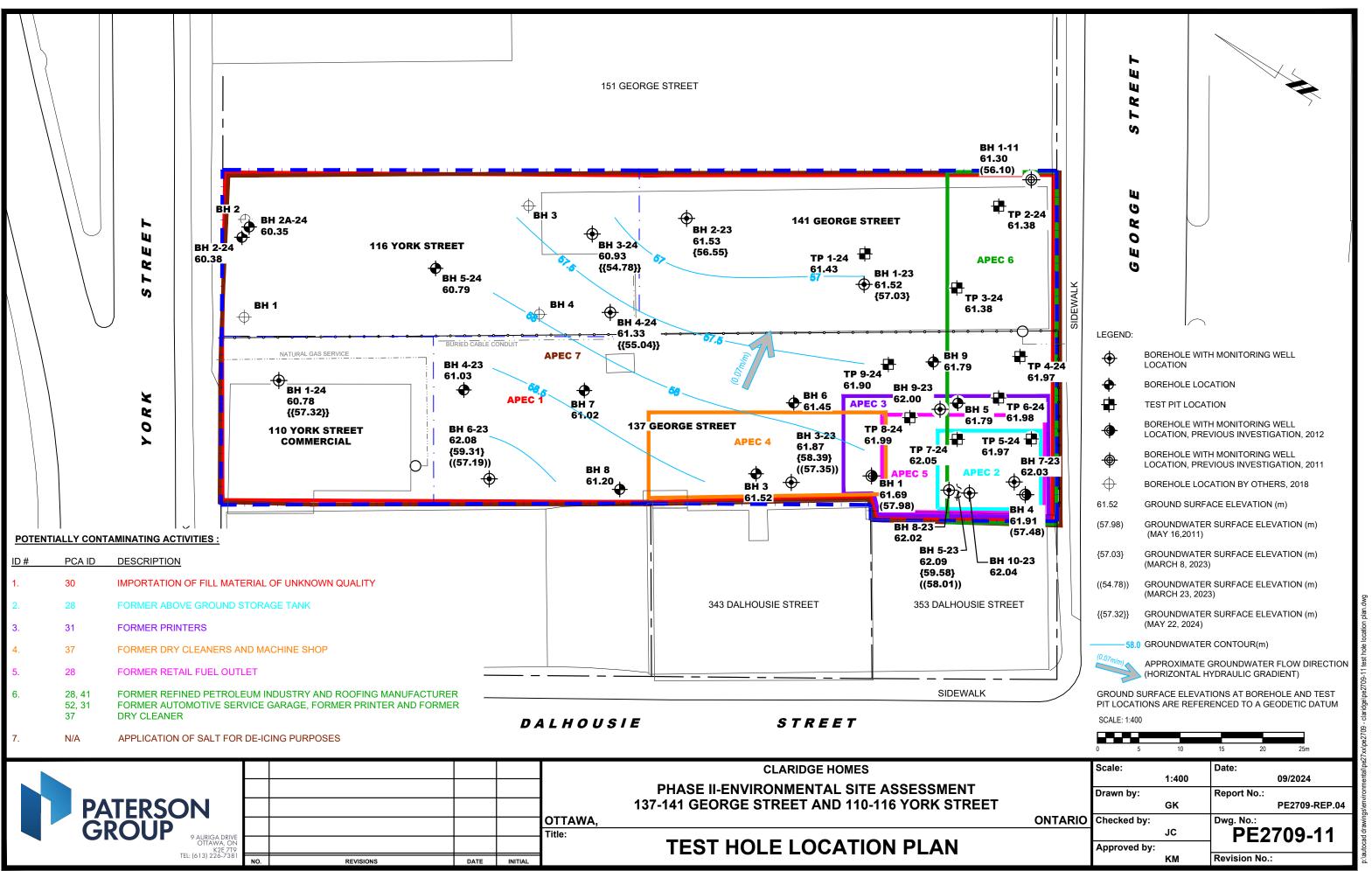
- DRAWING PE2709-12 ANALYTICAL TESTING PLAN SOIL-METALS DRAWING PE2709-12A CROSS SECTION A-A' SOIL-METALS DRAWING PE2709-12B CROSS SECTION B-B' SOIL-METALS DRAWING PE2709-12C CROSS SECTION C-C' SOIL-METALS
 - DRAWING PE2709-13 ANALYTICAL TESTING PLAN SOIL-PAHS DRAWING PE2709-13A CROSS SECTION A-A' SOIL-PAHS DRAWING PE2709-13B CROSS SECTION B-B' SOIL-PAHS DRAWING PE2709-13C– CROSS SECTION C-C' SOIL-PAHS
 - DRAWING PE2709-14 ANALYTICAL TESTING PLAN SOIL-PHCS DRAWING PE2709-14A CROSS SECTION A-A' SOIL-PHCS DRAWING PE2709-14B CROSS SECTION B-B' SOIL-PHCS DRAWING PE2709-14C CROSS SECTION C-C' SOIL-PHCS
 - DRAWING PE2709-15 ANALYTICAL TESTING PLAN SOIL-BTEX DRAWING PE2709-15A CROSS SECTION A-A' SOIL-BTEX DRAWING PE2709-15B CROSS SECTION B-B' SOIL-BTEX DRAWING PE2709-15C CROSS SECTION C-C' SOIL-BTEX
 - DRAWING PE2709-16 ANALYTICAL TESTING PLAN SOIL-VOCs DRAWING PE2709-16A CROSS SECTION A-A' SOIL-VOCs DRAWING PE2709-16B CROSS SECTION B-B' SOIL-VOCs DRAWING PE2709-16C CROSS SECTION C-C' SOIL-VOCs
 - DRAWING PE2709-17 ANALYTICAL TESTING PLAN GROUNDWATER-BTEX, VOCs
- DRAWING PE2709-17A CROSS-SECTION A A' GROUNDWATER-BTEX, VOCs
- DRAWING PE2709-17B CROSS-SECTION B B' GROUNDWATER-BTEX. VOCs
- DRAWING PE2709-17C CROSS-SECTION C C' GROUNDWATER-BTEX, VOCs

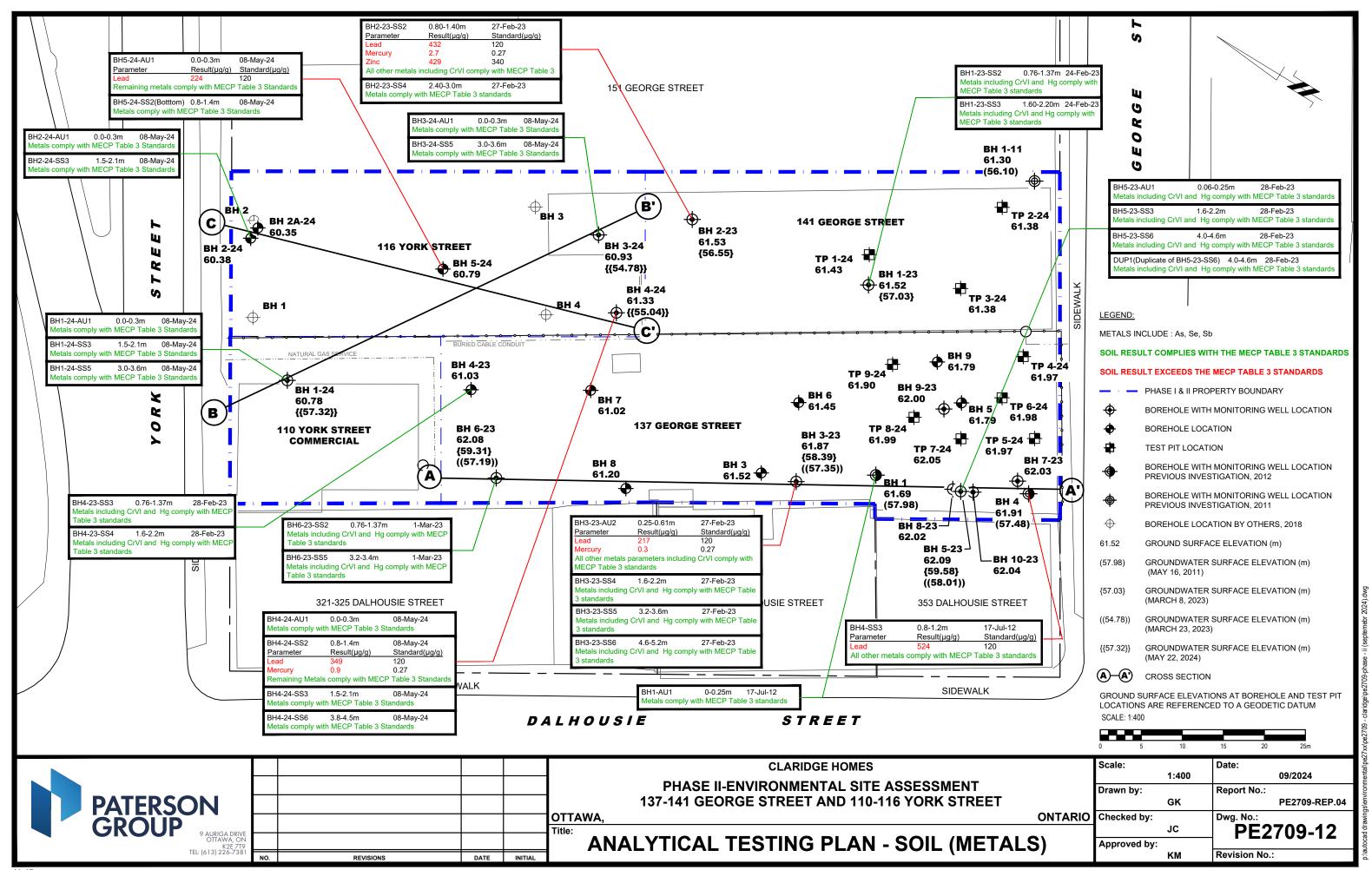
DRAWING PE2709-18 – ANALYTICAL TESTING PLAN – GROUNDWATER-METALS, HG, CRVI, PAHS

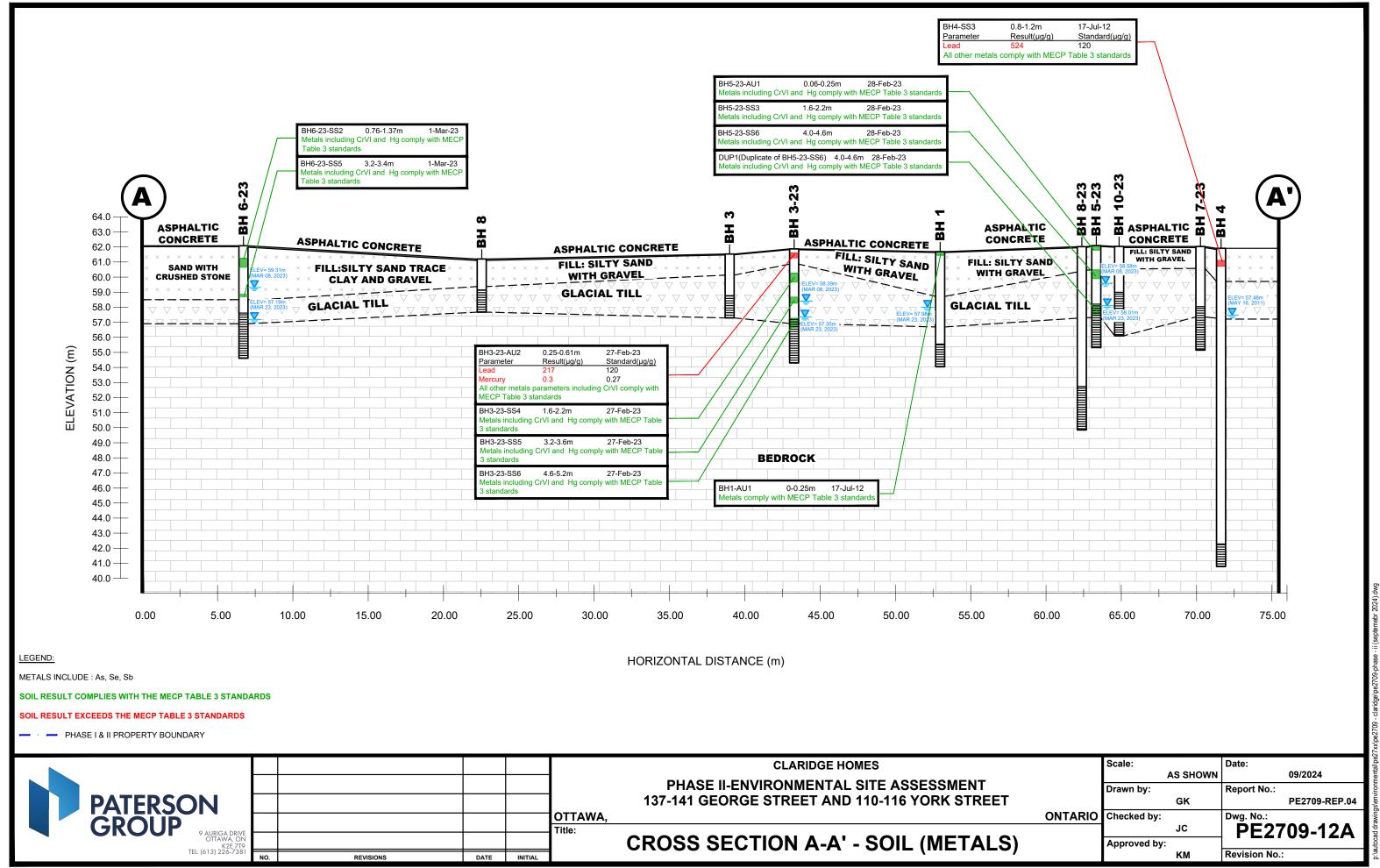
- DRAWING PE2709-18A CROSS-SECTION A A' GROUNDWATER-METALS, HG, CRVI, PAHS
- DRAWING PE2709-18B CROSS-SECTION B B' GROUNDWATER-METALS, HG, CRVI, PAHS
- DRAWING PE2709-18C CROSS-SECTION C C' GROUNDWATER-METALS, HG, CRVI, PAHS

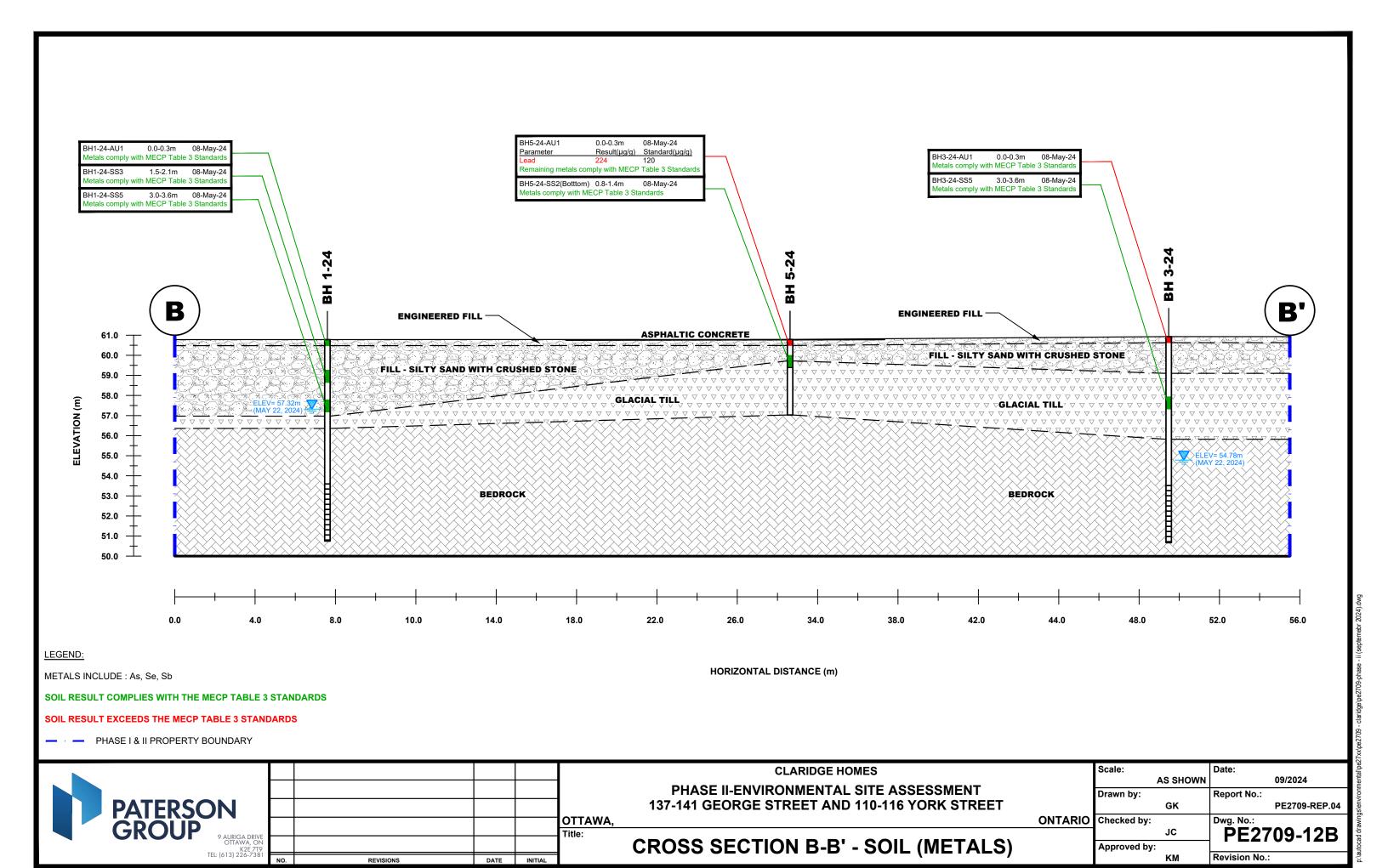
DRAWING PE2709-19 – ANALYTICAL TESTING PLAN – GROUNDWATER-PHCS

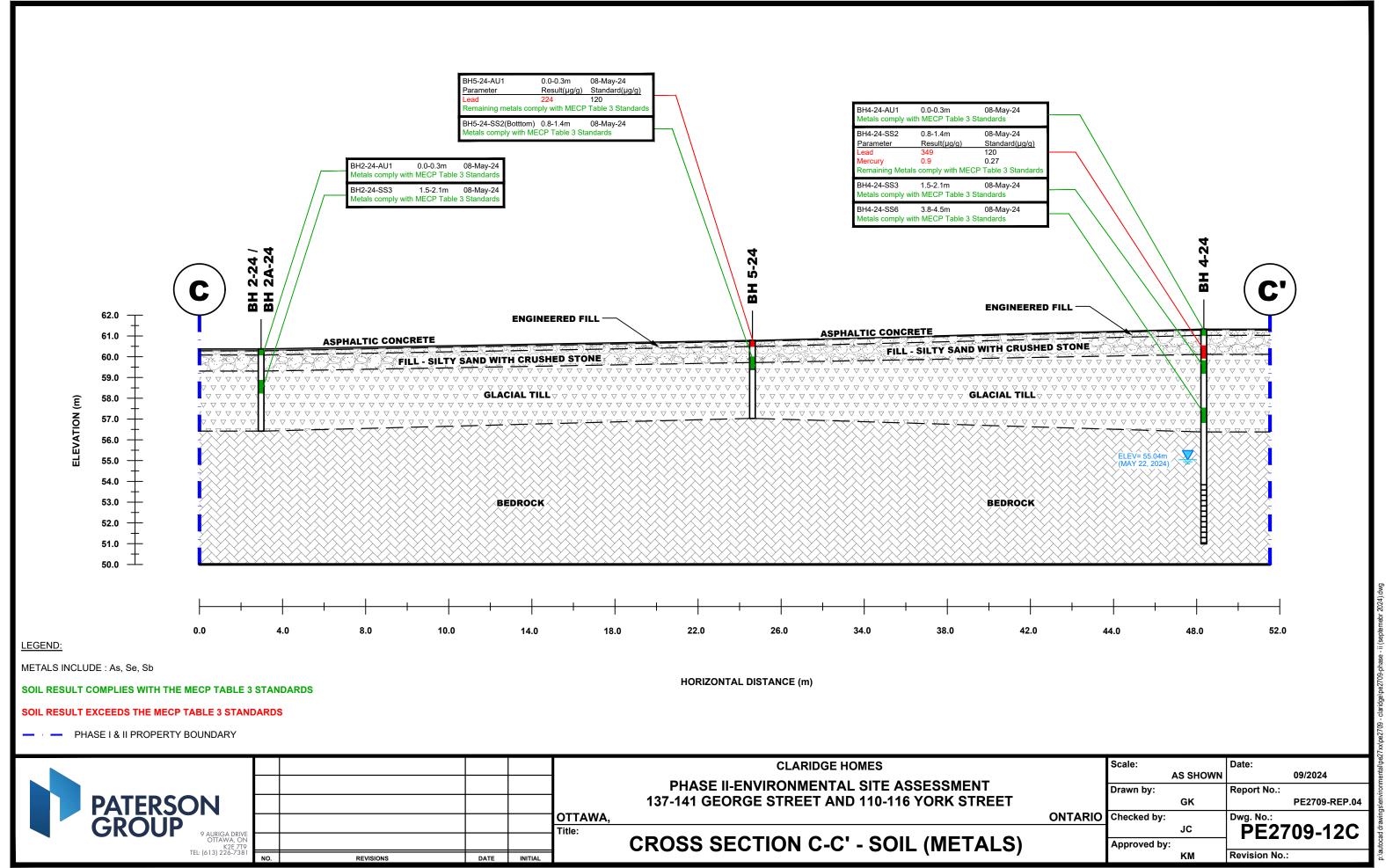
- DRAWING PE2709-19A CROSS-SECTION A A' GROUNDWATER-PHCS
- DRAWING PE2709-19B CROSS-SECTION B B' GROUNDWATER-PHCS
- DRAWING PE2709-19C CROSS-SECTION C C' GROUNDWATER-PHCS

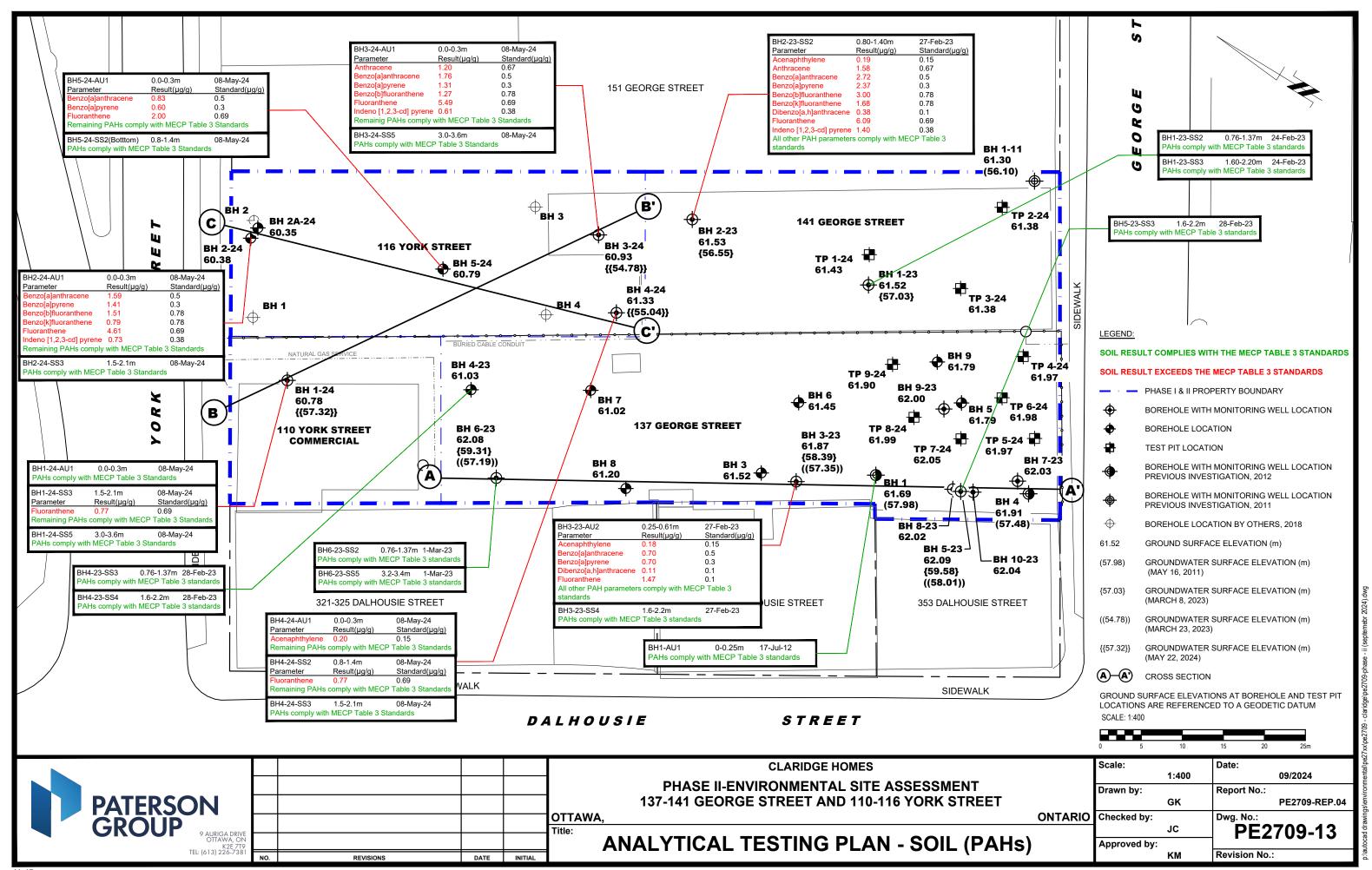


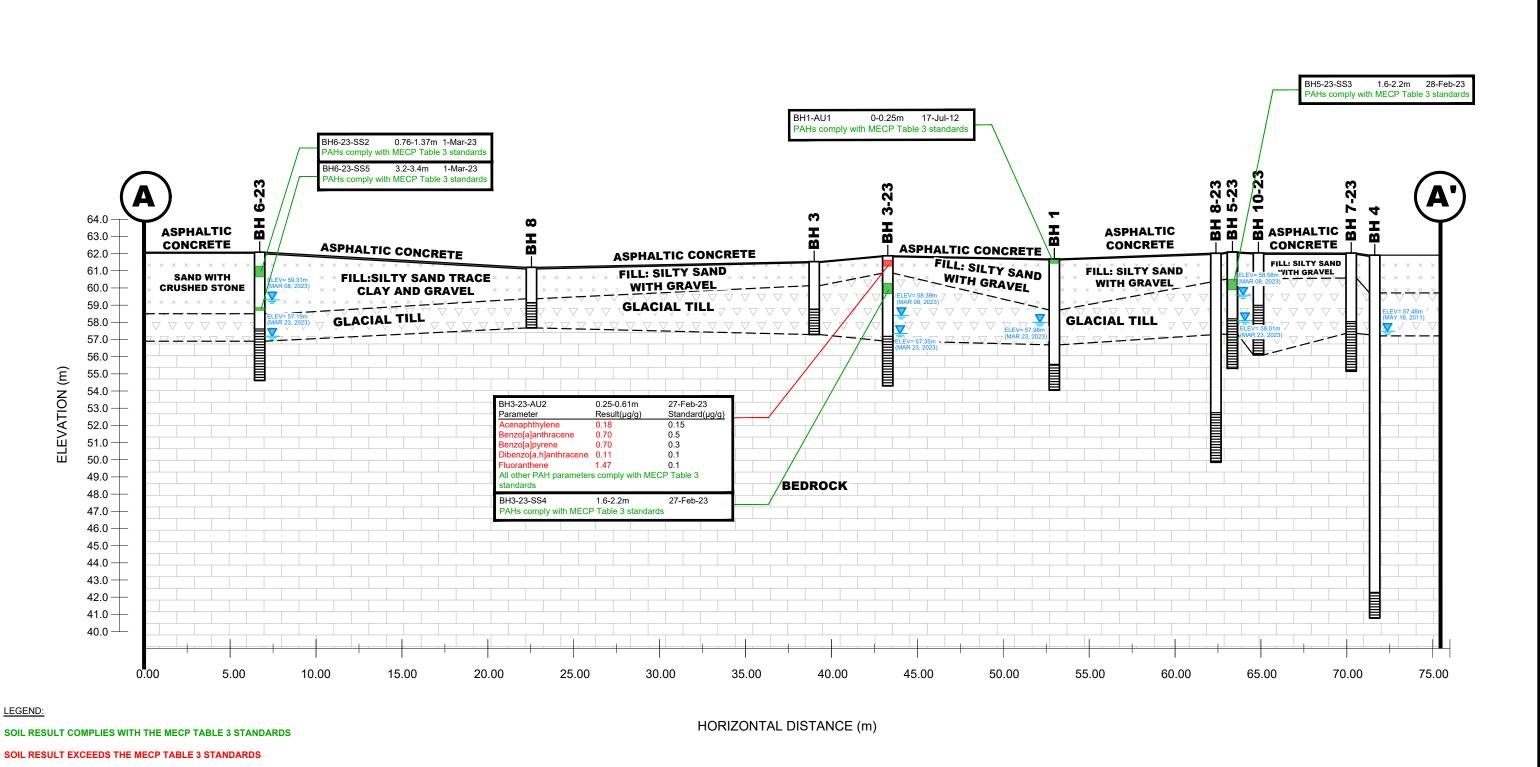






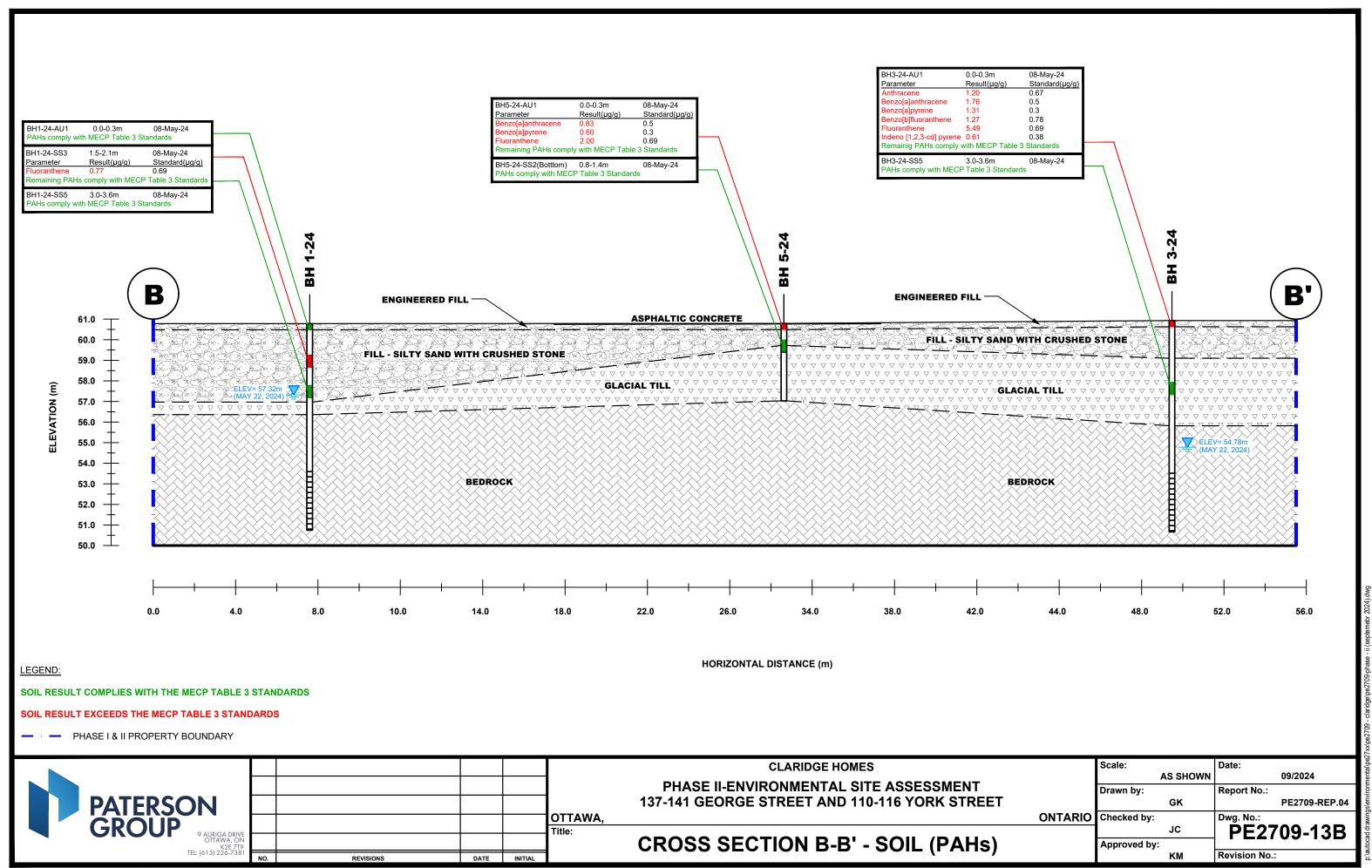


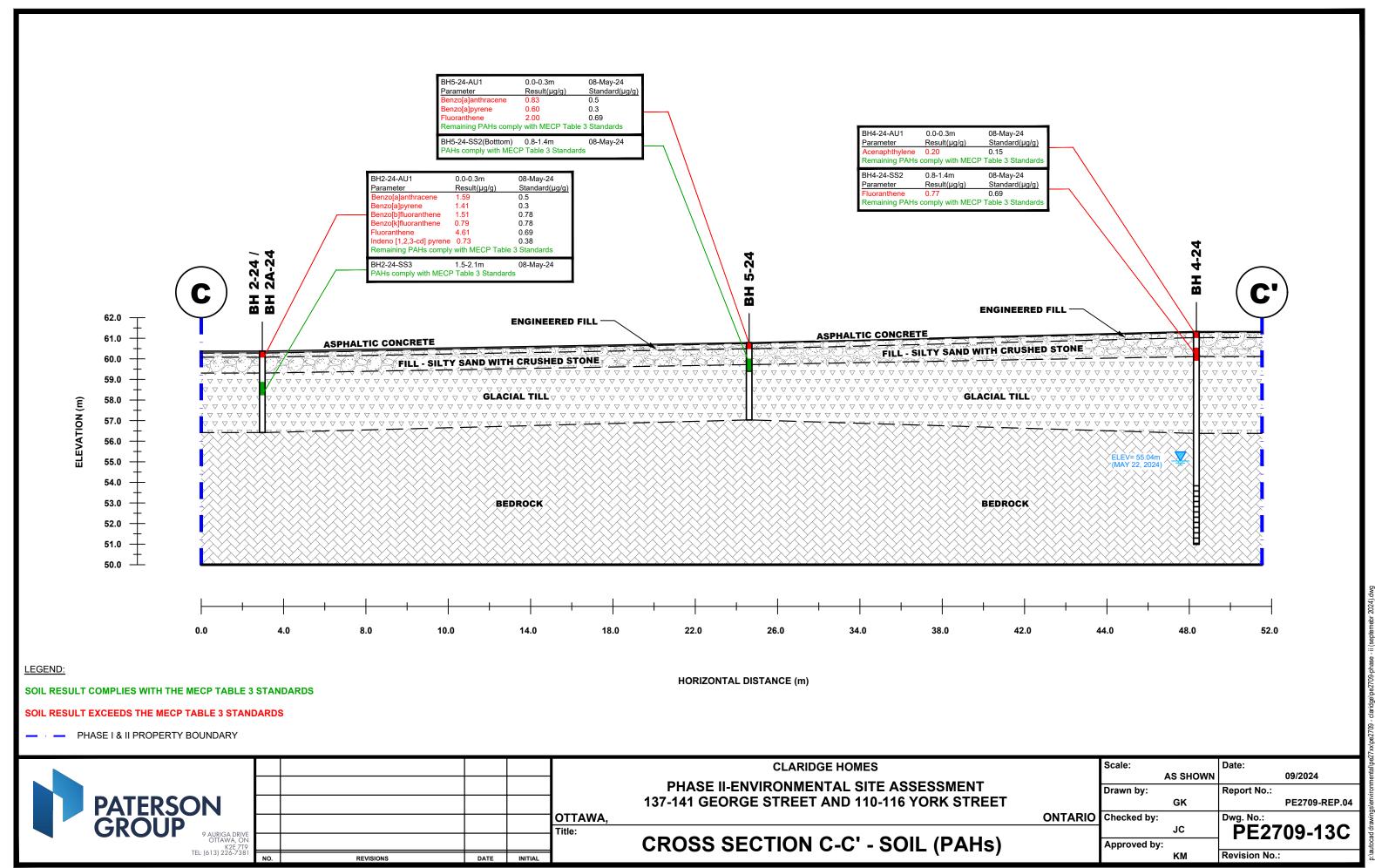


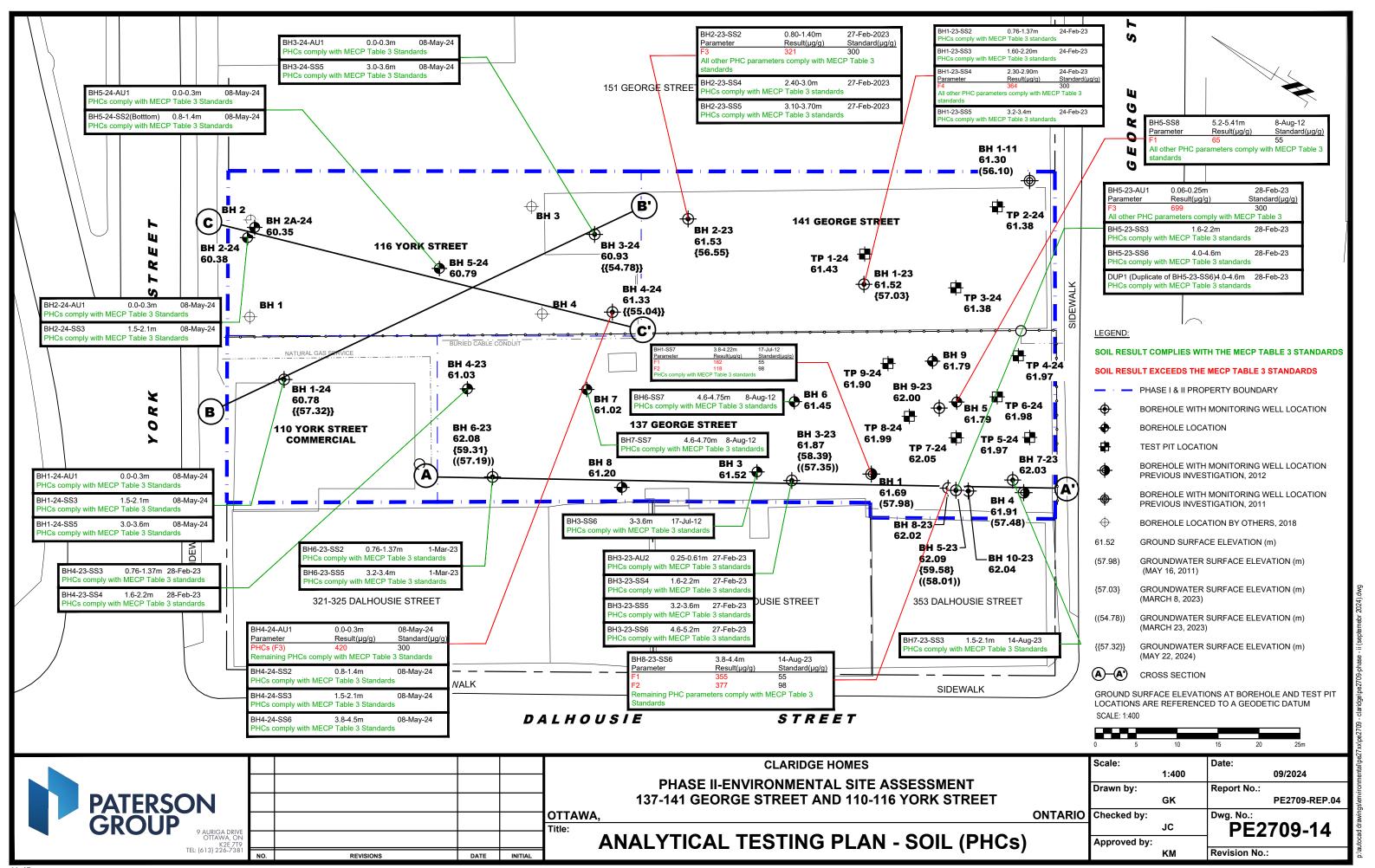


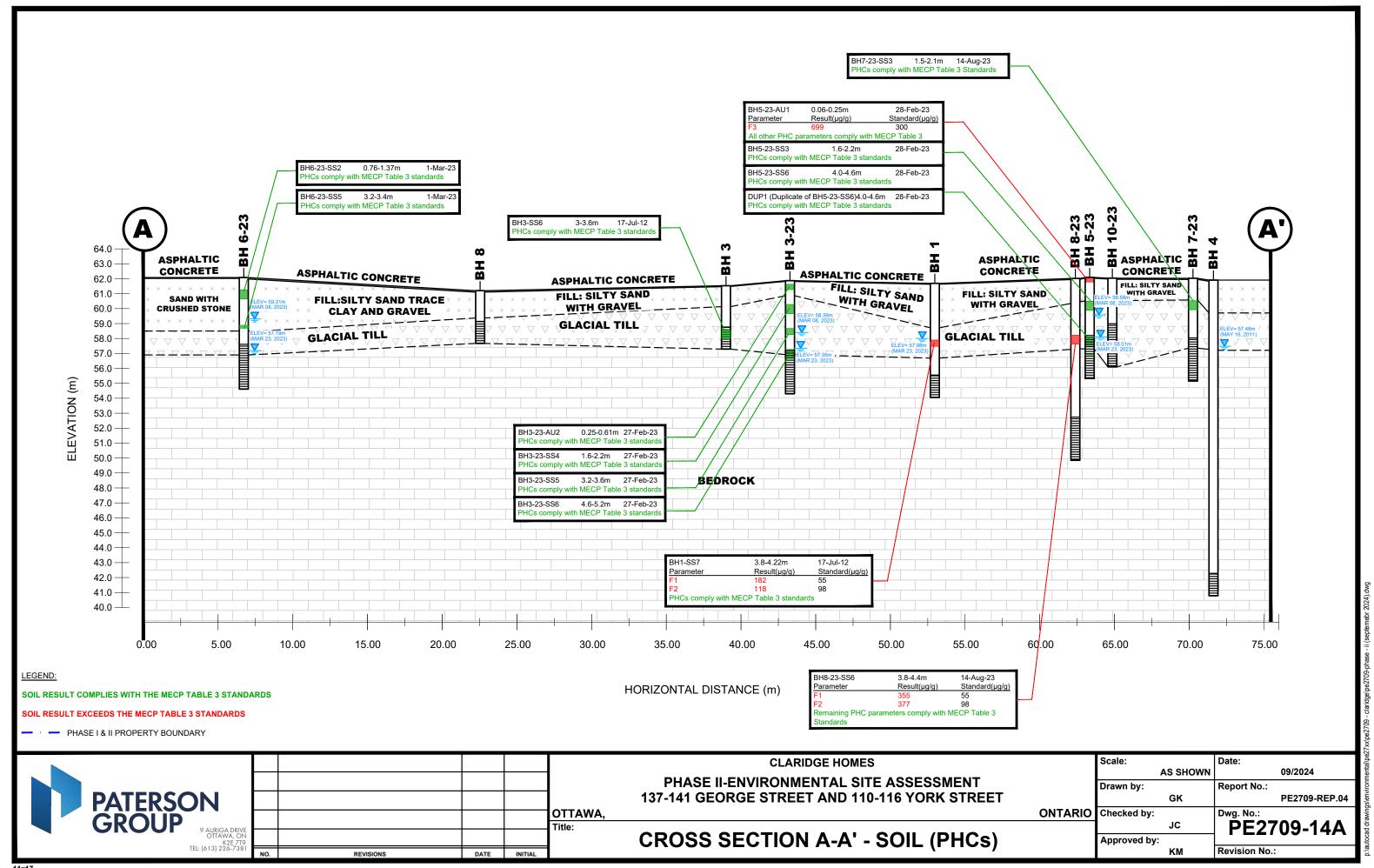
PHASE I & II PROPERTY BOUNDARY

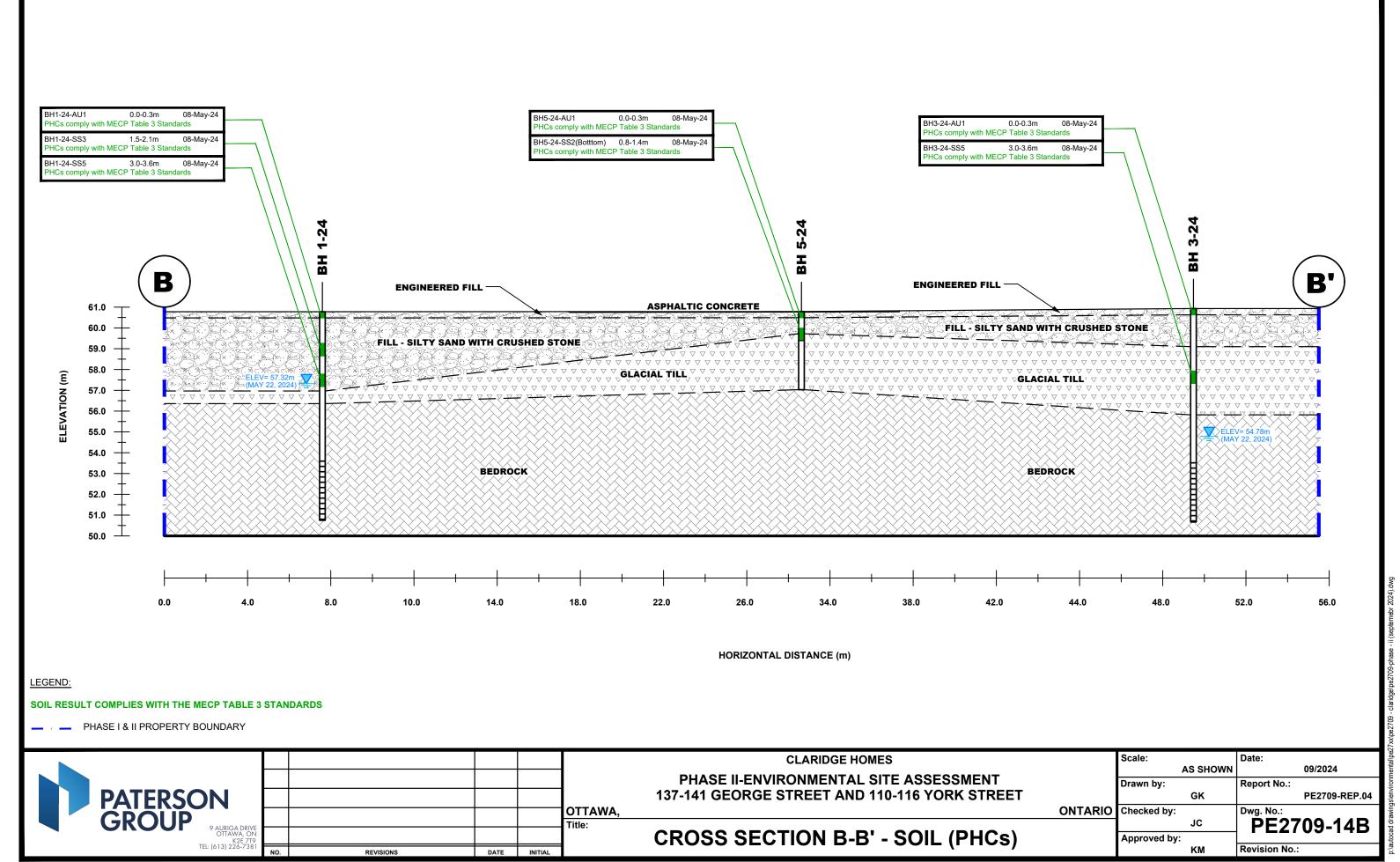
						CLARIDGE HOMES	Scale:	AS SHOWN	Date:	09/2024
PATERSON	+				OTTAWA, Title:	PHASE II-ENVIRONMENTAL SITE ASSESSMENT 137-141 GEORGE STREET AND 110-116 YORK STREET	Drawn by:	GK	Report No.:	PE2709-REP.04
GROUP 9 AURIGA DRIVE							Checked by:	JC	Dwg. No.: PE27	09-13A
OTTAWA, ON K2E 7T9 TEL: (613) 226-7381	NO.	REVISIONS	DATE	INITIAL		CROSS SECTION A-A' - SOIL (PAHs)	Approved by		Revision No.:	

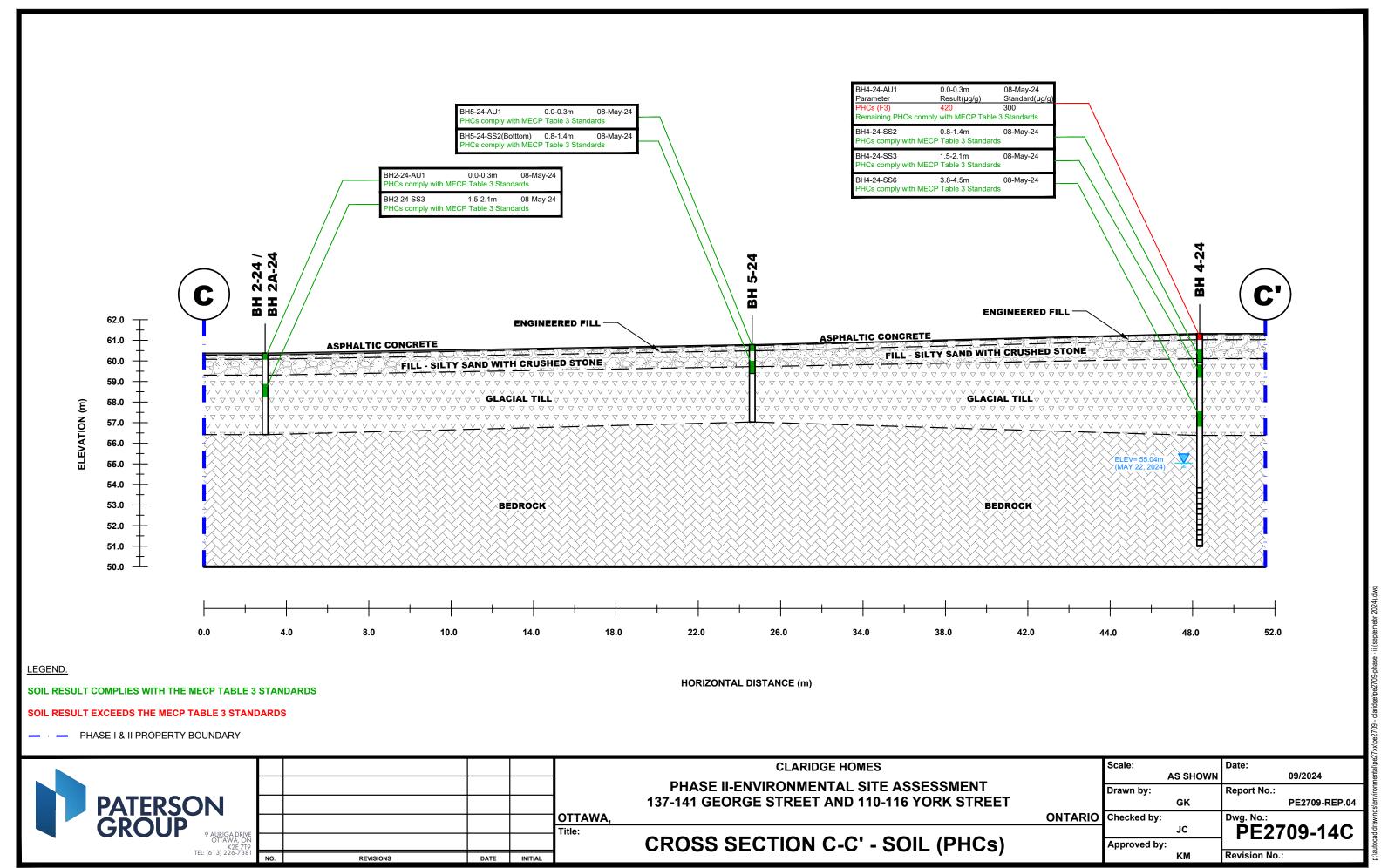


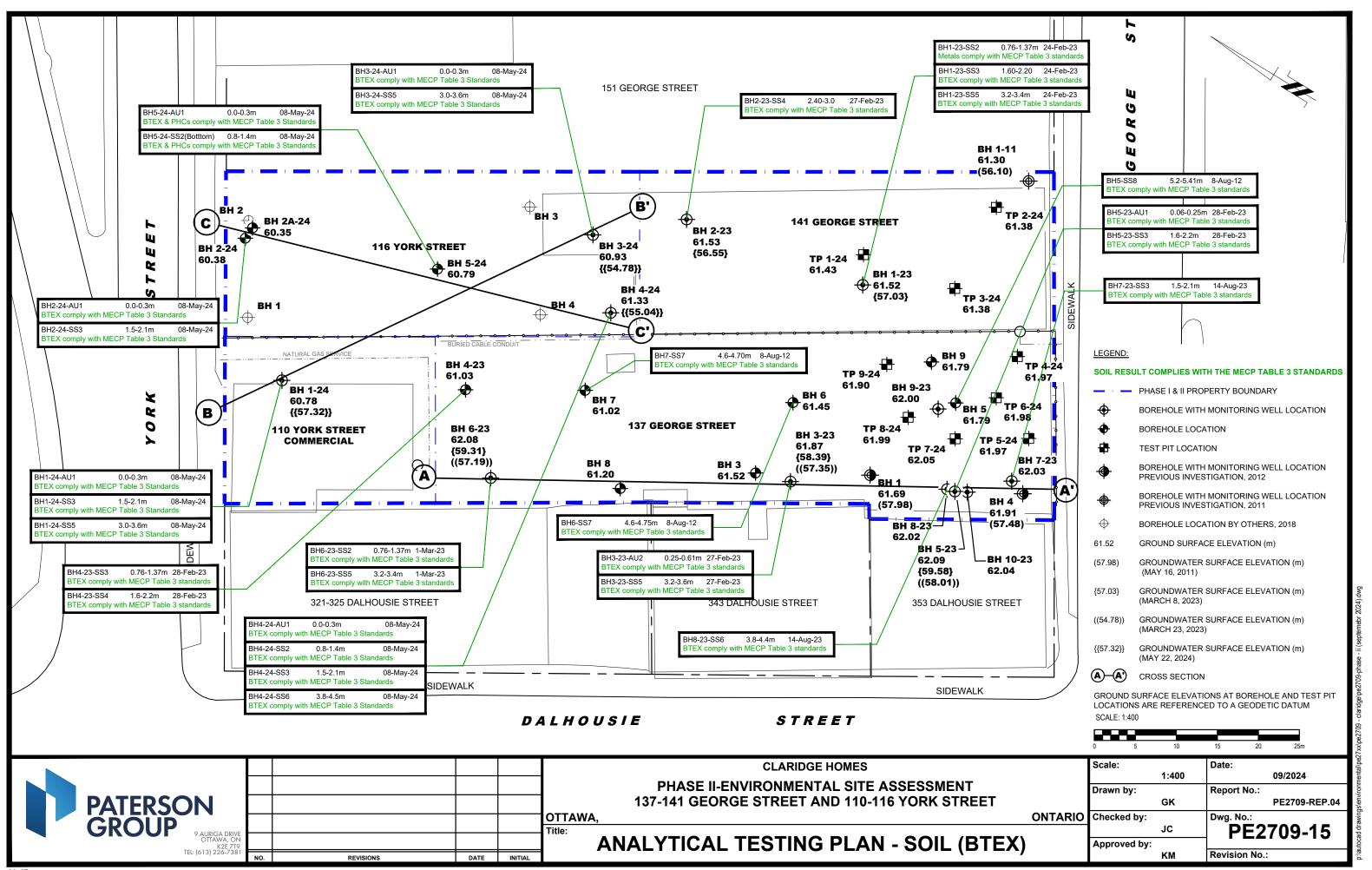


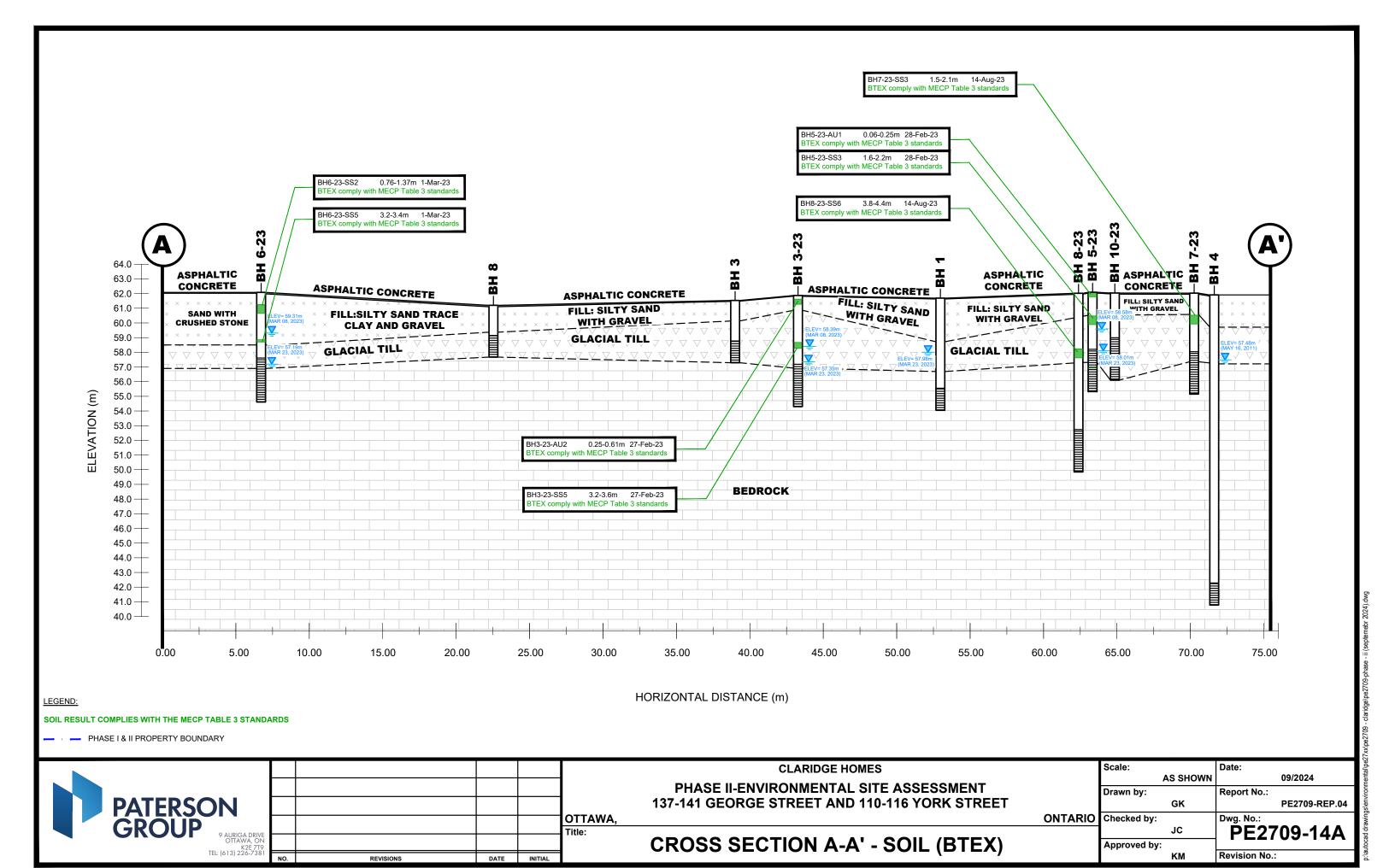


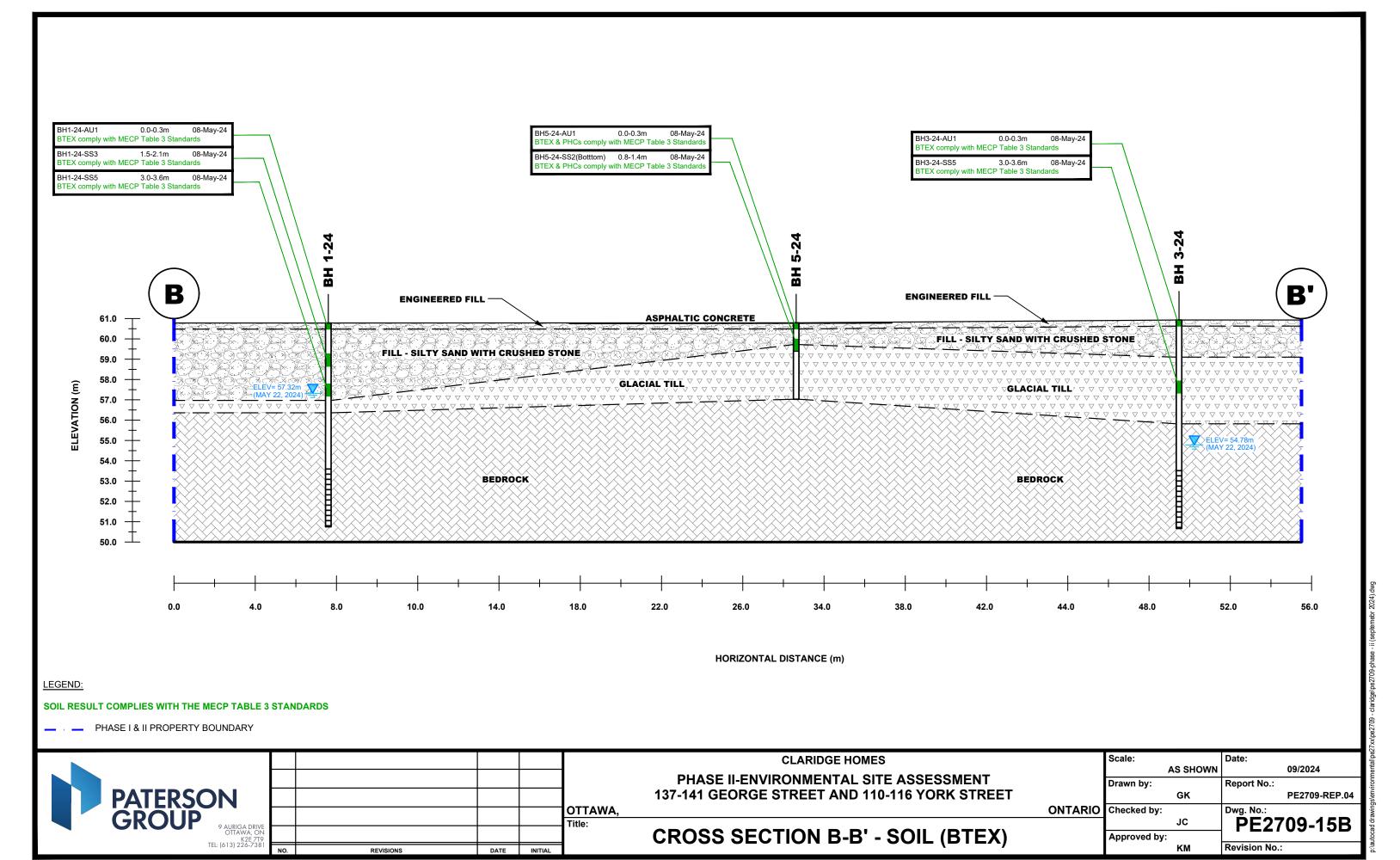


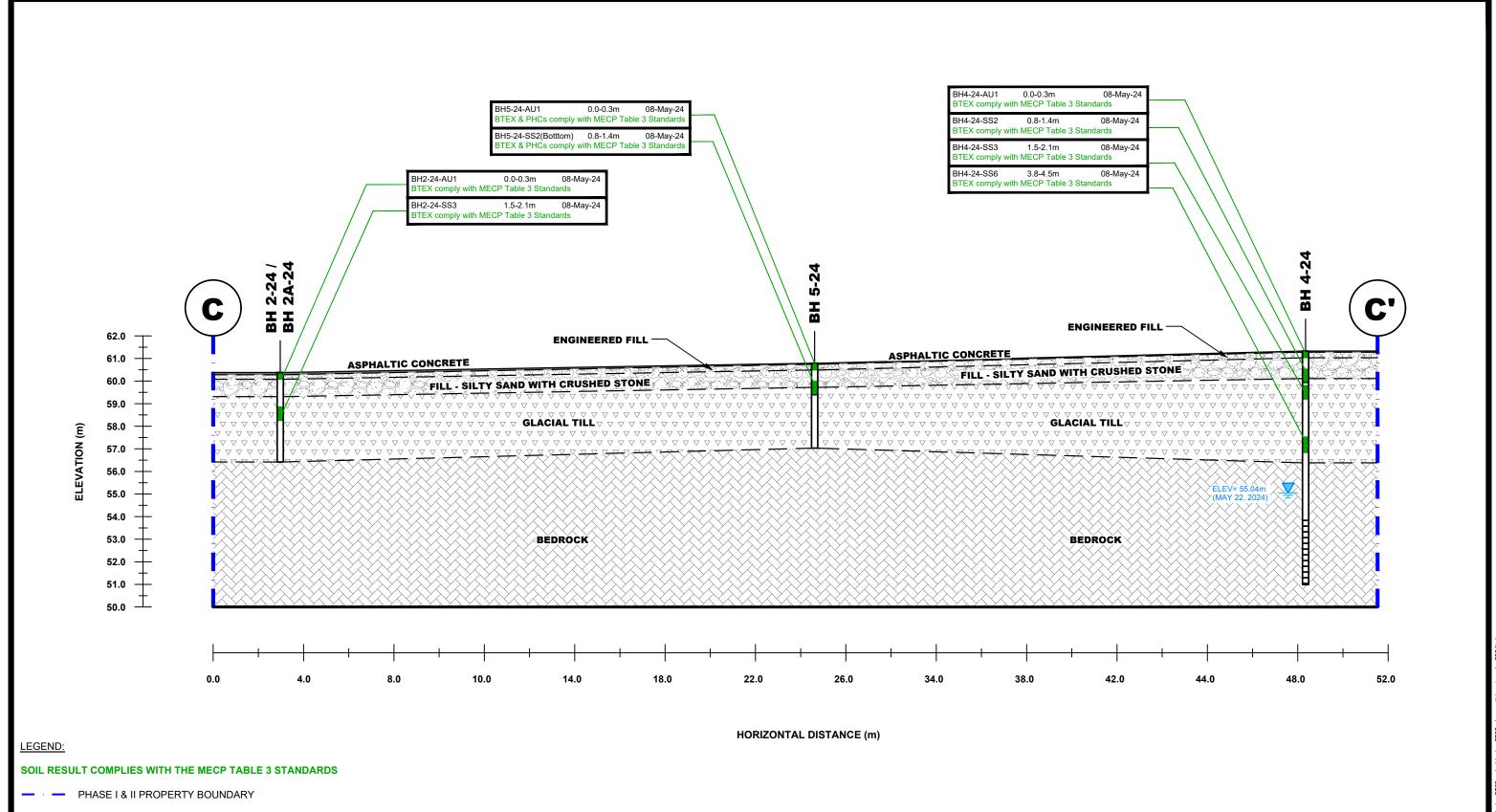




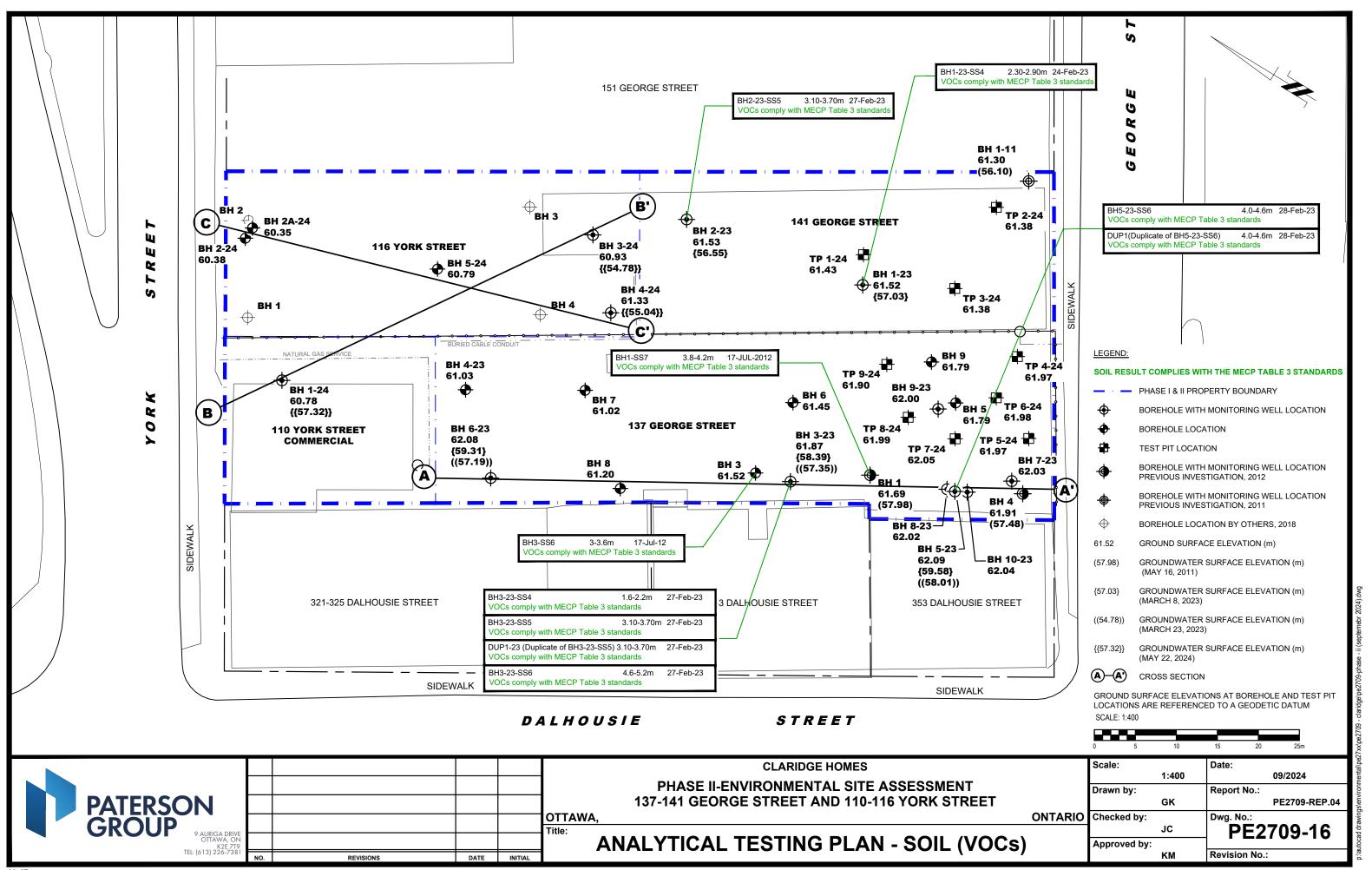


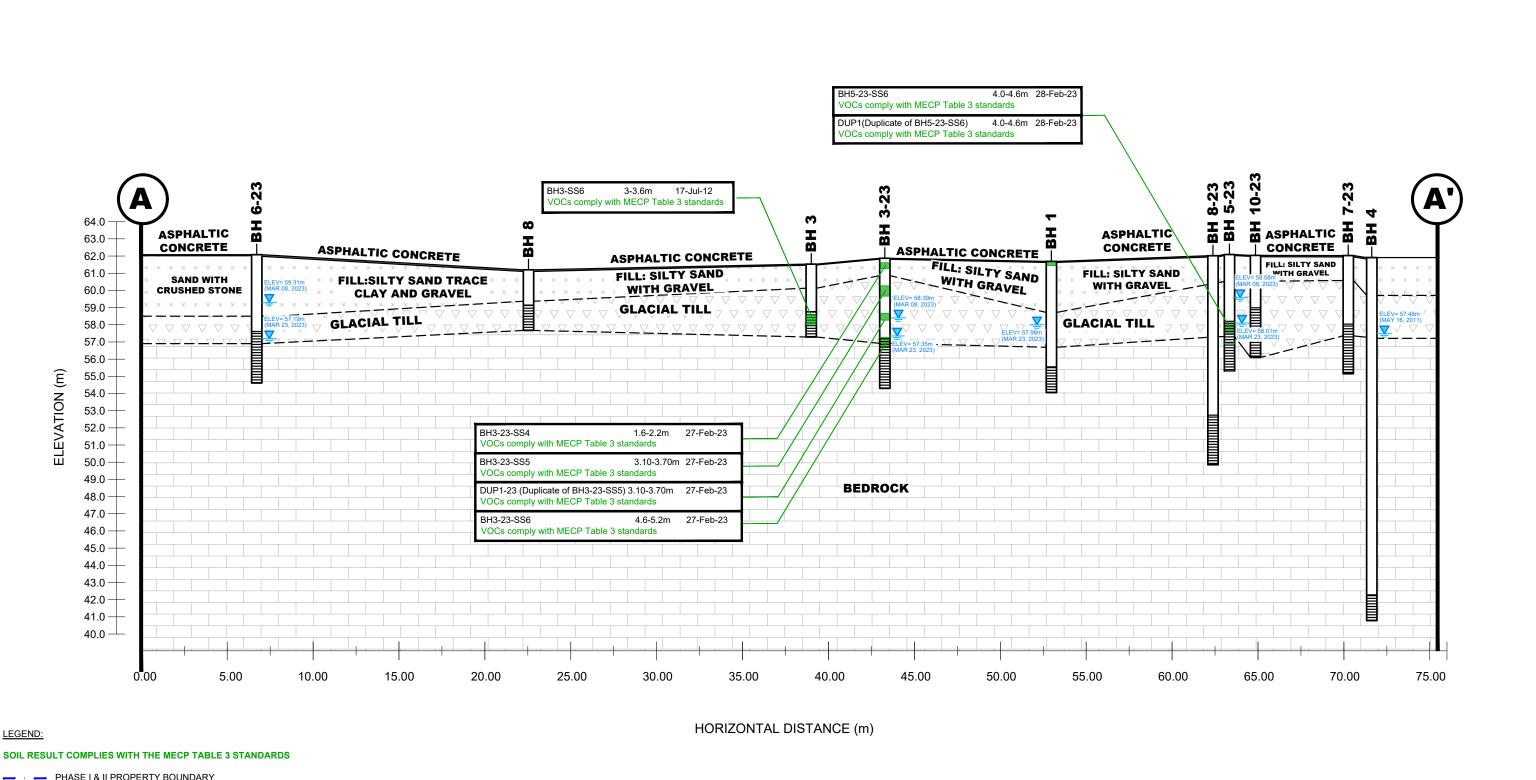






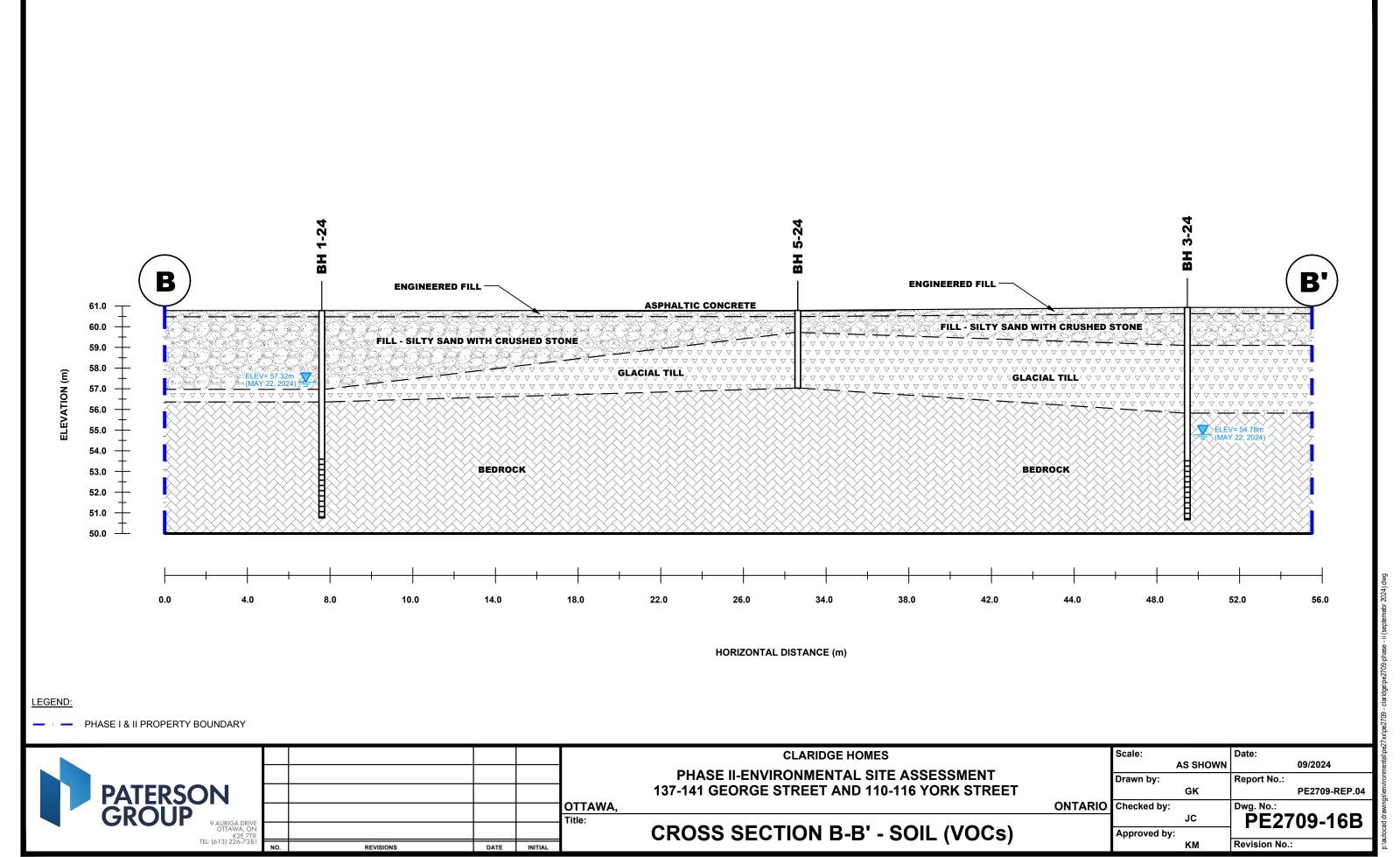
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PATERSON						PHASE II-ENVIRONMENTAL SITE ASSESSMENT 137-141 GEORGE STREET AND 110-116 YORK STREET	Drawn by:	GK	Report No.:	PE2709-REP.04
GROUP 9 AURIGA DRIVE					OTTAWA, Title:	ONTA	Checked by	/: JC	Dwg. No.: PE2709-15	′09_15C
OTTAWA, ON K2E 7T9 TEL: (613) 226-7381	NO.	REVISIONS	DATE INITIAL		CROSS SECTION C-C' - SOIL (BTEX)	Approved b	oy: KM	Revision No.:		

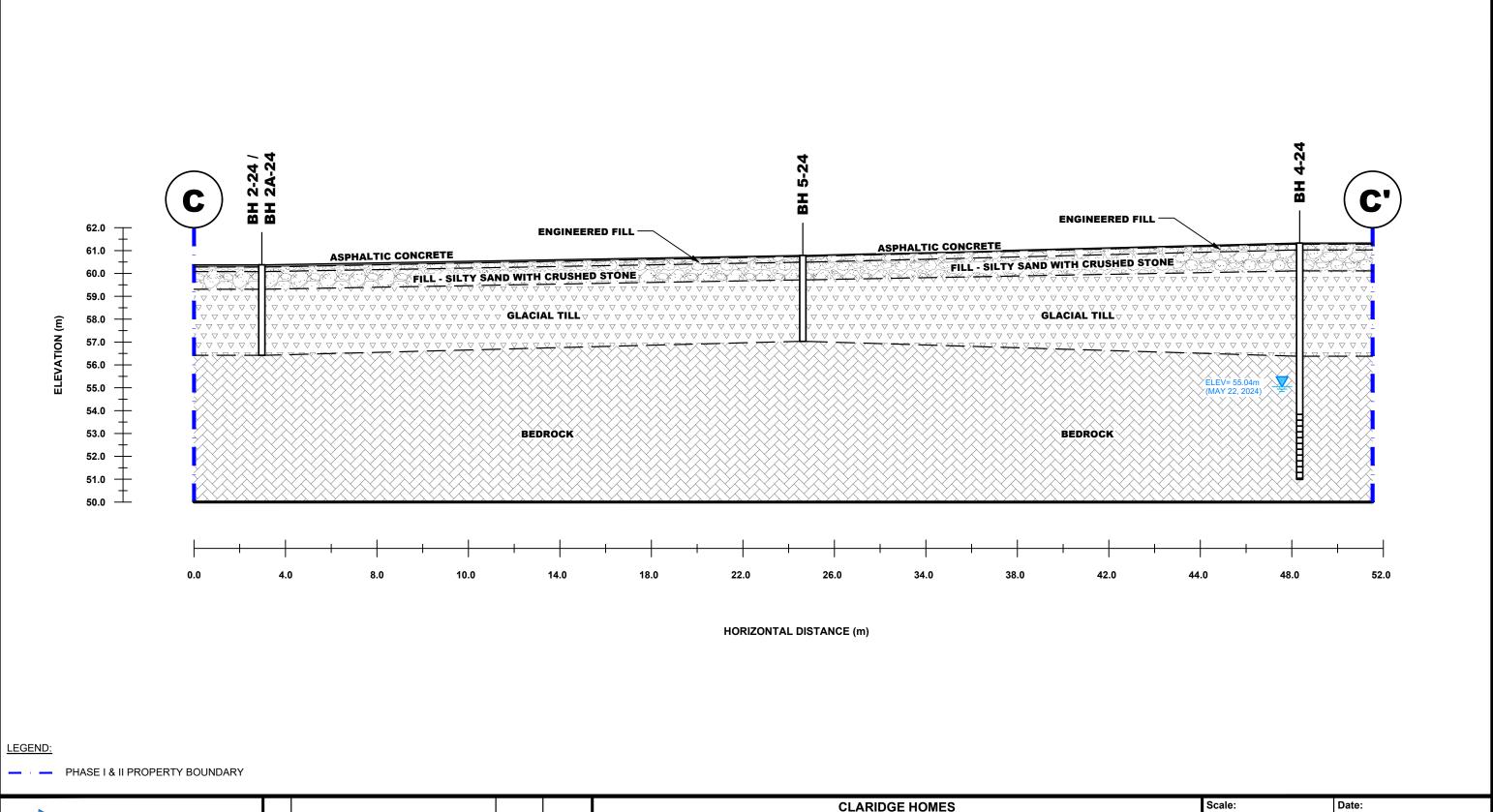




PHASE I & II PROPERTY BOUNDARY

							CLARIDGE HOMES	Scale:	AS SHOWN	Date:	09/2024
PATERSO	N F						PHASE II-ENVIRONMENTAL SITE ASSESSMENT 137-141 GEORGE STREET AND 110-116 YORK STREET	Drawn by:	GK	Report No.:	PE2709-REP.04
GROUP	9 AURIGA DRIVE					OTTAWA, Title:	ONTAI	Checked by		Dwg. No.: PE2709-16	'09_16Δ
	Y AURIGA DRIVE OTTAWA, ON K2E 7T9 TEL: (613) 226-7381	REVISIONS	DATE	INITIAL		CROSS SECTION A-A' - SOIL (VOCs)	Approved b	y:	Revision No.:		





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9 AURIGA DRIVE
OOTTAWA, ON
K22 ON
TEL: (613) 226-7381

NO.	REVISIONS	DATE	INITIAL

OTTAWA,

Title:

CLARIDGE HOMES

PHASE II-ENVIRONMENTAL SITE ASSESSMENT 137-141 GEORGE STREET AND 110-116 YORK STREET

ONTARIO Checked by:

Drawn by:

GK

Report No.:

PE2709-REP.04

Checked by:

JC

Approved by:

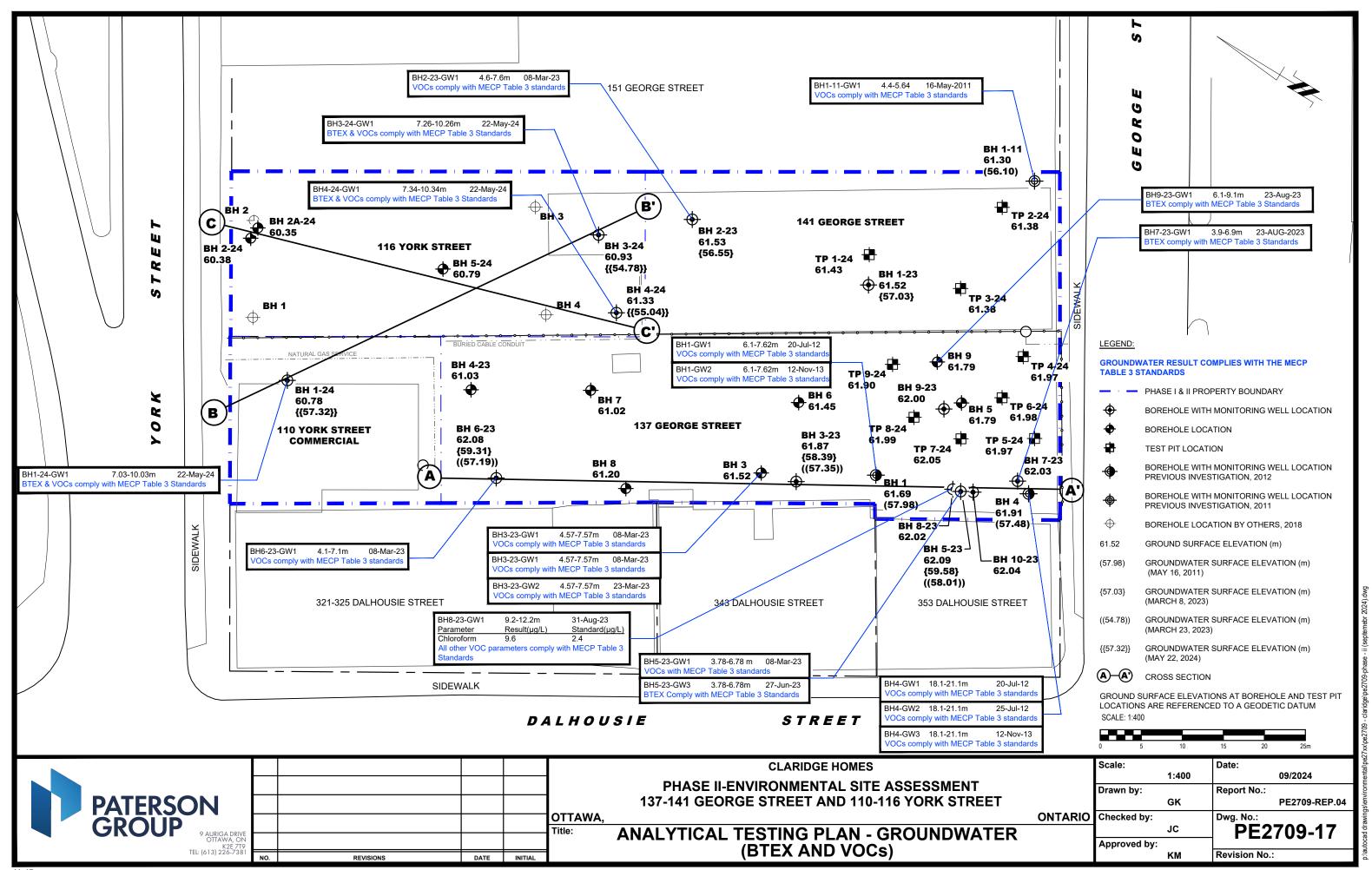
PE2709-16C

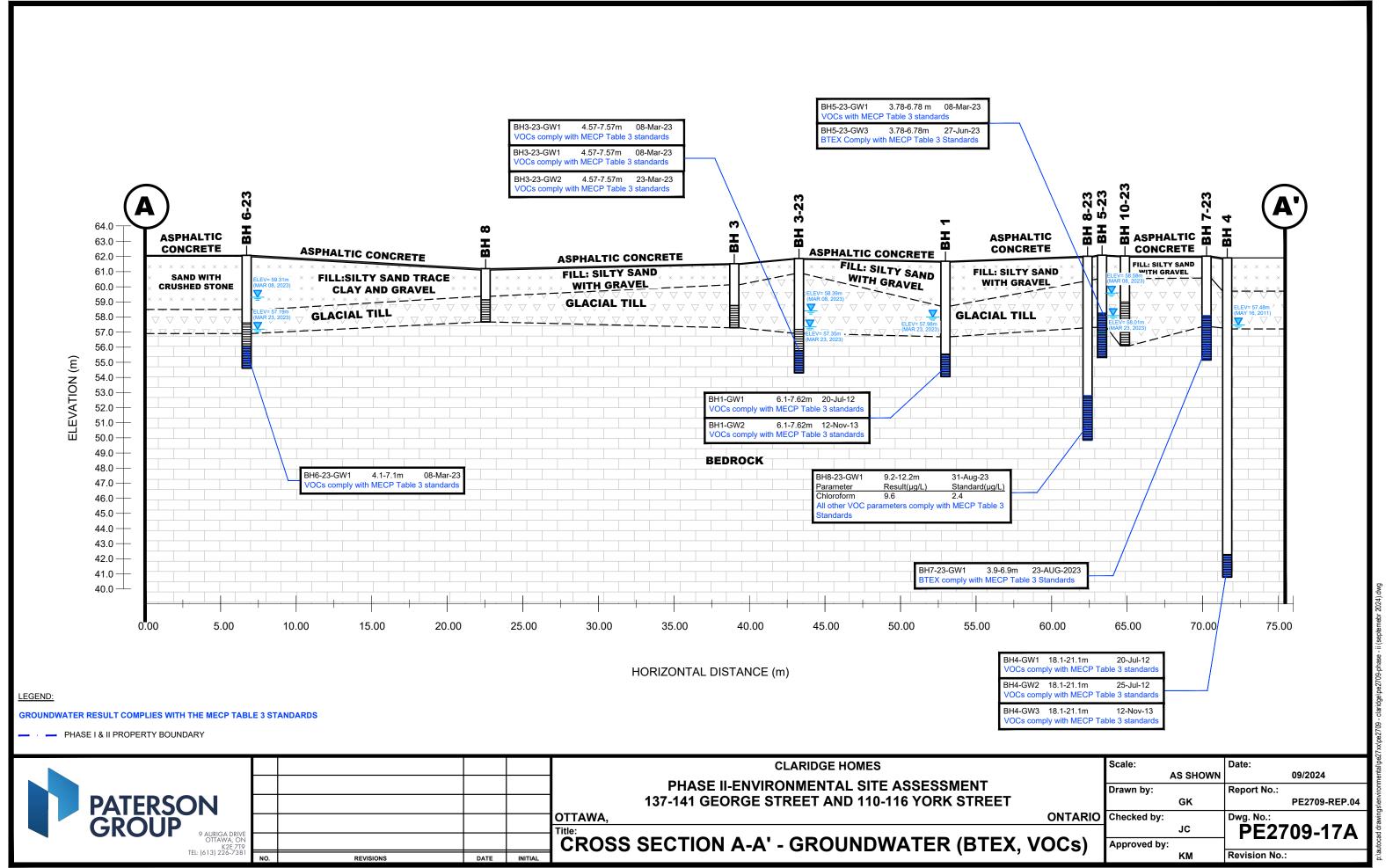
Revision No.:

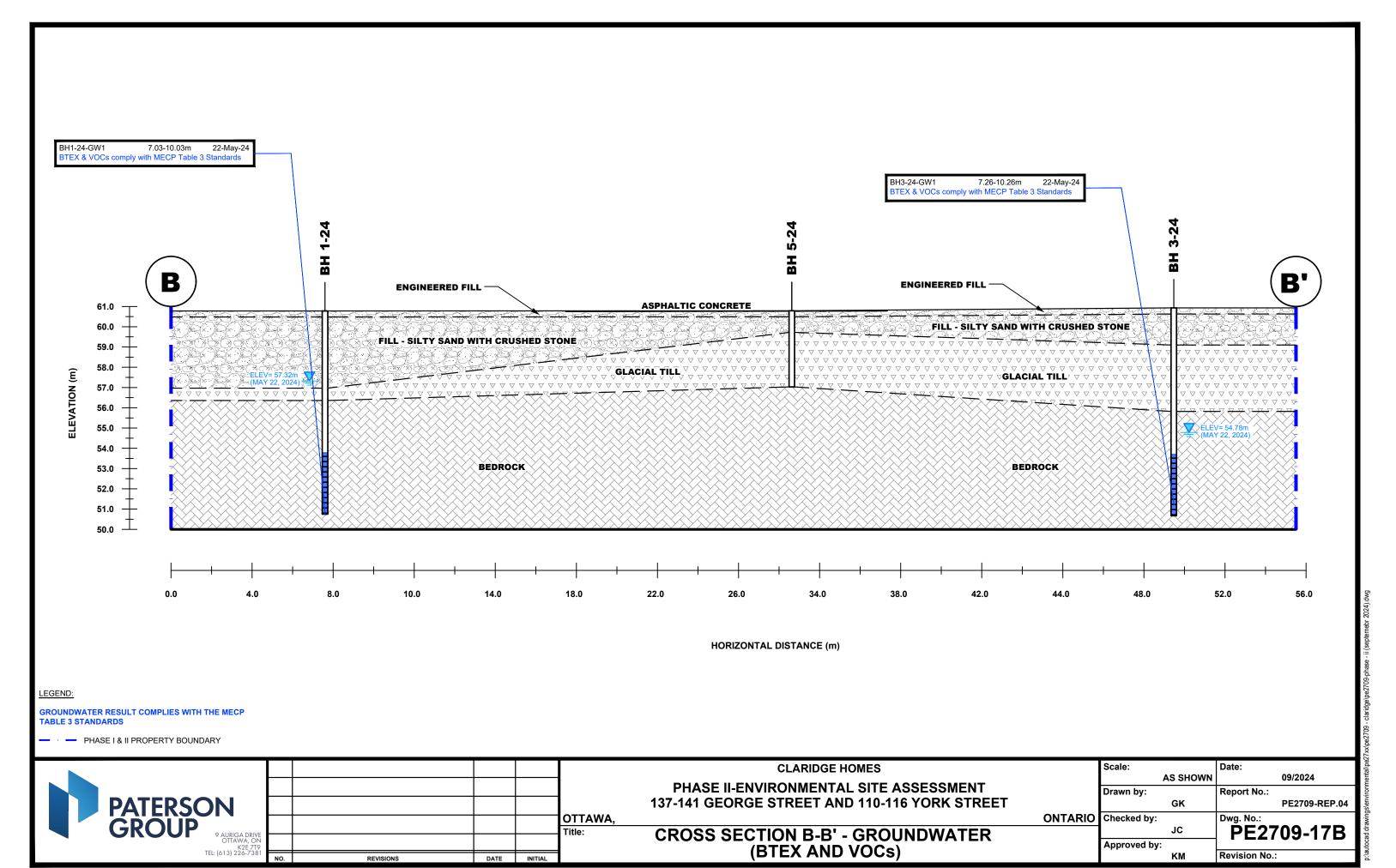
09/2024

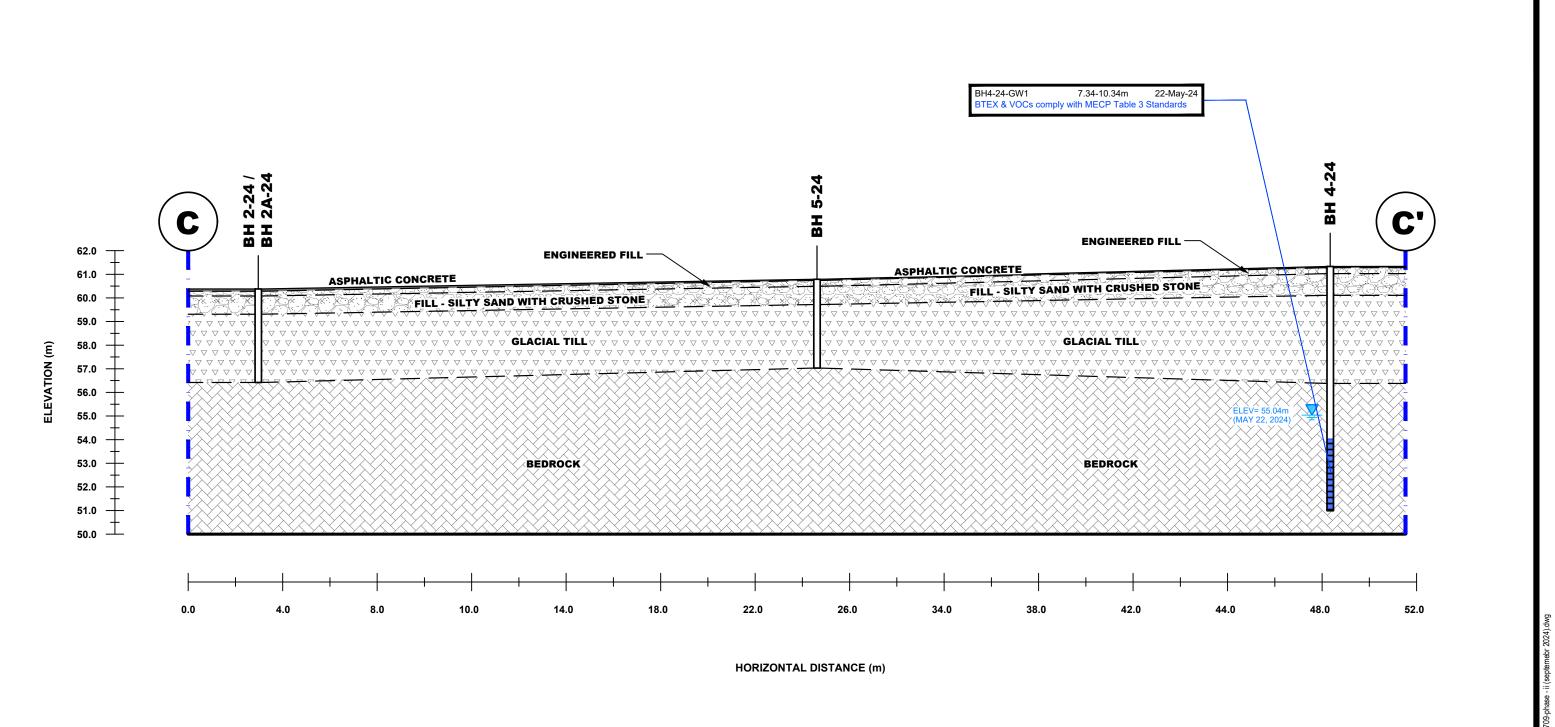
AS SHOWN

CROSS SECTION C-C' - SOIL (VOCs)







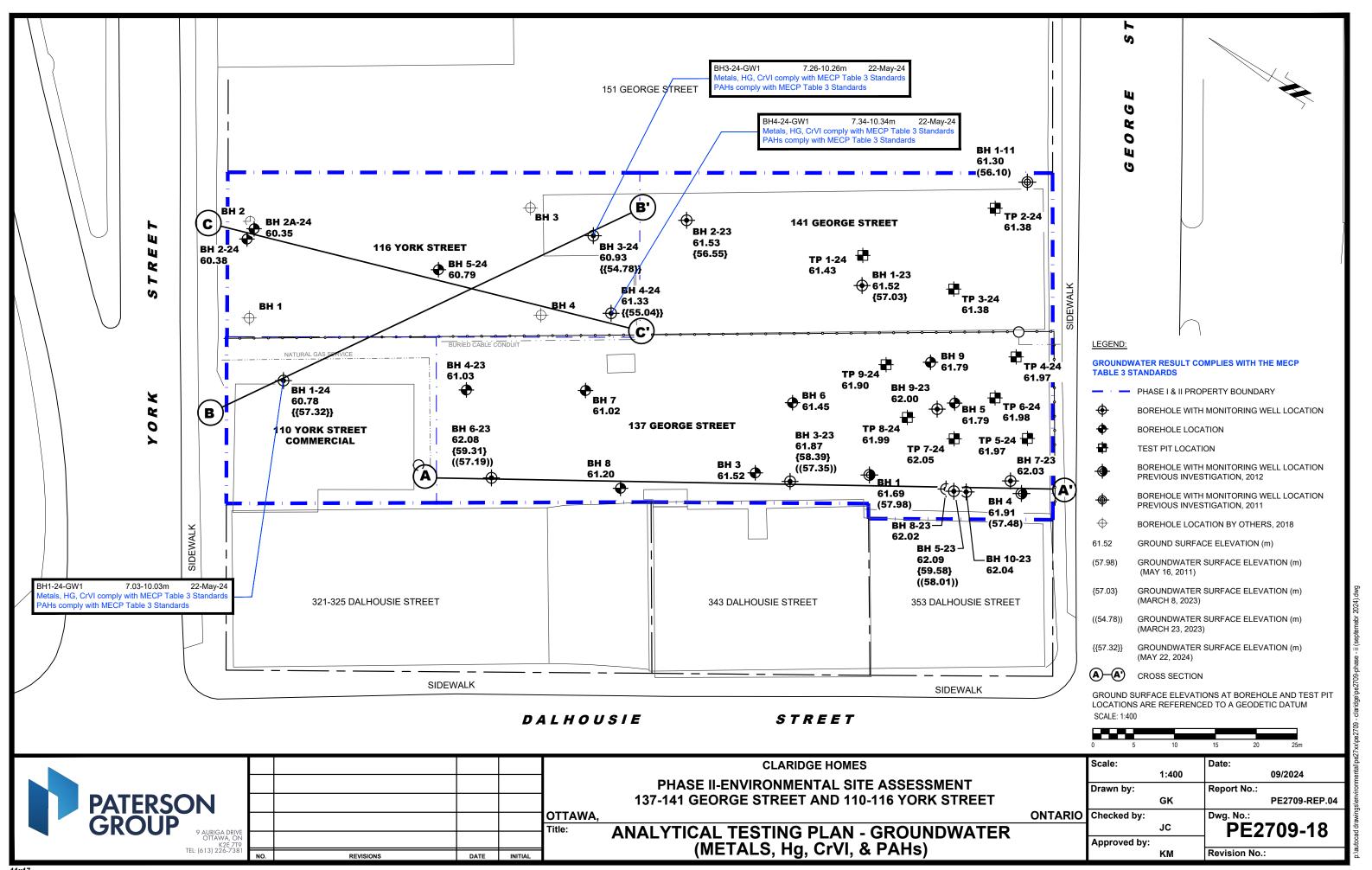


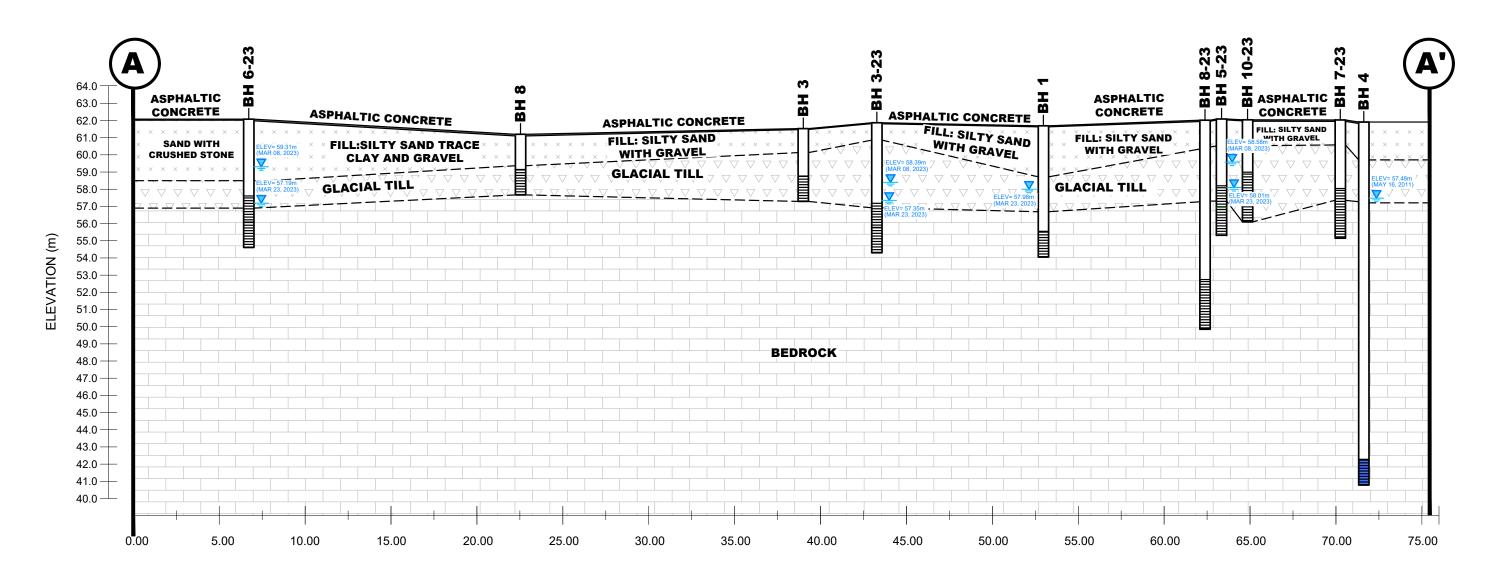
LEGEND:

GROUNDWATER RESULT COMPLIES WITH THE MECP TABLE 3 STANDARDS

- PHASE I & II PROPERTY BOUNDARY

					CLARIDGE HOMES	Scale:	40.011014/11	Date:	00/2024
					PHASE II-ENVIRONMENTAL SITE ASSESSMENT	Drawn by:	AS SHOWN	Report No.:	09/2024
PATERSON					137-141 GEORGE STREET AND 110-116 YORK STREET		GK		PE2709-REP.04
GROUP					OTTAWA, ONTARIO	Checked by:	JC	PE270	ng_17C
9 Auriga drive Ottawa, On K2E 719						Approved by			09-17 C
TEL: (613) 226-7381	NO.	REVISIONS	DATE	INITIAL	, , , , ,		KM	Revision No.:	





HORIZONTAL DISTANCE (m)

GROUNDWATER RESULT COMPLIES WITH THE MECP TABLE 3 STANDARDS

PHASE I & II PROPERTY BOUNDARY

PATERSO GROUP	
- OKOOI	9 AURIGA DRIVE OTTAWA, ON K2E 7T9 TEL: (613) 226-7381

NO.	REVISIONS	DATE	INITIAL	OTTAWA, Title:
NO.	KEVIGIONS	DAIL	INITIAL	

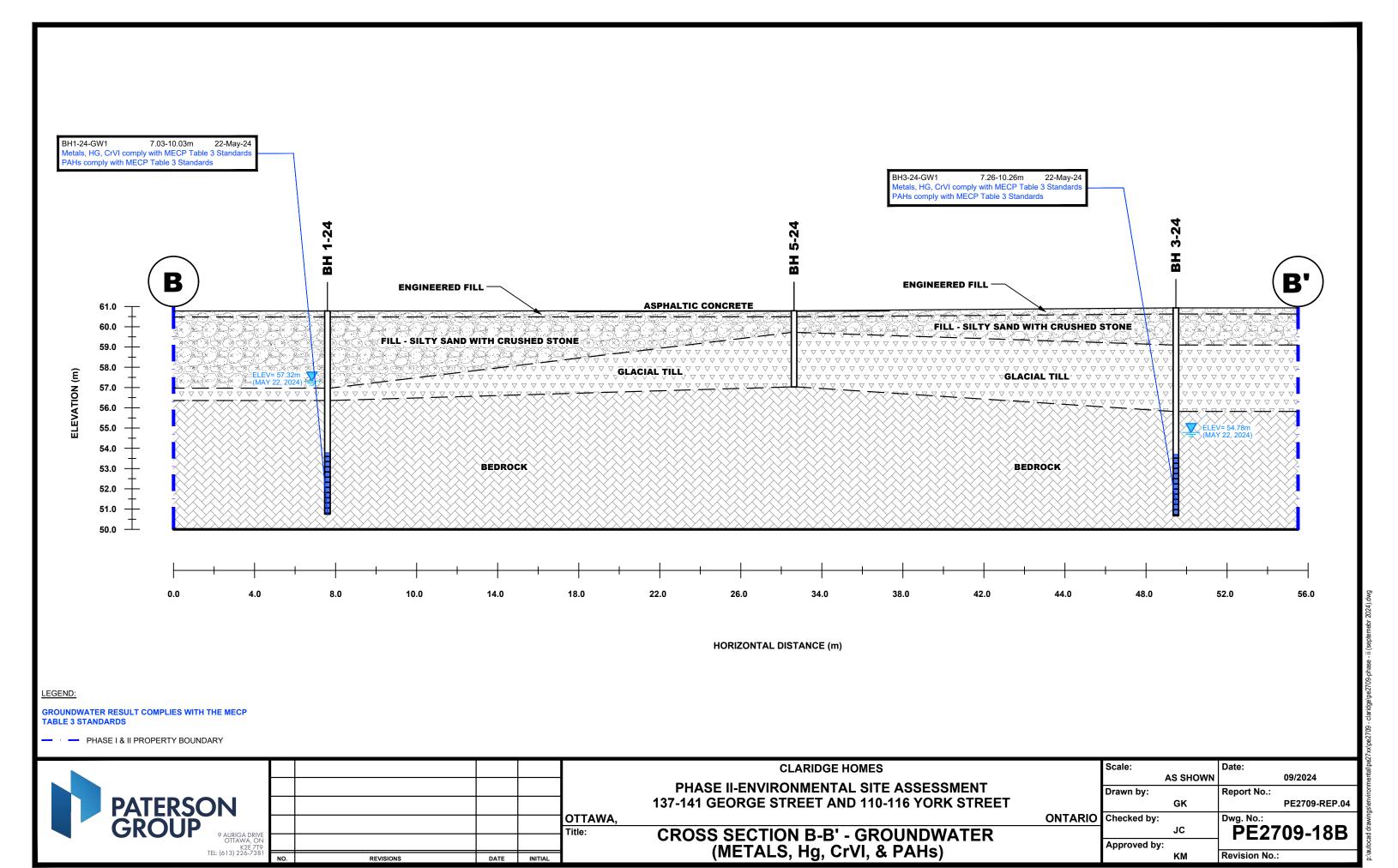
CLARIDGE HOMES PHASE II-ENVIRONMENTAL SITE ASSESSMENT 137-141 GEORGE STREET AND 110-116 YORK STREET

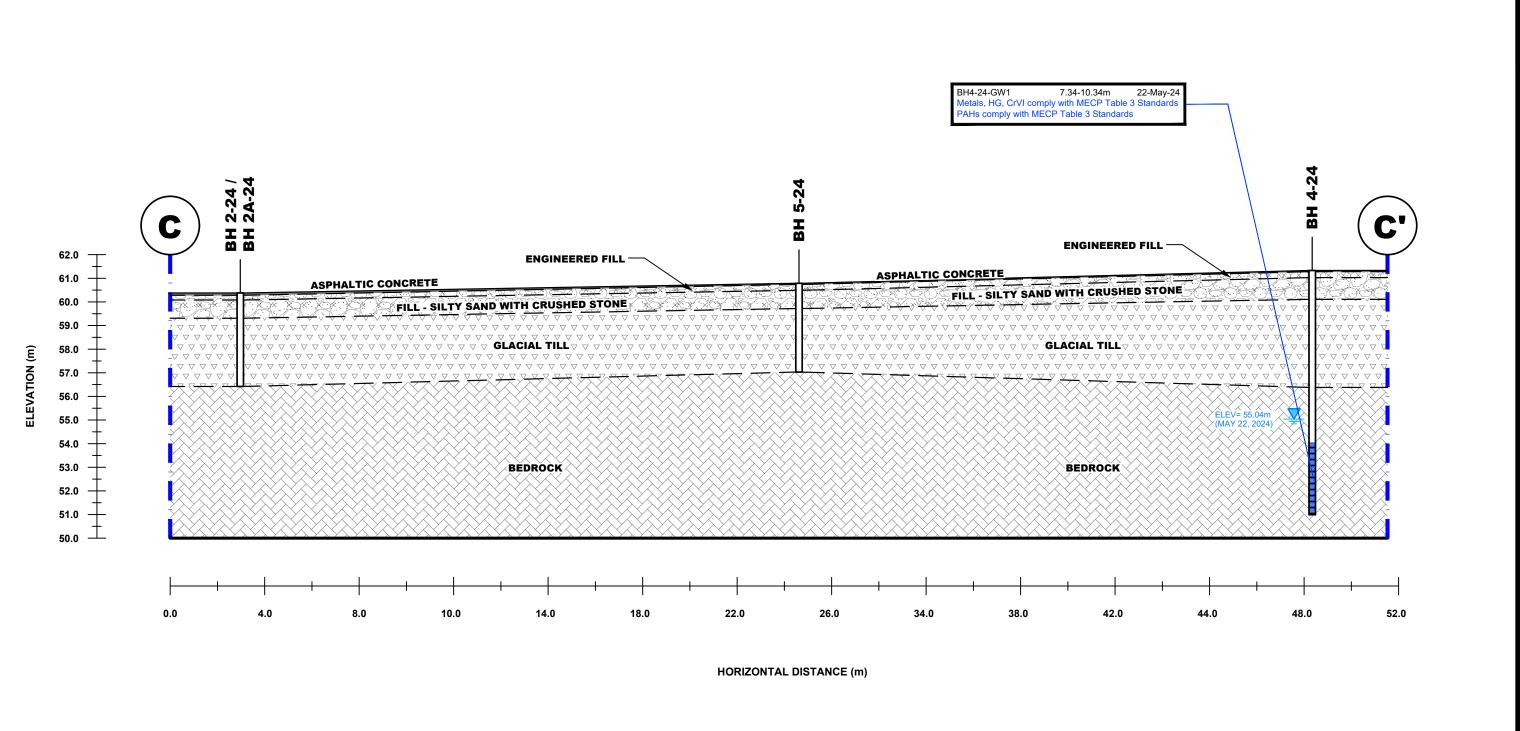
Approved by:

09/2024 **AS SHOWN** Drawn by: Report No.: PE2709-REP.04 ONTARIO Checked by: PE2709-18A

ANALYTICAL TESTING PLAN - GROUNDWATER (METALS, Hg, CrVI, & PAHs)

Revision No.:



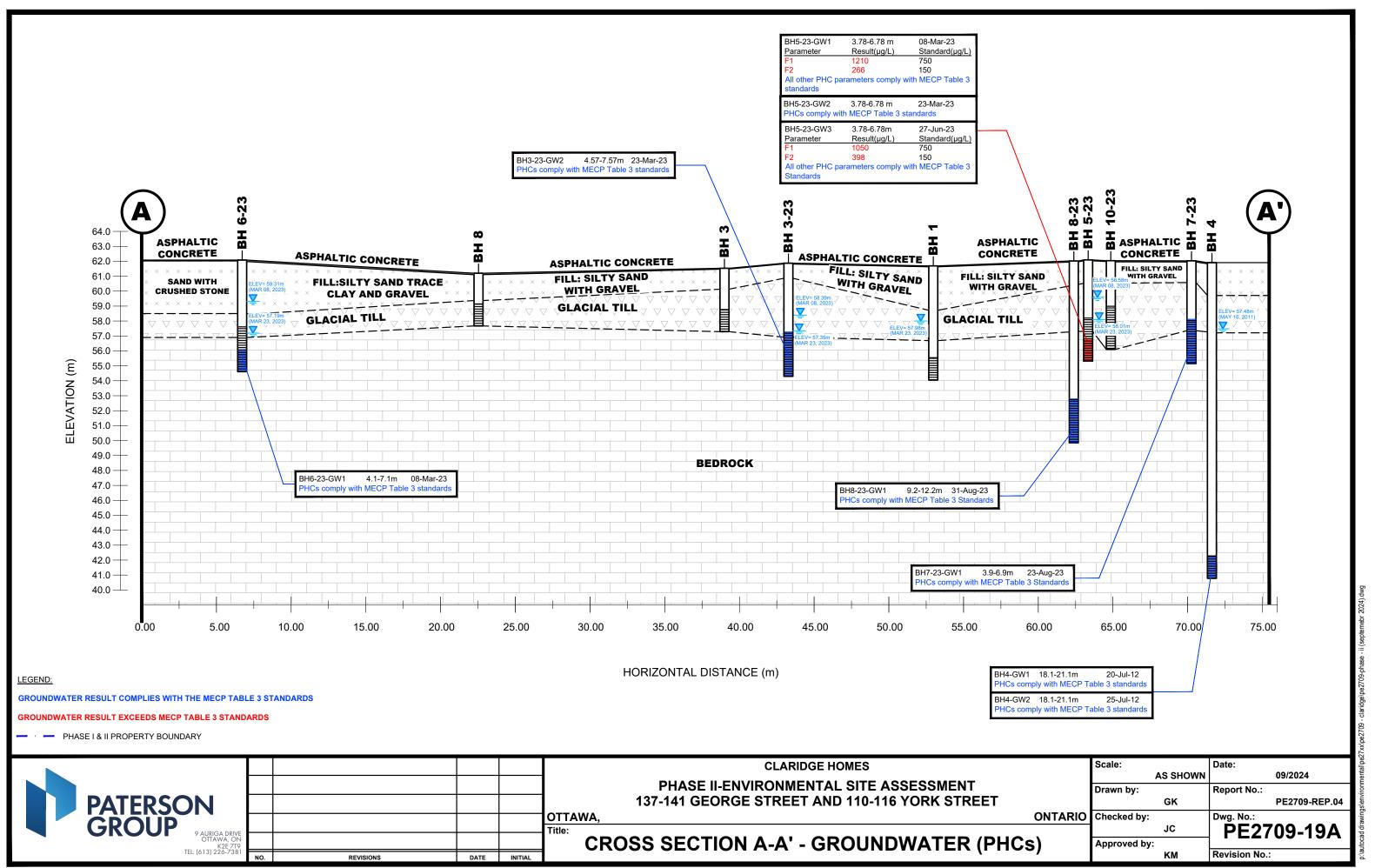


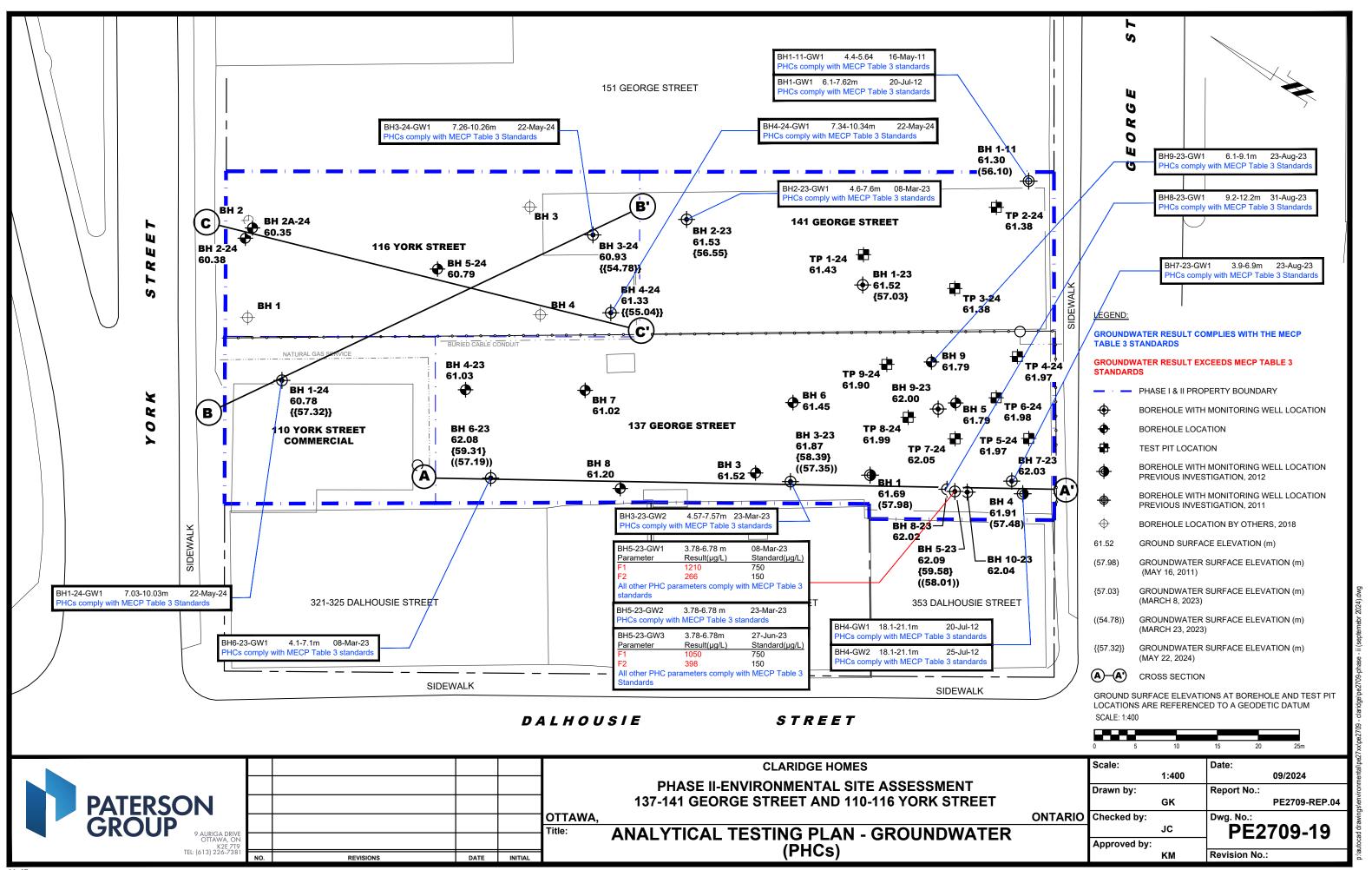
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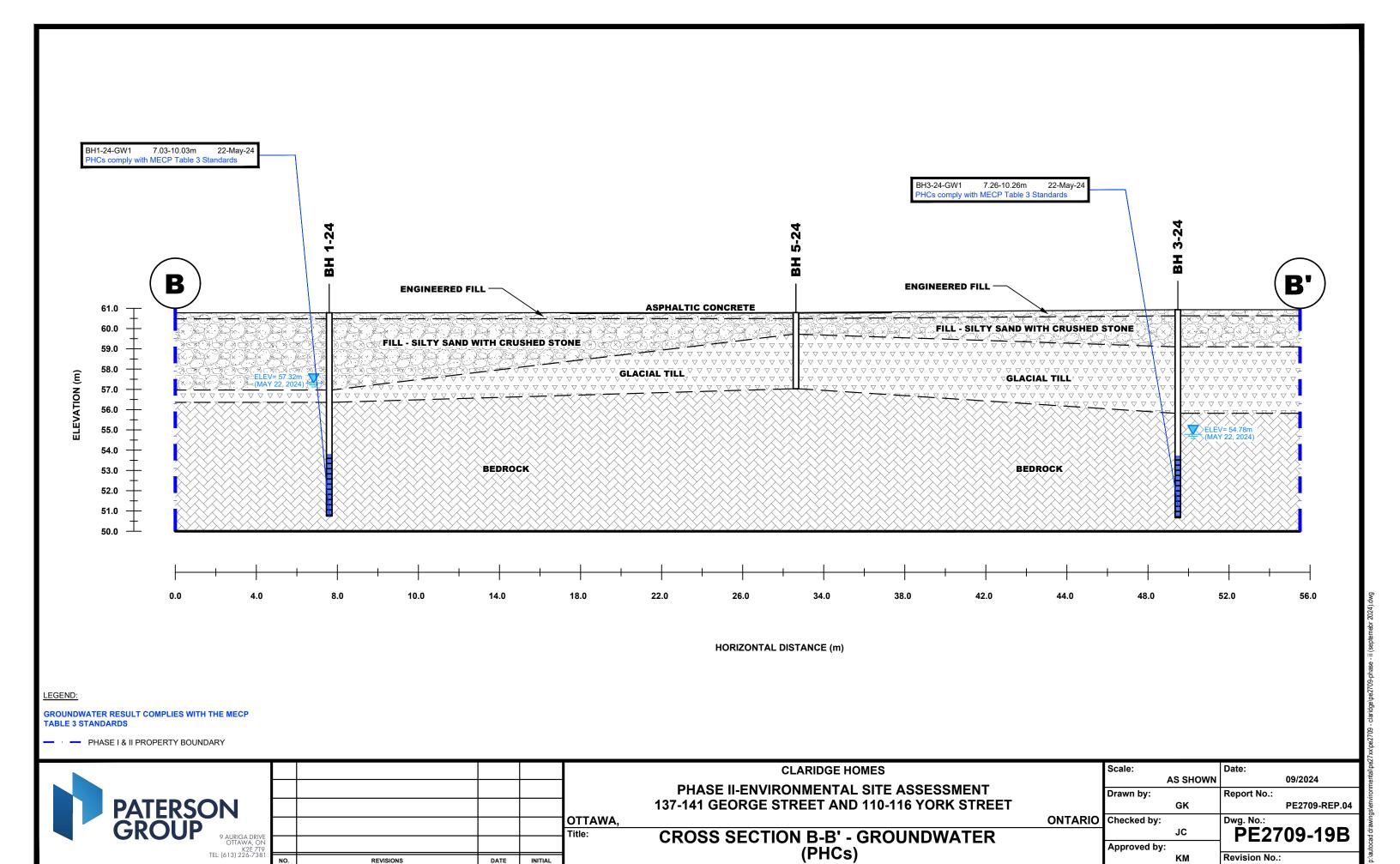
GROUNDWATER RESULT COMPLIES WITH THE MECP TABLE 3 STANDARDS

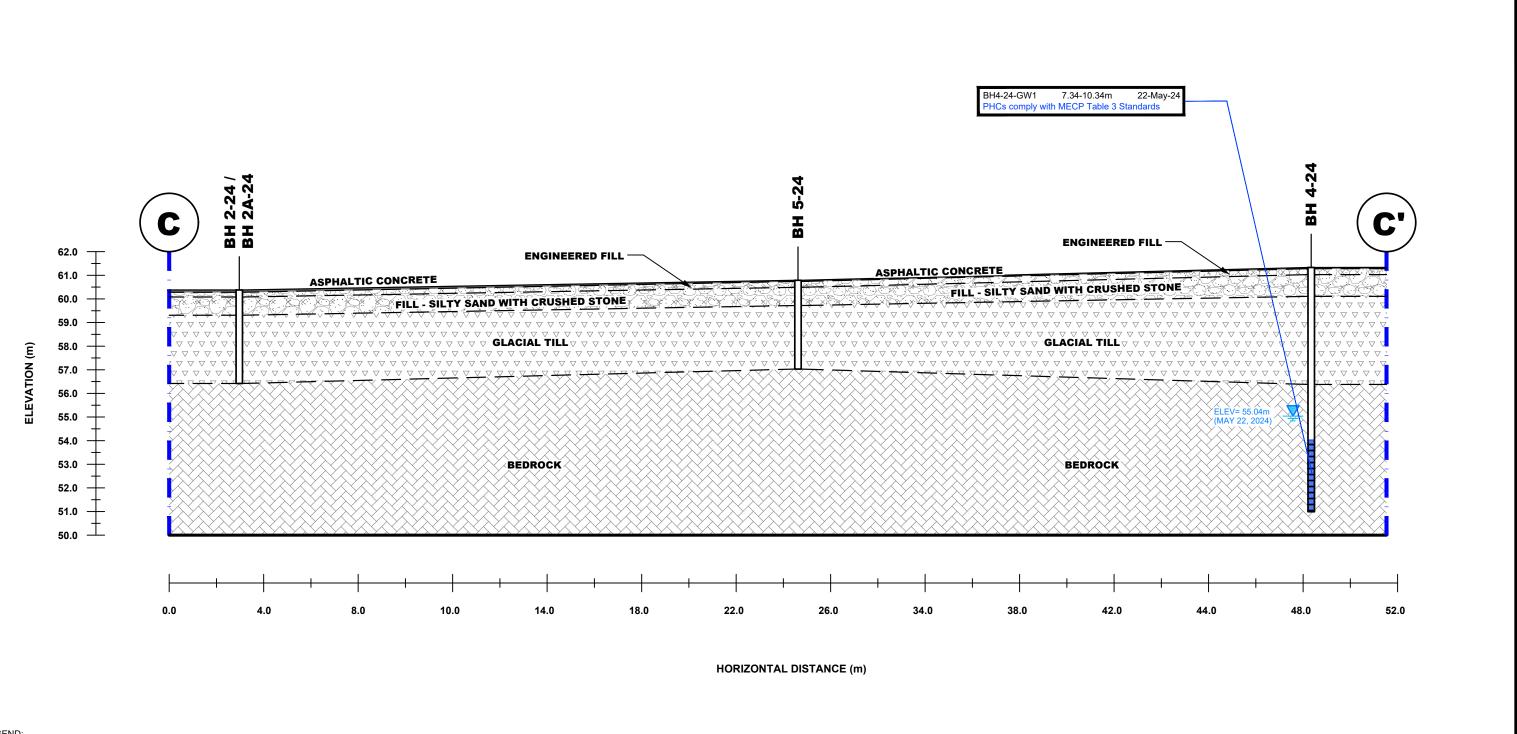
- PHASE I & II PROPERTY BOUNDARY

						CLARIDGE HOMES		Scale:	AS SHOWN	Date:	09/2024
PATERSON						PHASE II-ENVIRONMENTAL SITE ASSESSMENT 137-141 GEORGE STREET AND 110-116 YORK STREET		Drawn by:	GK	Report No.:	PE2709-REP.04
GROUP 9 AURIGA DRIVE					OTTAWA, Title:	CROSS SECTION C-C' - GROUNDWATER	ONTARIO	Checked by:	JC	Dwg. No.: PF27	09-18C
OTTAWA, ON K2E 7T9 TEL: (613) 226-7381	NO.	REVISIONS	DATE	INITIAL		(METALS, Hg, CrVI, & PAHs)		Approved by:		Revision No.:	









LEGEND:

GROUNDWATER RESULT COMPLIES WITH THE MECP TABLE 3 STANDARDS

PHASE I & II PROPERTY BOUNDARY

TEL: (613) 226-7381	NO.	REVISIONS	DATE	INITIAL	1
9 AURIGA DRIVE OTTAWA, ON K2E 7T9 TEL: (613) 226-7381					CROSS SEC
GROUP					OTTAWA, Title:
PATERSON					137-141 GI
					PHASE

CLARIDGE HOMES SE II-ENVIRONMENTAL SITE ASSESSMENT GEORGE STREET AND 110-116 YORK STREET

AS SHOWN 09/2024 Drawn by: Report No.: PE2709-REP.04

ONTARIO Checked by:

PE2709-19C Approved by:

Revision No.:

CTION C-C' - GROUNDWATER (PHCs)

TABLES

TABLE A1: SOIL ANALYTICAL RESULTS COMPARED TO MECP TABLE 3 STANDARDS RESIDENTIAL PROPERTY USE

TABLE A2: SOIL ANALYTICAL RESULTS COMPARED TO MECP TABLE 3 STANDARDS RESIDENTIAL PROPERTY USE



Parameter Sample Depth (m)	Units	MDL	Regulation Reg 153/04-Table 3 Residential,	BH1-AU1 1229217-01 0.00-0.25	BH1-SS7 1229139-01 3.80-4.20	BH3-SS6 1229139-03 3.00-3.60	BH4-SS3 1229217-02 0.80-1.20	BH5-SS8 1232102-01 5.20-5.41	BH6-SS7 1232102-02 4.60-4.75	BH7-SS7 1232102-03 4.60-4.70	BH1-23-SS2 2310245-12 0.76-1.37	BH1-23-SS3 2310245-01 1.60-2.20	BH1-23-SS4 2309081-02 2.29-2.90	BH1-23-SS5 2310245-13 3.20-3.40	BH2-23-SS2 2310245-02 0.80-1.40	BH2-23-SS4 2310245-03 2.40-3.00	BH2-23-SS5 2309472-01 3.05-3.66	BH3-23-AU2 2310245-04 0.25-0.61	BH3-23-SS4 2310245-05 1.60-2.20	BH3-23-SS5 2310245-14 3.20-3.60	DUP1-23 2309472-07 3.20-3.60	BH3-23-SS6 2310245-15 4.20-5.60	BH4-23-SS3 2309472-05 0.76-1.37	BH4-23-SS4 2310245-06 1.60-2.20	BH5-23-AU1 2310245-07 0.06-0.25	BH5-23-SS3 2310245-08 1.60-2.20	BH5-23-SS6 2310245-09 4.00-4.60	BH6-23-SS2 2309472-06 0.76-1.37	BH6-23-SS5 2310245-10 3.20-3.40
Sample Date		ı	coarse	17/Jul/2012	17/Jul/2012	17/Jul/2012	18/Jul/2012	8/Aug/2012	8/Aug/2012	8/Aug/2012	24/Feb/2023	24/Feb/2023	24/Feb/2023	24/Feb/2023	27/Feb/2023	27/Feb/2023	27/Feb/2023	27/Feb/2023	27/Feb/2023	27/Feb/2023	27/Feb/2023	27/Feb/2023	28/Feb/2023	28/Feb/2023	28/Feb/2023	28/Feb/2023	28/Feb/2023	28/Feb/2023	1/Mar/2023
Phsical Characteristics % Solids	% by Wt.	0.1		98.6	89.1	90.4	89.6	74.0	96.4	96.5	90.7	93.2	92.5	92.2	88.4	92.7	92.6	90.6	96.2	90.5	93.4	91.1	92.5	92.5	92.6	98.4	94.2	96.5	94.7
General Inorganics	N/A	0.01	5.0			_					1.9	0.8	_		1.61	1.77		2.16	1.19	1.54		2.07	4.39	4.44	0.9	1.32	1.25	4,65	1.88
Conductivity	uS/cm	5.00	700	-	-	-	-	-	-	-	943	469	- 1	-	1360	693	-	442	209	260	-	317	1160	1090	280	293	226	1770	3310
рН	pH Units	0.05	NV	-	-	-	-	-	-	-	-	-	-	-	7.69	7.76	-	7.66	-	-	-	7.88	7.59	-		7.89		7.89	-
Metals Antimony	ug/g dry	1.0	7.5	<1	-	-	2	-	-	-	ND (1.0)	ND (1.0)	-	-	2.2	ND (1.0)	-	2.2	ND (1.0)	ND (1.0)	-	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)
Arsenic	ug/g dry	1.0	18	2		-	2	-	-	-	3.5	3.8	-	-	8	2.6	-	6.8	4.1	2.9		2.5	3.4	3.5	5.9	2.8	4.1	2.2	10.2
Barium Beryllium	ug/g dry ug/g dry	1.0 0.5	390 4.0	20 <0.5	-	-	184 <0.5	-		-	62.5 ND (0.5)	30.7 ND (0.5)	-	-	321 0.6	27.1 ND (0.5)	•	126 0.6	38.6 ND (0.5)	87.3 ND (0.5)	-	68.1 ND (0.5)	31.3 ND (0.5)	27.1 ND (0.5)	24.1 ND (0.5)	38.1 ND (0.5)	129 ND (0.5)	211 ND (0.5)	64.7 ND (0.5)
Boron	ug/g dry	0.5	120	<5.0	-	-	<5.0	-		-	ND (5.0)	6.4	-	-	7	6.5		ND (5.0)	5.6	5.3	-	8.1	6.8	7.2	6.8	5.6	8	16.2	11.6
Chromium (VII)	ug/g dry	0.5 0.2	1.2	<0.5	-	-	0.5	-	-	-	ND (0.5) ND (0.2)	ND (0.5) ND (0.2)	-	-	0.6 ND (0.2)	ND (0.5) ND (0.2)	-	ND (0.5) ND (0.2)	ND (0.5) ND (0.2)	ND (0.5) ND (0.2)	-	ND (0.5) ND (0.2)	ND (0.5) ND (0.2)	ND (0.5) ND (0.2)	ND (0.5) ND (0.2)	ND (0.5)	ND (0.5) ND (0.2)	ND (0.5)	ND (0.5) ND (0.2)
Chromium (VI) Chromium	ug/g dry ug/g dry	5	8.0 160	7	-	-	14	-	-	-	14.9	12.2	-	-	29.2	10.1		16.7	11.3	9.9	-	10.8	13	12.5	15.9	ND (0.2) 10.5	13.5	ND (0.2) 11	13.6
Cobalt	ug/g dry	1	22	3	-	-	3	-		-	3.9	3.8	-	-	7.1	3.6		5.7	4.2	3.2	-	2.8	3.3	3.5	6.8	3.7	4.5	5.5	11.8
Copper Lead	ug/g dry ug/g dry	5	140 120	9	-	-	17 524	-	-	-	13 61.9	9.3		-	120 432	5.7 3.2		36.6 217	8.7 5.3	ND (5.0) 3.3	-	ND (5.0) 2.8	6.1 12.3	7 4.5	12.4 32.5	5.9 4.5	ND (5.0)	7.2 5.8	9.6 19.7
Mercury	ug/g dry	0.1	0.27	-	-	-	-	-	-	-	0.2	ND (0.1)	-	-	2.7	ND (0.1)	-	0.3	ND (0.1)	ND (0.1)	-	ND (0.1)	ND (0.1)	ND (0.1)	ND (0.1)	ND (0.1)	ND (0.1)	ND (0.1)	ND (0.1)
Molybdenum	ug/g dry	1	6.9 100	2 11	-	-	1 8	-	-	-	ND (1.0) 8.7	ND (1.0) 8.2	-	-	1.1	ND (1.0) 7.1	-	1.3 14.3	ND (1.0) 8.6	ND (1.0) 6.8	-	ND (1.0) 7.2	ND (1.0) 7.5	ND (1.0) 7.7	6.7 15.6	ND (1.0) 6.8	ND (1.0) 11.4	10.1	5.4 20.3
Selenium	ug/g dry ug/g dry	1	2.4	<1	-	-	<1	-	-	-	ND (1.0)	ND (1.0)	- 1	-	1.3	ND (1.0)	-	ND (1.0)	ND (1.0)	ND (1.0)	-	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)
Silver	ug/g dry	0.3	20	1.8	-	-	1.5	-		-	ND (0.3)	ND (0.3)	-	-	0.6	ND (0.3)		ND (0.3)	ND (0.3)	ND (0.3)	-	ND (0.3)	ND (0.3)	ND (0.3)	ND (0.3)	ND (0.3)	ND (0.3)	ND (0.3)	ND (0.3)
Thallium Uranium	ug/g dry ug/g dry	1	1.0	<1	-	-	<1	-	-	-	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)		-	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)		ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	-	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)	ND (1.0) ND (1.0)
Vanadium	ug/g dry	1	86	30	-	-	20	-	-	-	17.2	17.5	-	-	26.1	15.7	-	22.1	19.2	15.7	-	12.9	18.9	19.5	25.8	14.3	15.4	ND (10.0)	10.5
Zinc Methyl Mercury	ug/g dry	10 0.00005	340 0.0084	<20	-	-	191	-	-	-	55.3	21.8	-		429	ND (20.0)	-	74.8	20.5	ND (20.0)	-	ND (20.0)	32.1	23.8	26.3	ND (20.0)	ND (20.0)	ND (20.0)	ND (20.0)
Volatiles	ug/g dry	0.00005	0.0084	 	 	 		-	<u> </u>		<u> </u>			- +	-	-	-	-	- -	- -	- 1	-	-	-			+	+ -	+ -
Acetone	ug/g dry	0.5	16		ND (0.50)	ND (0.50)	-	-	-	-	-	-	ND (0.50)	-	-		ND (0.50)		ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	-		-		ND (0.50)	1	
Benzene Bromodichloromethane	ug/g dry	0.02	0.21 13	-	ND (0.02) ND (0.05)	ND (0.02) ND (0.05)	-	-	-	-		-	ND (0.02) ND (0.05)	- :	-		ND (0.02) ND (0.05)	-	ND (0.02) ND (0.05)	ND (0.02) ND (0.05)	ND (0.02) ND (0.05)	ND (0.02) ND (0.05)		-	- '	-	ND (0.02) ND (0.05)	+ :	1 -
Bromodichloromethane Bromoform	ug/g dry ug/g dry	0.05	0.27	i :	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-	-	ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	'	1	ND (0.05) ND (0.05)	+-:-	1 :
Bromomethane	ug/g dry	0.05	0.05		ND (0.05)	ND (0.05)	-	-	-	-	-		ND (0.05)	-	-		ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	-		ND (0.05)	<u> </u>	
Carbon Tetrachloride Chlorobenzene	ug/g dry	0.05 0.05	0.05	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-		-	ND (0.05) ND (0.05)	- :	-		ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)		-	- '	-	ND (0.05) ND (0.05)	+ :	1 -
Chlorobenzene Chloroform	ug/g dry ug/g dry	0.05	0.05	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-	-	ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	- '	<u> </u>	ND (0.05) ND (0.05)	 	1 -
Dibromochloromethane	ug/g dry	0.05	9.4	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-		-	ND (0.05)	<u> </u>	-
Dichlorodifluoromethane 1,2-Dichlorobenzene	ug/g dry	0.2	16 3.4	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-	-	ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	· ·	-	ND (0.05) ND (0.05)		-
1,3-Dichlorobenzene	ug/g dry ug/g dry	0.05	4.8	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-		-	ND (0.05)	-	-
1,4-Dichlorobenzene	ug/g dry	0.05	0.083	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	-	-	ND (0.05)	-	-
1,1-Dichloroethane	ug/g dry	0.05	3.5 0.05	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-	-	ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	· '	-	ND (0.05) ND (0.05)	-	-
1,2-Dichloroethane 1,1-Dichloroethylene	ug/g dry ug/g dry	0.05	0.05	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05) ND (0.05)	-	-		-	ND (0.05)	 	-
cis-1,2-Dichloroethylene	ug/g dry	0.05	3.4	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-		-	ND (0.05)	-	-
trans-1,2-Dichloroethylene	ug/g dry	0.05	0.084	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	· '	-	ND (0.05)	-	-
1,2-Dichloropropane cis-1,3-Dichloropropylene	ug/g dry ug/g dry	0.05	0.05 0.05	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-	-	ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-		-	ND (0.05) ND (0.05)	 	-
trans-1,3-Dichloropropylene	ug/g dry	0.05	0.05	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	- '	-	ND (0.05)	-	-
1,3-Dichloropropene, total	ug/g dry	0.05	0.05	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	· '	-	ND (0.05)	-	-
Ethylbenzene Ethylene dibromide (dibromoethane, 1,2-)	ug/g dry ug/g dry	0.05	2.0 0.05	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-		ND (0.05) ND (0.05)		ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	<u> </u>	-	0.1 ND (0.05)	-	-
Hexane	ug/g dry	0.05	2.8	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	-	-	ND (0.05)	-	-
Methyl Ethyl Ketone (2-Butanone) Methyl Isobutyl Ketone	ug/g dry	0.05	16 1.7	-	ND (0.50) ND (0.50)	ND (0.50) ND (0.50)	-	-	-	-	-	-	ND (0.50) ND (0.50)	-	-	-	ND (0.50) ND (0.50)	-	ND (0.50) ND (0.50)	ND (0.50) ND (0.50)	ND (0.50) ND (0.50)	ND (0.50) ND (0.50)	-	-	· · ·	-	ND (0.50) ND (0.50)		-
Methyl tert-butyl ether	ug/g dry ug/g dry	0.05	0.75	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-		-	ND (0.05)	-	-
Methylene Chloride	ug/g dry	0.5	0.1	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	- '	-	ND (0.05)	-	-
Styrene 1,1,1,2-Tetrachloroethane	ug/g dry ug/g dry	0.5	0.7 0.058	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-	-	ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	· ·	-	ND (0.05) ND (0.05)	 	-
1,1,2,2-Tetrachioroethane	ug/g dry	0.05	0.05	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-		-	ND (0.05)	+ -	
Tetrachloroethylene	ug/g dry	0.05	0.28	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	- '	-	ND (0.05)	-	-
Toluene 1,1,1-Trichloroethane	ug/g dry ug/g dry	0.05	2.3 0.38	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	-	-	-	-	ND (0.05) ND (0.05)	-	-		ND (0.05) ND (0.05)	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	-	· · · · · · · · · · · ·	-	ND (0.05) ND (0.05)		-
1,1,2-Trichloroethane	ug/g dry	0.05	0.05	-	ND (0.05)	ND (0.05)	-	-	-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	· - '	-	ND (0.05)	-	-
Trichloroethylene	ug/g dry	0.05	0.061		ND (0.05)	ND (0.05)	-	-	-	-	-		()															1	
Trichlorofluoromethane Vinyl Chloride	ug/g dry ug/g dry	0.05	4.0 0.02	-	ND (0.05)	ND (0.05)				1			ND (0.05)	-	-		ND (0.05)		ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	'	-	ND (0.05)		1 .
m/p-Xylene	ug/g dry	0.05			ND (0,02)	ND (0,02)			-	-	-	-	ND (0.05)	-	-	-	ND (0.05)	-	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	-	-	-	-	ND (0.05)	-	-
o-Xylene Vylene total			3.1	-	ND (0.02) ND (0.05)	ND (0.02) ND (0.05)	-		-	-		-	ND (0.05) ND (0.02) ND (0.05)	-	- - -	- - -	ND (0.05) ND (0.02) ND (0.05)	-	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	-	- - -	- - -		ND (0.05) ND (0.02) 0.22		-
Xylenes, total BTEX	ug/g dry	0.05	3.1 3.1	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	- - -		-	-	- - - -	- - -	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	- - - -	- - - -	- - - -	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	- - - -	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)		- - - -	- - - -		ND (0.05) ND (0.02) 0.22 ND (0.05)		-
Benzene			3.1	-	ND (0.05)	ND (0.05)	-					-	ND (0.05) ND (0.02) ND (0.05)	- - - - -	- - - -	- - - - -	ND (0.05) ND (0.02) ND (0.05)	- - - - -	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)					ND (0.05) ND (0.02) 0.22	-	-
Ethylbenzene	ug/g dry ug/g dry ug/g dry	0.05 0.05 0.02	3.1 3.1 3.1 0.21	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	- - - - ND (0.02)	- - - - - ND (0.02)	- - - - - ND (0.02)	- - - - - ND (0.02)	- - - - - ND (0.02)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	- - - - - ND (0.02)		- - - - - ND (0.02)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	- - - - - ND (0.02)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	- - - - - ND (0.02)	- - - - - - ND (0.02)	- - - - - - ND (0.02)	- - - - - - ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05)		- - - - ND (0.02)
	ug/g dry ug/g dry ug/g dry ug/g dry	0.05 0.05 0.02 0.05	3.1 3.1 3.1 0.21 2.0	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)		ND (0.02)	ND (0.02)	ND (0.02)	ND (0.05)	ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05)	ND (0.05) ND (0.02) 0.22 ND (0.05)	ND (0.05)	ND (0.05)
Toluene m/p-Xylene	ug/g dry ug/g dry ug/g dry	0.05 0.05 0.02	3.1 3.1 3.1 0.21	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	· · · · · · · · · · · · · · · · · · ·						ND (0.05) ND (0.02) ND (0.05) ND (0.05)				ND (0.05) ND (0.02) ND (0.05) ND (0.05)		ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)					ND (0.05) ND (0.02) 0.22 ND (0.05)		
Toluene m/p-Xylene o-Xylene	ug/g dry	0.05 0.05 0.02 0.05 0.05 0.05 0.05	3.1 3.1 0.21 2.0 2.3 3.1 3.1	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) 0.22 ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)
Toluene m/p-Xylene o-Xylene Xylenes, total	ug/g dry	0.05 0.05 0.02 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1		ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) 0.22 ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)
Toluene m/p-Xylene o-Xylene	ug/g dry	0.05 0.05 0.02 0.05 0.05 0.05 0.05	3.1 3.1 0.21 2.0 2.3 3.1 3.1		ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) 0.22 ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)
Toluene m/p-Xylene o-Xylene Xylenes, total Hydrocorbons F1 PHCS (C5-C10) F2 PHCS (C10-C16)	ug/g dry	0.05 0.05 0.02 0.05 0.05 0.05 0.05 0.05	3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 55		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4	ND (0.02) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.7) ND (77)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) 	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (80)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.02) 0.22 ND (0.05) 0.22 ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)
Toluene m/p-Xylene o-Xylene o-Xylene Vylene, total Hydrocarubons F1 PMCS (CG-CL0) F2 PMCS (CL0-CL6) F3 PMCS (CL0-CL6) F3 PMCS (CL0-CL4)	ug/g dry	0.05 0.05 0.02 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (40) 312	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 48	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (80) 699	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 50	ND (0.05) ND (7) ND (4) 15
Toluene m/p-Xylene o-Xylene Xylenes, total Hydrocorbons F1 PHCS (C5-C10) F2 PHCS (C10-C16)	ug/g dry	0.05 0.05 0.02 0.05 0.05 0.05 0.05 0.05	3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 55		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4	ND (0.02) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.7) ND (77)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05) 	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (80)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.02) ND (0.02) 0.22 ND (0.05) 0.22 ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4)
Toluene m/p-Xylene o-Xylene Xylenes, total Hydrocarchons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C10-C16) F4 PHCS (C34-C30) F4 PHCS (C34-C30) F4 PHCS (C34-C30) Semi-Volatilities	ug/g dry	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 55 98 300 2800 2800	-	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107 70	ND (0.05) ND (7) ND (4) ND (8) ND (6)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (7) ND (40) 312 298 -	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (4) 48 23	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (7) ND (4) ND (8) ND (6) -	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (7) ND (80) 699 1650	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (7) ND (4) 50 42	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.10) ND (10.05) ND (110 ND (1
Toluene m/p-Xylene o-Xylene Xylenes, total Hydrocarbons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C10-C34) F4 PHCS (C34-C50) F4 PHCS (C44-C50) F4 PHCS (C	ug/g dry	0.05 0.05 0.02 0.05 0.05 0.05 0.05 0.05	3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.5 98 300 2800 2800	- - - - - - - - - - - - -	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107 70 - 0.03	ND (0.05) ND (7) ND (4) ND (8) ND (6) - ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05)	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 48 23 - 0.07	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (7) ND (4) ND (8) ND (6) - ND (0.02)	ND (0.05) ND (7) ND (4) ND (8) ND (6) ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (80) 699 1650 - ND (0.40)	ND (0.05) ND (7) ND (4) ND (8) ND (6) - ND (6)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) S0 42 - ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 15 11 - ND (0.02)
Toluene m/p-Xylene o-Xylene Xylenes, total Hydrocarchons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C10-C16) F4 PHCS (C34-C30) F4 PHCS (C34-C30) F4 PHCS (C34-C30) Semi-Volatilities	ug/g dry	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 55 98 300 2800 2800	-	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107 70	ND (0.05) ND (7) ND (4) ND (8) ND (6)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (7) ND (40) 312 298 -	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (4) 48 23	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (7) ND (4) ND (8) ND (6) -	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (7) ND (80) 699 1650	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (7) ND (4) 50 42	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.10) ND (10.05) ND (110 ND (1
Toluene m/p-Xylene o-Xylene vylenes, total Mydrocarbons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C10-C34) F4 PHCS (C34-C30) F4 PHCS (C34-C30) F4 PHCS (C34-C30) F4 PHCS (C34-C30) F4 PHCS (C10-C34) F4 PHCS (C34-C30) F4 PHCS (C3	ug/g dry	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 7.5 98 300 2800 2800 2000		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107 70 - 0.03 0.05	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8) ND (6) ND (0.02) ND (0.02) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (7) ND (40) 312 298 - 0.76 0.19 1.58 2.72	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (77) ND (4) 48 23 0.07 0.18 0.29 0.7	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8) ND (6) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (80) 699 1650 - ND (0.40) ND (0.40) ND (0.40) ND (0.40) ND (0.40)	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (7) ND (4) 50 42 ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.05) 15 11 - ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02)
Toluene m/p-Xylene o-Xylene o-Xylene Xylenes, total Hydrocarrbons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C10-C16) F4 PHCS (C34-C30) F	ug/g dry	0.05 0.05 0.02 0.02 0.05 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 55 98 300 2800 2800 7.9 0.15 0.67 0.5 0.3		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107 70 0.03 0.05 0.09 0.23 0.26	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (40) 312 298 - - - 0.76 0.19 1.58 2.72 2.37	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (2) ND (2) ND (4) 48 23 0.07 0.18 0.29 0.7 0.7	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02)	ND (0.05)	ND (0.05) ND (7) ND (4) ND (8) ND (6) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 50 42 - ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 15 11 - ND (0.02)
Toluene m/p-Xylene 0-Xylene Xylene, Stotal Hydrocarbons F1 PHCS (C6-C10) F2 PHCS (C16-C3) F3 PHCS (C16-C3) F3 PHCS (C16-C3) F4G PHCS (C34-C5) F4G PHCS (C34-	ሀል/ይ dry	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.5 98 300 2800 2800 2800 7.9 0.15 0.67 0.5 0.3 0.78		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107 70 - 0.03 0.05 0.09 0.23 0.26	ND (0.05) ND (7) ND (7) ND (8) ND (6) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (7) ND (40) 312 298 - 0.76 0.19 1.58 2.72	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (7) ND (4) 48 23	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.02) ND (0.02) ND (0.02) ND (0.02)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (7) ND (4) ND (8) ND (6) ND (6) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (80) 699 1650 - ND (0.40) ND (0.40) ND (0.40) ND (0.40) ND (0.40)	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (7) ND (4) 50 42 - ND (0.02)	ND (0.05) 15 11
Toluene m/p-Xylene o-Xylene o-Xylene Xylenes, total Hydrocarrbons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C10-C16) F4 PHCS (C34-C30) F	ug/g dry	0.05 0.05 0.02 0.02 0.05 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 55 98 300 2800 2800 7.9 0.15 0.67 0.5 0.3		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	-	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) 17 107 70 0.03 0.05 0.09 0.23 0.26	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (40) 312 298 0.76 0.19 1.58 2.72 2.37	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (2) ND (2) ND (4) 48 23 0.07 0.18 0.29 0.7 0.7	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) ND (8) ND (6) ND (0.02)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (7) ND (4) ND (8) ND (6)	ND (0.05) ND (0.02)	ND (0.05) ND (0.40)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 50 42 - ND (0.02)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (4) 15 11 - ND (0.02)
Toluene m/p-Xylene o-Xylene Xylenes, total Hydrocarbons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C16-C34) F4 PHCS (C34-C50) F4 PHCS (C3	내용 여가 내용 이가	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 55 98 300 2800 2800 2800 0 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0	<0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) The control of the control	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (0.06) ND (7) ND (40) 312 298	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (7) ND (4) 48 23	ND (0.05) ND (7) ND (4) ND (8) ND (6) ND (0.02)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (8) ND (6) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (4) 50 42 ND (0.02)	ND (0.05) ND (0.02)
Toluene m/p-Xylene o-Xylene o-Xylene o-Xylene o-Xylene xylenes, total Hydrocarbons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C10-C16) F4 PHCS (C34-C50) F5 PHCS (C34-C50) F6 PHCS (F10-C30) F6	ሀይ/ይ dry	0.05 0.05 0.02 0.02 0.05 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 3.1 0.25 98 98 300 2800 2800 2800 0 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0 0.1	 - -	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) 17 107 70	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (40) 312 298 0.76 0.19 1.58 2.72 2.37 3 1.45 1.68 2.68 0.38	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.10) ND (4) 48 23	ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05) ND (0.40)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)
Toluene m/p-Xylene o-Xylene Xylenes, total Hydrocarbons F1 PHCS (C6-C10) F2 PHCS (C10-C16) F3 PHCS (C16-C34) F4 PHCS (C34-C50) F4 PHCS (C3	내용 여가 내용을 이가 내용을 이기를	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 55 98 300 2800 2800 2800 0 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0	<0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08 <0.08	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) The control of the control	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (0.06) ND (7) ND (40) 312 298	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (7) ND (4) 48 23	ND (0.05) ND (7) ND (4) ND (8) ND (6) ND (0.02)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (8) ND (6) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (4) 50 42 ND (0.02)	ND (0.05) ND (0.02)
Toluene m/p-Xylene O-Xylene Xylene, Stotal Mydrocarbons F1 PMcS (C6-C10) F2 PMcS (C16-C34) F3 PMcS (C16-C34) F3 PMcS (C16-C34) F46 PMCS (C34-C50)	내용 여가 내용 여기	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 3.1 3.1 55 98 300 2800 2800 2800 7.9 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0 0.1 0.69 62 0.38	 	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) 17 107 70	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (7) ND (40) 312 298 0.76 0.19 1.58 2.72 2.37 3 1.45 1.68 2.68 0.38 6.09 0.53 1.4	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (7) ND (7) ND (4) 48 23 0.07 0.18 0.29 0.7 0.73 0.38 0.42 0.71 0.11 1.47 0.08	ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05) ND (0.40)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (7) ND (4) 50 42 ND (0.02)	ND (0.05) ND (0.02)
Toluene m/p Xylene o Xylene o Xylene o Xylene o Xylene tylenes, total Mydrocarbons F1 PHCS (C16-C10) F2 PHCS (C10-C16) F3 PHCS (C16-C34) F4 PHCS (C16-C34) F	내를 여가 내를 여기 내를 여기 내를 이가 내를 이기 때를 이기	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 3.1 55 98 300 2800 2800 2800 7.9 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0 0.1 0.69 62 0.38 0.99		ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) 17 107 70	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (7) ND (40) 312 298	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (4) 48 23	ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (8) ND (6) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05) ND (0.40)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)
Toluene m/p-Xylene O-Xylene Xylene, Stotal Mydrocarbons F1 PMcS (C6-C10) F2 PMcS (C16-C34) F3 PMcS (C16-C34) F3 PMcS (C16-C34) F46 PMCS (C34-C50)	내용 여가	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 3.1 3.1 55 98 300 2800 2800 2800 7.9 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0 0.1 0.69 62 0.38	 	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) 17 107 70	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (7) ND (40) 312 298 0.76 0.19 1.58 2.72 2.37 3 1.45 1.68 2.68 0.38 6.09 0.53 1.4	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (7) ND (7) ND (4) 48 23 0.07 0.18 0.29 0.7 0.73 0.38 0.42 0.71 0.11 1.47 0.08	ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05) ND (0.40)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (7) ND (4) 50 42 ND (0.02)	ND (0.05) ND (0.02)
Toluene m/p-Xylene o-Xylene o-Xylene o-Xylene o-Xylene xylenes, total Hydrocarubons F3 PHCS (C5-C10) F3 PHCS (C16-C14) F3 PHCS (C16-C34) F4 PHCS (C16-C34) F	ug/g dry	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 3.1 55 98 300 2800 2800 2800 2800 0 7.9 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0 0.1 0.69 62 0.38 0.99 0.99	 - -	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) 17 107 70	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (10.05) N	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (10.05)	ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05) ND (0.40)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)
Toluene m/p-Xylene o-Xylene o-Xylene o-Xylene o-Xylene Xylenes, total Hydrocarbons F1 PHCS (C5C-10) F2 PHCS (C16-C3) F3 PHCS (C16-C3) F4 PHCS (C36-C3) F4 PHCS (C34-C5) F4 PHCS (C34-C5) F4 PHCS (C34-C5) F4 PHCS (C34-C5) F4 PHCS (F34-C3) F4 PHCS	내용 여가 내용 이가	0.05 0.05 0.05 0.05 0.05 0.05 0.05 0.05	3.1 3.1 3.1 3.1 0.21 2.0 2.3 3.1 3.1 3.1 3.1 3.1 3.1 55 98 300 2800 2800 7.9 0.15 0.67 0.5 0.3 0.78 6.6 0.78 7.0 0.1 0.69 62 0.38 0.99 0.99	 	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)		ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 65 74 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) 8 <4 <8	ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (0.02) ND (7) <4 62	ND (0.05) 17 107 70	ND (0.05) ND (0.02)	NO (0.05) ND (0.02) ND (0.02) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (7) ND (7) ND (4) ND (8)	ND (0.05) ND (0.	ND (0.05) ND (7) ND (4) ND (8)	ND (0.05) ND (0.02) ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.	ND (0.05)	ND (0.05) ND (0.02) ND (0.05)	ND (0.05) ND (0.02) ND (0.05) ND (0.05)	ND (0.05) ND (0.02) ND (0.02) ND (0.05)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)	ND (0.05) ND (0.40)	ND (0.05) ND (0.02)	ND (0.05) ND (0.02) 0.22 ND (0.05) 0.22	ND (0.05) ND (0.02)	ND (0.05) ND (0.02)

2.00 Result exceeds Reg 153/04-Table 3 Residential, coarse Standards
ND (0.2) MDL exceeds Reg 153/04-Table 3 Residential, coarse Standards
ND (0.2) No concentrations identified above the MDL
Parameter not analysed
NV No value given for indicated parameter



Parameter Sample Depth (n	Units n)	MDL	Regulation Reg 153/04-Table 3 Residential,	DUP 2310245-11 4.00-4.60	BH7-23-SS3 2333190-01 1.52-2.13	BH8-23-SS6 2333190-02 3.81-4.42	BH1-24-AU1 2420225-01 0.00-0.30	BH1-24-SS3 2420225-02 1.52-2.13	BH1-24-SS5 2420225-03 3.05-3.66	BH2-24-AU1 2420225-04 0.10-0.30	BH2-24-SS3 2420225-05 1.52-2.13	BH3-24-AU1 2420225-06 0.00-0.46	BH3-24-SS5 2420225-07 3.05-3.66	BH4-24-AU1 2420225-08 0.00-0.46	BH4-24-SS2 2420225-09 0.76-1.37	BH4-24-SS3 2420225-10 1.52 - 2.13	BH4-24-SS6 2420225-11 3.81-4.42	BH5-24-AU1 2420225-12 0.05-0.46	BH5-24-SS2(BOTTOM) 2420225-13 0.76-1.37	DUP1 2420225-14 3.05-3.66	DUP2 2420225-15 3.81-4.42
Sample Date	,		coarse	28/Feb/2023	14/Aug/2023	14/Aug/2023	8/May/2024	8/May/2024	8/May/2024	8/May/2024	8/May/2024	9/May/2024	9/May/2024	9/May/2024	9/May/2024	9/May/2024	9/May/2024	9/May/2024	9/May/2024	8/May/2024	9/May/2024
Phsical Characteristics % Solids	% by Wt.	0.1		96.0	91.8	93.6	91.5	91.5	90.4	96.4	91.8	91.6	92.1	93.6	80.8	93.6	94.5	93.4	92.9	92	93.7
General Inorganics	70 by ** t.	0.1		30.0	31.0	33.0	31.3	31.3	30.4	30.4	31.0	31.0	32.1	33.0	00.0	33.0	34.3	33.4	32.3		33.7
SAR	N/A uS/cm	0.01 5.00	5.0 700	1.43 304	-	-	2.19 2330	1.6 583	1.87 402	4.11 544	6.83 944	2.29 406	2.08 459	0.28 156	2.3 272		2.21 357	2 363	1.54 353	2.18 433	1.75 347
Conductivity pH	pH Units	0.05	NV	-	-	-	N/A	N/A	8.79	7.28	N/A	406 N/A	7.79	N/A	N/A	-	N/A	7.4	N/A	9.6	N/A
Metals																					
Antimony Arsenic	ug/g dry ug/g dry	1.0	7.5 18	ND (1.0) 4.6	-	-	<1.0 2.6	<1.0 2	<1.0 2.1	<1.0 2.7	<1.0 2.3	<1.0 2.8	<1.0 2.9	<1.0 4	<1.0 5.2	<1.0 4.8	<1.0	1.7 4.2	<1.0 2.1	<1.0 2.2	<1.0 2.7
Barium	ug/g dry	1.0	390	211	-	-	99.5	29.9	35.8	87.8	41.1	59.5	21.3	98.2	177	36.2	36.7	137	30.2	35.9	43
Beryllium Boron	ug/g dry	0.5	4.0 120	0.5 8.9	-	-	<0.5	<0.5 6.2	<0.5 <5.0	<0.5	<0.5 <5.0	<0.5 6.5	<0.5 7.4	<0.5	<0.5 5.3	<0.5 6.7	<0.5 7.4	<0.5 5.2	<0.5 <5.0	<0.5 5.2	<0.5 7.7
Cadmium	ug/g dry ug/g dry	0.5	1.2	ND (0.5)	-	-	11 <0.5	<0.5	<0.5	11.1 <0.5	<0.5	<0.5	<0.5	7.3 <0.5	<0.5	<0.5	<0.5	0.5	<0.5	<0.5	<0.5
Chromium (VI)	ug/g dry	0.2	8.0	ND (0.2)	-	-	0.8	0.3	<0.2	<0.2	<0.2	0.2		<0.2	<0.2	N/A	-	0.4	-	<0.2	
Chromium Cobalt	ug/g dry ug/g dry	1	160 22	15.5 4.9	-	-	20.5 5.7	10.5 4.9	11.8 4.6	14.8 5.4	11.7 4.6	13 4.7	11.7 3.6	10.4 5.1	18.5 4.6	13.9 6.4	9.5 2.9	14.2 4.8	10.1 3.6	12.2 4.8	10.5
Copper	ug/g dry	5	140	5.5	-	-	16.1	9.9	6.8	12.8	8.3	12.4	5.5	11.8	22.7	10.2	<5.0	53.8	6.6	7.8	<5.0
Lead	ug/g dry	1	120 0.27	5.1	-	-	32.9	10.5	4 <0.1	32.1 <0.1	3.7	43.4	5.2	50.7	349 0.9	8.5	5.6	0.2	3.7	4.2 <0.1	5.6
Mercury Molybdenum	ug/g dry ug/g dry	0.1	6.9	ND (0.1) ND (1.0)	-	-	<0.1 1.5	<0.1 <1.0	<1.0	<1.0	<0.1 <1.0	0.1 <1.0	<1.0	<0.1	<1.0	<0.1 1.1	<1.0	1.1	<1.0	<1.0	<1.0
Nickel	ug/g dry	5	100	12.5	-	-	13.2	10.8	8.3	11.5	8.4	9.5	8.3	11.3	10.2	10.4	6.5	14.9	6.9	8.7	7.1
Selenium Silver	ug/g dry ug/g dry	0.3	2.4	ND (1.0) ND (0.3)	-	-	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3	<1.0 <0.3
Thallium	ug/g dry	1	1.0	ND (1.0)	-	-	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Uranium	ug/g dry	1	23	ND (1.0)	-	-	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0
Vanadium Zinc	ug/g dry ug/g dry	1 10	86 340	18 ND (20.0)	-	-	27.5 35.1	14.9 <20.0	18.7 <20.0	25.1 34.7	20.6 <20.0	19.8 45.1	16.7 <20.0	27.6 26.6	21.1 94.5	21.1 75.7	13.6	28.8 228	16.2 <20.0	19.1	14.9 <20.0
Methyl Mercury	ug/g dry	0.00005	0.0084	- (20.0)	-	-	-	-		-	-	5.1		-	-	ND(0.00005)	-		-	-	
Volatiles	no to a			No (0.5-)																	
Acetone Benzene	ug/g dry ug/g dry	0.5 0.02	16 0.21	ND (0.50) ND (0.02)		-	- :			-	-	-		- :				-	-		+ :
Bromodichloromethane	ug/g dry	0.05	13	ND (0.05)	-		-		-	-	-	-	-				-	-	-	-	
Bromoform Bromomethano	ug/g dry	0.05	0.27	ND (0.05)	-	-	-	-	-	-	-	-	-		-	-	-	-	-		-
Bromomethane Carbon Tetrachloride	ug/g dry ug/g dry	0.05 0.05	0.05 0.05	ND (0.05) ND (0.05)	-					-	-	-					-	-	-		
Chlorobenzene	ug/g dry	0.05	2.4	ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	
Chloroform Dibromochloromethane	ug/g dry	0.05 0.05	0.05 9.4	ND (0.05) ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-		
Dichlorodifluoromethane	ug/g dry ug/g dry	0.03	16	ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
1,2-Dichlorobenzene	ug/g dry	0.05	3.4	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	- 1	-
1,3-Dichlorobenzene 1,4-Dichlorobenzene	ug/g dry ug/g dry	0.05 0.05	4.8 0.083	ND (0.05) ND (0.05)	-	-	-	-	-	- :	-	-	-	-	-	-	-	-	-	-	
1,1-Dichloroethane	ug/g dry	0.05	3.5	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	-	-
1,2-Dichloroethane	ug/g dry	0.05	0.05	ND (0.05)	-	-		-	-		-	-	-		-		-	-	-	-	-
1,1-Dichloroethylene cis-1,2-Dichloroethylene	ug/g dry ug/g dry	0.05	0.05 3.4	ND (0.05) ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-		
trans-1,2-Dichloroethylene	ug/g dry	0.05	0.084	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	-	-
1,2-Dichloropropane	ug/g dry	0.05	0.05	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-		-
cis-1,3-Dichloropropylene trans-1,3-Dichloropropylene	ug/g dry ug/g dry	0.05	0.05	ND (0.05) ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-		
1,3-Dichloropropene, total	ug/g dry	0.05	0.05	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	-	-
Ethylbenzene Ethylene dibromide (dibromoethane, 1,2-)	ug/g dry	0.05 0.05	2.0 0.05	0.08 ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-		
Hexane	ug/g dry ug/g dry	0.05	2.8	ND (0.05)	-	-	-	-	-		-	-	-		-		-	-	-	-	
Methyl Ethyl Ketone (2-Butanone)	ug/g dry	0.05	16	ND (0.50)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	- 1	-
Methyl Isobutyl Ketone Methyl tert-butyl ether	ug/g dry ug/g dry	0.05 0.05	1.7 0.75	ND (0.50) ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-		
Methylene Chloride	ug/g dry	0.5	0.1	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	-	-
Styrene	ug/g dry	2	0.7	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-		-
1,1,2,2-Tetrachloroethane 1,1,2,2-Tetrachloroethane	ug/g dry ug/g dry	0.5 0.05	0.058	ND (0.05) ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-		
Tetrachloroethylene	ug/g dry	0.05	0.28	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	-	-
Toluene 1,1,1-Trichloroethane	ug/g dry	0.05 0.05	2.3 0.38	ND (0.05) ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-		
1,1,2-Trichloroethane	ug/g dry ug/g dry	0.05	0.05	ND (0.05)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Trichloroethylene	ug/g dry	0.05	0.061	ND (0.05)	-	-	-	-	-		-	-	-	-	-	-	-	-	-	-	-
Trichlorofluoromethane Vinyl Chloride	ug/g dry ug/g dry	0.05	4.0 0.02	ND (0.05) ND (0.02)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-		
m/p-Xylene	ug/g dry	0.05	3.1	0.21	-	-	-		-	-	-	-	-	-		-	-	-	-	-	
o-Xylene Yylenes total	ug/g dry	0.05	3.1	ND (0.05)	-	-	-	-	-	-	-	-	-		-	-	-	-	-		-
Xylenes, total BTEX	ug/g dry	0.05	3.1	0.21	-	-	-	 		-	-	-	-	-	-	-		-	-		
Benzene	ug/g dry	0.02	0.21	-	ND (0.02)	ND (0.02)	<0.02	<0.02	<0.02	<0.02	<0.02	<0.02	<0.02	<0.02	<0.02		<0.02	<0.02	<0.02	<0.02	<0.02
Ethylbenzene Toluene	ug/g dry ug/g dry	0.05 0.05	2.0	-	ND (0.05) ND (0.05)	0.28 0.48	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	-	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05	<0.05 <0.05
m/p-Xylene	ug/g dry	0.05	3.1	-	ND (0.05)	0.46	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05	<0.05		<0.05	<0.05	<0.05	<0.05	<0.05
o-Xylene	ug/g dry	0.05	3.1	-	ND (0.05)	ND (0.05)	<0.05	<0.05	<0.05	<0.05	<0.05	0.07	<0.05	<0.05	<0.05	-	<0.05	<0.05	<0.05	<0.05	<0.05
Xylenes, total Hydrocarbons	ug/g dry	0.05	3.1	-	ND (0.05)	0.36	<0.05	<0.05	<0.05	<0.05	<0.05	0.07	<0.05	<0.05	<0.05	-	<0.05	<0.05	<0.05	<0.05	<0.05
F1 PHCs (C6-C10)	ug/g dry	7	55	8	ND (7)	355	<7	<7	<7	<7	<7	<7	<7	<7	<7		<7	<7	<7	<7	<7
F2 PHCs (C10-C16)	ug/g dry	4	98	28	ND (4)	377	<4	<4	<4	<40	39	<4	<4	<80	6	-	9	<40	<4	<4	<4
F3 PHCs (C16-C34) F4 PHCs (C34-C50)	ug/g dry ug/g dry	8 6	300 2800	ND (8) ND (6)	ND (8) ND (6)	74 10	105 263	44 74	<8 <6	164 1040	58 80	169 135	<8 <6	420 1800	54 67	- :	14 25	269 777	<8 <6	15 13	<8 <6
F4G PHCs (gravimetric)	ug/g dry	50	2800			-	382			1680	-	131		2540	87	-		707		<u> </u>	
Semi-Volatiles																					
Acenaphthene Acenaphthylene	ug/g dry ug/g dry	0.02	7.9 0.15	-	-	-	<0.02 <0.02	0.08 <0.02	0.02 <0.02	<0.40 <0.40	<0.02 <0.02	<0.40 <0.40	<0.02 <0.02	<0.02	0.03	<0.02 <0.02	-	0.1	<0.02 <0.02	0.02 <0.02	-
Anthracene	ug/g dry	0.02	0.67	-	-	-	0.03	0.2	0.06	0.54	<0.02	1.2	<0.02	0.15	0.13	<0.02	-	0.26	<0.02	0.06	-
Benzo[a]anthracene	ug/g dry	0.02	0.5	-	-	-	0.05	0.22	0.05	1.59	<0.02	1.76	<0.02	0.27	0.31	<0.02	-	0.83	<0.02	0.05	-
Benzo[a]pyrene Benzo[b]fluoranthene	ug/g dry ug/g dry	0.02	0.3 0.78	-	-	-	0.04	0.15 0.16	0.04	1.41	<0.02 <0.02	1.31	<0.02 <0.02	0.2	0.26 0.23	<0.02 <0.02	-	0.6	<0.02 <0.02	0.04	-
Benzo[g,h,i]perylene	ug/g dry	0.02	6.6	-	-	-	0.03	0.08	0.02	0.87	<0.02	0.65	<0.02	0.14	0.14	<0.02	-	0.26	<0.02	<0.02	-
Benzo[k]fluoranthene	ug/g dry	0.02	0.78	-	-	-	0.03	0.12	0.03	0.79	<0.02	0.78	<0.02	0.12	0.16	<0.02	-	0.42	<0.02	0.03	
Chrysene Dibenzo(a,h)anthracene	ug/g dry ug/g dry	0.02	7.0 0.1	-	-	-	0.04 <0.02	0.2	0.05 <0.02	1.92 <0.40	<0.02 <0.02	1.78	<0.02 <0.02	0.24	0.25	<0.02 <0.02	-	0.76	<0.02 <0.02	0.05 <0.02	-
Fluoranthene	ug/g dry	0.02	0.69	-	-		0.11	0.77	0.2	4.61	0.02	5.49	<0.02	0.51	0.77	<0.02	-	2	<0.02	0.18	-
Fluorene	ug/g dry	0.02	62	-	-	-	<0.02	0.09	0.03	<0.40	<0.02	0.43	<0.02	<0.02	0.02	<0.02	-	0.09	<0.02	0.02	-
Indeno [1,2,3-cd] pyrene	ug/g dry ug/g dry	0.02	0.38	-	-	-	0.02 <0.02	0.08 <0.02	<0.02 <0.02	0.73 <0.40	<0.02 <0.02	0.61 <0.40	<0.02 <0.02	0.07 <0.02	0.13 <0.02	<0.02 <0.02	-	0.25	<0.02 <0.02	<0.02 <0.02	-
1-Methylnaphthalene						-	<0.02	<0.02	<0.02	<0.40	<0.02	<0.40	<0.02	0.02	<0.02	<0.02	-	0.02	<0.02	<0.02	
1-Methylnaphthalene 2-Methylnaphthalene	ug/g dry	0.02	0.99	-	-																
2-Methylnaphthalene Methylnaphthalene (1&2)	ug/g dry ug/g dry	0.04	0.99	-	-	-	<0.04	<0.04	<0.04	<0.80	<0.04	<0.80	<0.04	<0.04	<0.04	<0.04	-	0.04	<0.04	<0.04	-
2-Methylnaphthalene	ug/g dry				-	-		<0.04 0.06 0.67	<0.04 0.04 0.19	<0.80 <0.20 2.4	<0.04 <0.01 <0.02	<0.80 0.38 4.17	<0.04 <0.01 <0.02	<0.04 0.01 0.19	<0.04 0.04 0.37	<0.04 <0.01 <0.02		0.04 0.03 1.02			

2.00 Result exceeds Reg 153/04-Table 3 Residential, coarse Standards
ND (0.2) MDL exceeds Reg 153/04-Table 3 Residential, coarse Standards
ND (0.2) No concentrations identified above the MDL
Parameter not analysed
NV No value given for indicated parameter



Parameter	Units	MDL	Regulation	BH1-GW1 1121005-01	BH1-GW1 1229249-01	BH4-GW1 1229249-03	BH4-GW2 1230199-01	BH1-GW2 1346233-01	BH4-GW3	BH2-23-GW1	BH3-23-GW1	BH5-23-GW1	BH6-23-GW1	DUP1-23-GW1	BH3-23-GW2	BH5-23-GW2		BH7-23-GW1	BH9-23-GW1	BH8-23-GW1	BH1-24-GW1	BH3-24-GW1	BH4-24-GW1
Sample Dept	pth (m)		Reg 153/04-Table 3 Non-Potable	4.4-5.64	6.1-7.62	1229249-03 18.1-21.1	1230199-01 18.1-21.1	6.1-7.62	1346233-03 18.1-21.1	2310387-01 4.6-7.6	2310387-02 4.57-7.57	2310387-03 3.78-6.78	2310387-04 4.1-7.1	2310387-05 4.57-7.57	2312554-01 4.57-7.57	2312554-02 3.78-6.78	2326362-01 3.78-6.79	2334476-01 3.83-6.88	2334476-02 6.02-9.07	2335491-01 9.12-12.17	2422099-01 7.03-10.03	2422099-02 7.26-10.26	2422099-03 7.34-10.34
Sample Da			Groundwater, coarse	16/May/2011	20/Jul/2012	20/Jul/2012	25/Jul/2012	12/Nov/2013	12/Nov/2013	8/Mar/2023	8/Mar/2023	8/Mar/2023	8/Mar/2023	8/Mar/2023	23/Mar/2023	23/Mar/2023	27/Jun/2023	23/Aug/2023	23/Aug/2023	31/Aug/2023	22/May/2024	22/May/2024	22/May/2024
Metals Mercury	ug/L	0.1	0.29			_	_					_	_	_	_	_	_	_	_	_	ND (0.1)	ND (0.1)	ND (0.1)
Antimony	ug/L	0.5	20000	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.5	0.8	0.7
Arsenic	ug/L	1	1900	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (1)	ND (1)	1
Barium Beryllium	ug/L ug/L	0.5	29000 67	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	831 ND (0.5)	882 ND (0.5)	377 ND (0.5)
Boron	ug/L	10	45000	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	403	164	172
Cadmium	ug/L	0.1	2.7	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.1)	ND (0.1)	ND (0.1)
Chromium Chromium (VI)	ug/L	1 10	810 140	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (1) ND (10)	ND (1) ND (10)	ND (1) ND (10)
Cobalt	ug/L ug/L	0.5	66	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.5)	ND (10) ND (0.5)	ND (0.5)
Copper	ug/L	0.5	87	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	2.3	2.7	2.6
Lead Molybdenum	ug/L	0.1	25 9200	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	1.1 5.7	0.4 9.7	0.6 12.1
Nickel	ug/L ug/L	1	490	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	1	2	1
Selenium	ug/L	1	63	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (1)	ND (1)	4
Silver	ug/L	0.1 200	1.5	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.1)	ND (0.1)	ND (0.1)
Sodium Thallium	ug/L ug/L	0.1	2300000 510	-	· :	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	908000	1430000 0.1	783000 0.1
Uranium	ug/L	0.1	420	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	1.6	3.9	2.5
Vanadium	ug/L	0.5	250	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	1.2	0.6	0.7
Zinc Volatiles	ug/L	5	1100	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (5)	ND (5)	ND (5)
Acetone	ug/L	5.0	130000	ND (5.0)	82.4	40.6	104	ND (5.0)	117	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	-	-	-	-	ND (5.0)	-	-	-
Benzene	ug/L	0.5	44	ND (0.5)	ND (0.5)	1.1	ND (0.5)	ND (0.5)	8	ND (0.5)	0.6	ND (0.5)	ND (0.5)	0.6	ND (0.5)	ND (0.5)	1.2	3.6	ND (0.5)	ND (0.5)	ND (0.5)	-	-
Bromodichloromethane Bromoform	ug/L ug/l	0.5	85000 380	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	3.8 ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	- :	-	-	ND (0.5) ND (0.5)	-	-	<u> </u>
Bromomethane	ug/L ug/L	0.5	5.6	ND (0.5)	ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)	<u> </u>	-	1
Carbon Tetrachloride	ug/L	0.2	0.79	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	ND (0.2)	-	-	-	-	ND (0.2)	-	-	-
Chlorobenzene Chloroform	ug/L	0.5	630 2.4	ND (0.5) ND (0.5)	ND (0.5) 16.1	ND (0.5)	ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	ND (0.5) 2.6	ND (0.5) ND (0.5)	-	-	-	-	ND (0.5) 9.6	-	-	-
Dibromochloromethane	ug/L ug/L	0.5	2.4 82000	ND (0.5) ND (0.5)	16.1 ND (0.5)	21.8 ND (0.5)	3.1 ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	2.5 ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	2.6 ND (0.5)	ND (0.5) ND (0.5)	-	-	-	-	9.6 ND (0.5)	-	-	<u> </u>
Dichlorodifluoromethane	ug/L	1.0	4400	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	-	-	-	-	ND (1.0)	-	-	-
1,2-Dichlorobenzene 1.3-Dichlorobenzene	ug/L	0.5	4600 9600	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	0.7 ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	-	-	-	ND (0.5) ND (0.5)	-	-	-
1,3-Dichlorobenzene 1,4-Dichlorobenzene	ug/L ug/L	0.5	9600	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	-	-	-	ND (0.5) ND (0.5)	-	-	-
1,1-Dichloroethane	ug/L	0.5	320	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-		-	-	ND (0.5)		-	-
1,2-Dichloroethane 1.1-Dichloroethylene	ug/L	0.5 0.5	1.6	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	-	-	-	ND (0.5) ND (0.5)	-	-	-
1,1-Dichloroethylene cis-1,2-Dichloroethylene	ug/L ug/L	0.5	1.6 1.6	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-		-	-	ND (0.5) ND (0.5)	-	-	-
trans-1,2-Dichloroethylene	ug/L	0.5	1.6	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)	-	-	-
1,2-Dichloropropane	ug/L	0.5	16	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)	-	-	-
cis-1,3-Dichloropropylene trans-1,3-Dichloropropylene	ug/L ug/L	0.5	5.2 5.2	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	-	-	-	ND (0.5) ND (0.5)	-		1
1,3-Dichloropropene, total	ug/L	0.5	5.2	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)	-	-	-
Ethylbenzene	ug/L	0.5	2300	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	1.2	ND (0.5)	ND (0.5)	17	ND (0.5)	ND (0.5)	ND (0.5)	9.2	10.7	0.9	ND (0.5)	ND (0.5)	ND (0.5)	-	-
Ethylene dibromide (dibromoethane, 1, Hexane	ug/L ug/L	0.2 1.0	0.25 51	- ND (1.0)	2.1	- ND (1.0)	- ND (1.0)	- ND (1.0)	- ND (1.0)	ND (0.2) ND (1.0)	ND (0.2) ND (1.0)	ND (0.2) ND (1.0)	ND (0.2) ND (1.0)	ND (0.2) ND (1.0)	ND (0.2) ND (1.0)	-		-	-	ND (0.2) ND (1.0)		-	-
Methyl Ethyl Ketone (2-Butanone)	ug/L	5.0	470000	ND (5.0)	10.7	8.4	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	-	-	-		ND (5.0)	-	-	-
Methyl Isobutyl Ketone	ug/L	5.0	140000	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	ND (5.0)	-	-	-	-	ND (5.0)	-	-	-
Methyl tert-butyl ether Methylene Chloride	ug/L ug/L	2.0 5.0	190 610	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	ND (2.0) ND (5.0)	-	-	-	-	ND (2.0) ND (5.0)	-	-	-
Styrene	ug/L	0.5	1300	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)	-	-	-
1,1,1,2-Tetrachloroethane	ug/L	0.5	3.3	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)	-	-	-
1,1,2,2-Tetrachloroethane Tetrachloroethylene	ug/L ug/L	0.5	3.2 1.6	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	-	-	-	ND (0.5) ND (0.5)	-	-	-
Toluene	ug/L	0.5	18000	ND (0.5)	6.2	5.6	ND (0.5)	ND (0.5)	1.4	ND (0.5)	3.4	1	ND (0.5)	3.5	ND (0.5)	2.4	1	1.5	ND (0.5)	ND (0.5)	ND (0.5)	-	-
1,1,1-Trichloroethane	ug/L	0.5	640	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)	-	-	-
1,1,2-Trichloroethane Trichloroethylene	ug/L ug/L	0.5	4.7 1.6	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-			-	ND (0.5) ND (0.5)		-	-
Trichlorofluoromethane	ug/L	1.0	2500	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	-	-	-		ND (1.0)	-	-	-
Vinyl Chloride	ug/L	0.5	0.5	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	-	-	-	-	ND (0.5)		-	-
m/p-Xylene o-Xylene	ug/L ug/L	0.5	4200 4200	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	1.6	ND (0.5) ND (0.5)	3.2	22	ND (0.5) ND (0.5)	3.2	ND (0.5)	11.5	12.4 0.9	6 0.8	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	<u> </u>
Xylenes, total	ug/L	0.5	4200	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	2.8	ND (0.5)	4.1	24	ND (0.5)	4.2	ND (0.5)	12.9	13.2	6.8	ND (0.5)	ND (0.5)	ND (0.5)	-	<u> </u>
BTEX																							
Benzene Ethylbenzene	ug/L ug/l	0.5	44 2300	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	1.1 ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	8 1.2	ND (0.5) ND (0.5)	0.6 ND (0.5)	ND (0.5) 17	ND (0.5) ND (0.5)	0.6 ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) 9.2	1.2	3.6 0.9	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	-
Toluene	ug/L ug/L	0.5	18000	ND (0.5)	6.2	ND (0.5) 5.6	ND (0.5) ND (0.5)	ND (0.5)	1.4	ND (0.5)	3.4	1	ND (0.5) ND (0.5)	3.5	ND (0.5)	2.4	1	1.5	ND (0.5)	ND (0.5)	ND (0.5)	-	
m/p-Xylene	ug/L	0.5	4200	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	ND (0.5)	1.6	ND (0.5)	3.2	22	ND (0.5)	3.2	ND (0.5)	11.5	12.4	6	ND (0.5)	ND (0.5)	ND (0.5)	-	-
o-Xylene Xylenes, total	ug/L	0.5	4200 4200	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	1.1	ND (0.5) ND (0.5)	4.1	2 24	ND (0.5) ND (0.5)	4.2	ND (0.5) ND (0.5)	1.4	0.9	0.8	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	ND (0.5) ND (0.5)	-	-
Xylenes, total Hydrocarbons	ug/L	U.5	4200	ND (U.5)	ND (U.5)	ND (U.5)	ND (U.5)	ND (U.5)	2.8	ND (U.5)	4.1	24	ND (U.5)	4.2	IND (U.5)	12.9	13.2	6.8	ND (U.5)	ND (U.5)	ND (U.5)	-	<u> </u>
F1 PHCs (C6-C10)	ug/L	25	750	ND (25)	45	ND (25)	ND (25)	-	-	ND (25)	-	1210	ND (25)	-	ND (25)	742	1050	450	ND (25)	ND (25)	ND (25)	-	-
F2 PHCs (C10-C16)	ug/L	100	150	ND (100)	ND (100)	ND (100)	ND (100)	-	-	ND (100)	-	266	ND (100)	-	ND (100)	ND (100)	398	ND (100)	ND (100)	ND (476)	ND (100)	-	-
F3 PHCs (C16-C34) F4 PHCs (C34-C50)	ug/L ug/L	100 100	500 500	ND (100) ND (100)	ND (100) ND (100)	ND (100) ND (100)	ND (100) ND (100)		-	ND (100) ND (100)	-	ND (100) ND (100)	ND (100) ND (100)	-	ND (100) ND (100)	ND (476) ND (476)	ND (100) 282	-	-				
Semi-Volatiles	ug/L	200	300	145 (100)	(100)	(100)	(100)		<u> </u>	(100)	 	.15 (100)	(100)	l -	(100)	(100)	(100)	.15 (100)	.45 (100)	(470)	202	<u> </u>	
Acenaphthene	ug/L	0.05	600	-	-	-	-	-		-	-	-	-	-	-	-	-	-	-	-	ND (0.05)	ND (0.05)	ND (0.05)
Acenaphthylene Anthracene	ug/L ug/L	0.05	1.8 2.4	-		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.05) ND (0.01)	ND (0.05) ND (0.01)	ND (0.05) ND (0.01)
Benzo[a]anthracene	ug/L ug/L	0.01	4.7	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.01)	ND (0.01) ND (0.01)	ND (0.01)
Benzo[a]pyrene	ug/L	0.01	0.81	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.01)	ND (0.01)	ND (0.01)
	ug/L	0.05	0.75	-	-	-	-	-	•	-	-	-	-	-	-	-	-	-	-	-	ND (0.05)	ND (0.05)	ND (0.05)
Benzo[b]fluoranthene	ug/L	0.05	0.2	-	1	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)	ND (0.05) ND (0.05)
Benzo[b]fluoranthene Benzo[g,h,i]perylene	ug/L		1.0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.05)	ND (0.05)	ND (0.05)
Benzo[b]fluoranthene Benzo[g,h,i]perylene Benzo[k]fluoranthene Chrysene	ug/L ug/L	0.05				-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.05)	ND (0.05)	ND (0.05)
Benzo[b]fluoranthene Benzo[g,h,i]perylene Benzo[k]fluoranthene Chrysene Dibenzo[a,h]anthracene	ug/L ug/L	0.05	0.52						-	-	-	-	-	-	-	-	-	-	-	-			ND (0.01)
Benzo[b]fluoranthene Benzo[g_h,i]perylene Benzo[k]fluoranthene Chrysene Dibenzo[a,h]anthracene Fluoranthene	ug/L ug/L ug/L	0.05 0.01	130	-	-	-	-						-								ND (0.01)	ND (0.01)	
Benzo[b]fluoranthene Benzo[g,h,i]perylene Benzo[k]fluoranthene Chrysene Dibenzo[a,h]anthracene	ug/L ug/L	0.05			-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND (0.01) ND (0.05) ND (0.05)	ND (0.01) ND (0.05) ND (0.05)	ND (0.05) ND (0.05)
Benzo[pliuoranthene Benzo[pliuoranthene Benzo[pliuoranthene Chrysene Dibenzo[a,h]anthracene Fluoranthene Fluoranthene Indeno [1,2,3-cd] pyrene 1-Methyinaphthalene	ug/L ug/L ug/L ug/L ug/L ug/L	0.05 0.01 0.05 0.05 0.05	130 400 0.2 1800	-	-					-			- - -			-	-	-		-	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)
Benzo[p/huoranthene Benzo[p/huoranthene Benzo[p/huoranthene Chrysene Dibenzo[a,h]anthracene Fluoranthene Fluorene Indeno [1,2,3-cd] pyrene 1-Methylnaphthalene 2-Methylnaphthalene	ug/L ug/L ug/L ug/L ug/L ug/L ug/L	0.05 0.01 0.05 0.05 0.05 0.05	130 400 0.2 1800		-	- - - -		-			-	-			-		-		- - -		ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05) ND (0.05)
Benzo[g.h]perylene Benzo[g.h]perylene Benzo[g.h]perylene Benzo[g.h]perylene Chrysene Dibenzo[a,h]anthracene Fluoranthene Fluoranthene Fluoranthene L-Methynaphthalene 2-Methylnaphthalene Methylnaphthalene Methylnaphthalene	ug/L ug/L ug/L ug/L ug/L ug/L ug/L ug/L	0.05 0.01 0.05 0.05 0.05	130 400 0.2 1800		-				- - - -	-		-			-	-	-				ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)	ND (0.05) ND (0.05) ND (0.05)
Benzolg/Huoranthene Benzolg/H, i)perylene Benzolg/Huoranthene Chrysene Dibenzolg/Hanthracene Fluoranthene Fluoranthene Indeno (1,2,3-cd) pyrene 1-Mettylnaphthalene 2-Mettylnaphthalene	ug/L ug/L ug/L ug/L ug/L ug/L ug/L	0.05 0.01 0.05 0.05 0.05 0.05 0.05	130 400 0.2 1800 1800			- - - -		-			-			-		-	-		- - -		ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.10)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.10)	ND (0.05) ND (0.05) ND (0.05) ND (0.05) ND (0.10)

2.00 Result exceeds Reg 153/04-Table 3 Non-Potable Groundwater, coarse Standards ND (0.2) MDL exceeds Reg 153/04-Table 3 Non-Potable Groundwater, coarse Standards ND (0.2) No concentrations identified above the MDL Parameter not analysed NV No value given for indicated parameter

APPENDIX 1

SAMPLING AND ANALYSIS PLAN
SOIL PROFILE AND TEST DATA SHEETS
SYMBOLS AND TERMS
LABORATORY CERTIFICATES OF ANALYSIS



Sampling & Analysis Plan

137-141 George Street and 110-116 York Street Ottawa, Ontario

Prepared for Claridge Homes





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1.0 SAMPLING PROGRAM

Paterson Group Inc. (Paterson) was commissioned by Claridge Homes, to conduct a Phase II – Environmental Site Assessment (Phase II ESA) for the property addressed 141 George Street, in the City of Ottawa, Ontario.

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

Borehole	Location & Rationale	Proposed Depth & Rationale
BH1-23	Eastern portion of the Phase II Property; to assess potential impacts resulting from the presence of fill material of unknown quality as well as general coverage for the geotechnical program.	12-15 m; to intercept the groundwater table for the purpose of installing a monitoring well as part of the geotechnical program.
BH2-23	Northern portion of the Phase II Property; to assess potential impacts resulting from the importation of fill material of unknown quality.	5-8 m; to intercept the groundwater table for the purpose of installing a monitoring well.
BH3-23	Western portion of the Phase II Property; to assess potential impacts resulting from former off-site dry cleaner and machine shop.	5-8 m; to intercept the groundwater table for the purpose of installing a monitoring well.
BH4-23	Northwestern portion of the Phase II Property; to assess potential impacts resulting from the importation of fill material of unknown quality.	12-15 m to provide general coverage for the environmental and geotechnical programs.
BH5-23	Southwestern portion of the Phase II Property; to assess potential impacts resulting from former onsite aboveground storage tank (AST), off-site retail fuel outlet and various former off-site industries.	5-8 m; to intercept the groundwater table for the purpose of installing a monitoring well.
BH6-23	Northwestern portion of the Phase II Property; to assess potential impacts resulting from the presence of fill material of unknown quality.	5-8 m; to intercept the groundwater table for the purpose of installing a monitoring well.
BH7-23	Southwestern portion of the Phase II Property; to delineate soil and/or groundwater impacts	5-7m; to intercept groundwater table for the purpose of installing a monitoring well and laterally delineate impacts identified at BH5-23
BH8-23		9-12m; to vertically delineate groundwater impacts identified at BH5-23
BH9-23		6-9m; to laterally delineate impacts identified at BH5-23
BH10-23		3-6m; shallow monitoring well installation, to determine if groundwater present in overburden
BH1-24	To provide general coverage of the site fill and proposed building footprint for environmental and	7-10m; to access groundwater table if required
BH2-24	geotechnical purposes.	4m – assess fill and underlying native
BH3-24		7-10m; to access groundwater table if
BH4-24		required
BH5-24		4m – assess fill and underlying native

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Borehole locations are shown on Drawing PE2709-11 – 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. 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 three boreholes to allow for the collection of groundwater samples.

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2.0 ANALYTICAL TESTING PROGRAM

The analytical testing program for soil at the Phase II Property 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 Phase I Property 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.

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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:

J	Glass soil sample jars	
J	two buckets	
	cleaning brush (toilet brush works well)	
	dish detergent	
	methyl hydrate	
]	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.

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Drilling Procedure

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

	Continuous split spoon samples (every 0.6 m or 2') or semi-continuous (every
	0.76 m or 2'6") are required.
	Make sure samples are well sealed in plastic bags with no holes prior to screening and are kept cool but unfrozen.
	If sampling for VOCs, BTEX, or PHCs F ₁ , 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.
Sp	oon Washing Procedure
	sampling equipment (spilt spoons, etc.) must be washed between samples in der 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 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.
J	Turn instrument on and allow to come to zero - calibrate if necessary
J	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.
J	Insert probe into soil bag, creating a seal with your hand around the opening.
J	Gently manipulate soil in bag while observing instrument readings.
J	Record the highest value obtained in the first 15 to 25 seconds
J	Make sure to indicate scale (ppm or LEL); also note which instrument was used
	(RKI Eagle 1 or 2, or MiniRae).
J	Jar samples and refrigerate as per Sampling and Analysis Plan.

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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 ½" 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

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surface.



3.3 Monitoring Well Sampling Procedure

Εq	uipment
	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 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 pH/Temperature/Conductivity combo pen Laboratory-supplied sample bottles
Sa	mpling 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.

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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 laboratory-provided 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.

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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

body of the Phase II ESA repor

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

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1 2022

May 1, 2023

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

TYPE ON TYPE	NUMBER WAS			July 17, 2 DEPTH			PE270 HOLE NO BH 1 Onization) .	etor	
<u></u> AU		IPLE		DEPTH				n Detec	tor	
<u></u> AU			图口	ł I	ELEV.		onization	Detec	tor	l —
<u></u> AU	NUMBER	% OVER	5 0	(m)	ELEV. (m)	Photo Ionization Detector ■ Volatile Organic Rdg. (ppm)				
1 . / 1		ы	N VALUE or RQD				Explos			Monitoring Well Construction
1 . / 1	1	α.	1	0-	61.69	20	40 6	60 8 	0 : : : 	
<u> </u>	2									
ss	3	42	9	1-	-60.69					
ss	4	50	26	2-	-59.69					
ss	5	33	9							
ss S	6	100	50+	3-	-58.69					
ss	7	50	12	4-	-57.69					
⊔ ≍ SS	8	67	50+							
				5-	56.69					
RC	1	92	67	6-	-55.69					
RC 	2	100	100	7-	-54.69					
							agle Rd	g. (ppn	1)	000
	SS SS SS SS RC	SS 4 SS 5 SS 6 SS 7 SS 8 RC 1	SS 4 50 SS 5 33 SS 6 100 SS 7 50 SS 8 67 RC 1 92	SS 4 50 26 SS 5 33 9 SS 6 100 50+ SS 7 50 12 SS 8 67 50+ RC 1 92 67	SS 4 50 26 2- SS 5 33 9 SS 6 100 50+ SS 7 50 12 SS 8 67 50+ RC 1 92 67 6- 7	SS 4 50 26 2-59.69 SS 5 33 9 SS 6 100 50+ 3-58.69 SS 7 50 12 4-57.69 SS 8 67 50+ 5-56.69 RC 1 92 67 6-55.69	SS 4 50 26 2-59.69 SS 5 33 9 SS 6 100 50+ 3-58.69 SS 7 50 12 4-57.69 SS 8 67 50+ 5-56.69 RC 1 92 67 6-55.69 RC 2 100 100 7-54.69	SS 4 50 26 SS 5 33 9 SS 6 100 50+ SS 7 50 12 SS 8 67 50+ RC 1 92 67 RC 2 100 100 7-54.69 100 200 3 RKI Eagle Rd	SS 4 50 26 2-59.69 SS 5 33 9 SS 6 100 50+ 3-58.69 SS 7 50 12 4-57.69 SS 8 67 50+ 5-56.69 RC 1 92 67 6-55.69 RC 2 100 100 7-54.69	SS 4 50 26 2-59.69 SS 5 33 9 SS 6 100 50+ 3-58.69 SS 7 50 12 4-57.69 SS 8 67 50+ 5-56.69 RC 1 92 67 6-55.69 RC 2 100 100 7-54.69

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SOIL PROFILE AND TEST DATA

DATUM Geodetic								FILE NO. PE2709			
REMARKS								HOLE NO.			
BORINGS BY CME 55 Power Auger					DATE .	July 17, 2 	BH 2				
SOIL DESCRIPTION		PLOT		SAMPLE		DEPTH (m)	ELEV. (m)	Photo Ionization Detector ■ Volatile Organic Rdg. (ppm)			
		TYPE	NUMBER	RECOVERY	N VALUE or RQD			O Lower Explosive Limit %	Monitoring Well Construction		
GROUND SURFACE	STRATA			2	Z	0-	-62.02	20 40 60 80	≥		
Asphaltic concrete 0.05		AU	1 2				0=:0=				
FILL: Brown silty sand with gravel		₩ ΛΟ	_								
1.22	\bowtie	<u>∦</u> ss	3	29	6	1-	61.02				
GLACIAL TILL: Brown silty clay		ss	4	21	23	2-	-60.02				
with sand, gravel, cobbles, boulders		ss	5	42	13						
		N 22	5	42	13		50.00				
3.30	\^^^^	∑.ss	6	36	50+	3-	-59.02				
	\^^^^/										
GLACIAL TILL: Brown silty sand	\^^^^	⊠ SS	7	80	50+	4-	58.02				
with gravel, cobbles, boulders											
4.90		ss	8								
		D0		100	47	5-	-57.02				
		RC	1	100	47		6+56.02				
						6-					
		D 0		100	100	0 00.02	00.02				
BEDROCK: Grey limestone		RC	2	100	100		7-55.02				
						7-					
		RC	3	100	100		E4.00				
		no	3	100	100	0-	-54.02				
8.76											
End of Borehole											
(GWL @ 4.50m-July 20, 2012)											
									500		
								RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim	١.		

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SOIL PROFILE AND TEST DATA

DATUM Geodetic									FILE NO. PE2709		
REMARKS									HOLE NO.		
BORINGS BY CME 55 Power Auger				D	ATE .	July 17, 2	012		BH 3		
	TO	SAMPLE SAMPLE				DEPTH	ELEV.	Photo Ionization Detector			
SOIL DESCRIPTION			۲ ×		臣〇	(m)	(m)	● Vola	tile Organic Rdg	. (ppm)	ing
	STRATA	TYPE	NUMBER	% RECOVERY	N VALUE or RQD			O Lowe	r Explosive L	.imit %	Monitoring Well Construction
GROUND SURFACE			ų	꿆	Z O		-61.53	20	40 60	80	Σ
Asphaltic concrete 0.05		AU AU	1				01.00				
FILL: Brown silty sand with gravel, crushed stone		iiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiii	2	77	50+	1-	-60.53				
1.40		<u></u>					00.00				
	\^,^,^,	ss	4	0	15	2-	-59.53				
		Π									
GLACIAL TILL: Brown silty sand with gravel, cobbles, boulders	\^^^^	ss	5	50	17						
with gravel, cobbles, boulders	\^^^^	_ V	_			3-	-58.53				
		ss	6	50	34						
4.24	\^^^^	- SS	7	0	50+	4-	-57.53				
End of Borehole											
Practical refusal to augering at 4.24m depth								100	200 300	400 50	00
								RKI E	Eagle Rdg. (p as Resp. △ Met	pm)	υU

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SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 110 York St., 321 Dalhousie St. & 167-141 George St. Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. BH 4 BORINGS BY CME 55 Power Auger **DATE** July 18, 2012 **SAMPLE Photo Ionization Detector** STRATA PLOT DEPTH ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD NUMBER **Lower Explosive Limit %** N o v **GROUND SURFACE** 80 0+61.9125mm Asphaltic concrete over 0.30K 1 crushed stone 2 1 + 60.91SS 3 25 3 FILL: Brown silty sand with gravel, cobbles, brick and concrete SS 4 42 27 2 + 59.912.21 SS 5 33 24 3+58.91GLACIAL TILL: Brown silty sand SS 6 50 33 with gravel, cobbles, boulders \times SS 7 40 50 +4+57.914.70 5+56.91RC 1 100 72 6+55.91RC 2 100 100 7 + 54.918+53.91**BEDROCK:** Grey limestone RC 3 100 100 9+52.91RC 4 100 100 10+51.91 11 + 50.91RC 5 100 87 12+49.91 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

SOIL PROFILE AND TEST DATA

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DATUM Geodetic									PE27		
REMARKS									HOLE N		
BORINGS BY CME 55 Power Auge	er 				DATE	July 18, 2	2012		BH 4		Τ
SOIL DESCRIPTION			SAMPLE			i I	ELEV.	Photo Vol	Monitoring Well Construction		
		TYPE	NUMBER	% RECOVERY	VALUE r RQD	(m)	(,	○ Lower Explosive Limit %			
GROUND SURFACE	STRATA	Ţ	NOM	RECO	N N			20 Low		60 80	₹ S S
	1 1 1 1 1 1 1 1 1 1 1 1 1 1					12-	49.91				
		RC	6	100	100	13-	-48.91				
							47.04				
	1 1 2 1 2 1 2 1 1 1 1 1 1 1 2 1 1 2 1 1 2 1 1 3 1 1	RC	7	100	100	14-	-47.91				
						15-	46.91				
BEDROCK: Grey limestone		RC	8	100	98	16-	-45.91				ներհանդմուկունդներն անդուներն հանդուներն հանդուներն անդուներն հանդուներն առուսան անդանություն անդուների անահայան հանդում անդուների անդուների անդուների
,		RC 9	9 100		17-	44.91					
				100	18-	-43.91					
		RC	10	100	100	10	40.04				
						19-	42.91				
		RC	11	100	100	20-	41.91				
End of Borehole	21.13					21-	-40.91				
(GWL @ 4.43m-July 20, 2012)											
									Eagle Rd		500

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SOIL PROFILE AND TEST DATA

DATUM Geodetic					'				LE NO. E2709		
REMARKS						August 8,		Н	HOLE NO.		
BORINGS BY CME 55 Power Auger						BH 5					
SOIL DESCRIPTION				IPLE 문	ы	DEPTH (m)	ELEV. (m)	Photo Ionization Detector Volatile Organic Rdg. (ppm)			
		TYPE	NUMBER	% RECOVERY	N VALUE or RQD				xplosive Limit %	Monitoring Well Construction	
GROUND SURFACE	.· ∧ . ∧. ∧	~		2	4	0-	61.79	20 4	0 60 80	_	
Asphaltic concrete 0.10		AU	1								
FILL: Dark brown silty clay, some sand, gravel, crushed stone		ss	2	17	9	1-	-60.79				
- some mortar by 1.5m depth		∆ Vss	3	21	3						
2.21		ss	4	50	23	2-	-59.79				
	\^,^,^, \^,^,^,	∑ ss	5	67	33	3-	-58.79				
GLACIAL TILL: Brown silty clay with sand, gravel, cobbles		∑ ss	6	62	53	4-	-57.79				
		Δ 7									
	\^^^^	ss	7	67	61	5-	-56.79				
<u>5.41</u> End of Borehole	1^^^^	×.SS	8		50+						
Practical refusal to augering at 5.41m depth								100 20	0 300 400 50		
								RKI Eag	0 300 400 50 I le Rdg. (ppm) Resp. △ Methane Elim.	00	

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SOIL PROFILE AND TEST DATA

DATUM Geodetic									FILE NO. PE2709	
REMARKS									HOLE NO.	
BORINGS BY CME 55 Power Auger					ATE	August 8,	2012	BH 6	_	
SOIL DESCRIPTION	PLOT		SAMPLE		EJ.	DEPTH (m)	ELEV. (m)	Photo I ● Vola	ng Wel uction	
	STRATA	TYPE	NUMBER	% RECOVERY	N VALUE		-61.45	O Lowe	Monitoring Well Construction	
GROUND SURFACE		×.		2	z °	0-		20	40 60 80	2
Asphaltic concrete 0.10) W	AU	1							
FILL: Brown silty sand with gravel, crushed stone, trace clay		ss	2	42	4	1-	60.45			
		ss	3	54	8	2-	2-59.45			
2.	4 () () () () () () () () () (ss	4	58	40					
GLACIAL TILL: Brown silty clay with sand, gravel, cobbles		ss	5	86	50+	3-	-58.45			
with Sand, gravel, cobbies		× SS	6	80	50+	4-	-57.45			
End of Borehole	5 ^^^^	∑.SS	7	43	50+					
Practical refusal to augering at 4.75m depth										
									200 300 400 5 Eagle Rdg. (ppm) as Resp. △ Methane Elim.	00

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SOIL PROFILE AND TEST DATA

DATUM Geodetic					•				FILE NO. PE270	9	
REMARKS									HOLE NO		
BORINGS BY CME 55 Power Auger					ATE /	August 8,	2012		BH 7		
SOIL DESCRIPTION	PLOT	SAMPLE DEPTH ELEV. (m) (m)					1	Photo Ionization Detector ■ Volatile Organic Rdg. (ppm)			
		TYPE	NUMBER	% RECOVERY	N VALUE or RQD			O Lowe	r Explosiv	ve Limit %	Monitoring Well Construction
GROUND SURFACE	STRATA		ı	R	Z	0-	-61.02	20	40 60	0 80	Ν
Asphaltic concrete 0.10		§ AU	1				01.02				
		\$ 70	'								
		ss	2	17	11	1-	-60.02				
FILL: Brown silty sand with gravel,		7									
cobbles, concrete, asphalt, trace organics		SS	3	25	17	2-	2-59.02				
		∇									
		SS	4	75	26					- 1 - 2 - 1 - 2 - 2 - 2 - 2 - 2 - 2 - 2	
3.20		₩- <u>-</u>	5	21	0.1	3-	-58.02				
		ss	Э	21	21						
GLACIAL TILL: Grey-brown silty clay with sand, gravel, cobbles		\sqrt{ss}	6	29	23	4-	57.02				
	\^^^^ \^^^^	ΔΙ									
End of Borehole		×.55	7	80	50+						
Practical refusal to augering at											
4.70m depth											
								100	200 30		00
									agle Rdg as Resp. △	J. (ppm) Methane Elim.	

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 110 York St., 321 Dalhousie St. & 167-141 George St. Ottawa, Ontario

DATUM Geodetic									PE2709	
REMARKS									HOLE NO.	
BORINGS BY CME 55 Power Auger				D	ATE /	August 8,	2012		BH 8	
	Ħ		SAN	IPLE				Photo Id	onization Detector	ا ا
SOIL DESCRIPTION	PLOT			l		DEPTH (m)	ELEV. (m)	Volat	tile Organic Rdg. (ppm)	Monitoring Well Construction
	T.	둳	ER	% RECOVERY	N VALUE or RQD	(,	(111)			orin
	STRATA	TYPE	NUMBER	% O	YA.			O Lowe	r Explosive Limit %	Sonit
GROUND SURFACE			4	滋	z °	0-	61.20	20	40 60 80	Σ
Asphaltic concrete 0.10		≱ AU	1				01.20			
FILL Prown silty aand trace slav			-							
FILL: Brown silty sand, trace clay and gravel		∦ ss	2	29	11	1-	60.20			
		<u>'</u>								
1.83		∦-ss	3	58	30	2-	-59.20			
	^^^^	7				_	00.20			
GLACIAL TILL: Brown to grey silty clay with sand, gravel, cobbles	\^^^^ \^^^^	\ ss	4	46	69					
ciay with sand, graver, cobbles		- V 00	_	F0	F0	3-	-58.20			
3.53 End of Borehole	^^^^	SS	5	52	59					
Practical refusal to augering at 3.53m depth										
·										
								100		oo
									Eagle Rdg. (ppm) as Resp. △ Methane Elim.	

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 110 York St., 321 Dalhousie St. & 167-141 George St. Ottawa, Ontario

DATUM Geodetic									FILE NO. PE2709	
REMARKS							0010		HOLE NO.	
BORINGS BY CME 55 Power Auger					ATE /	August 8,	2012		BH 9	
SOIL DESCRIPTION	A PLOT			IPLE	шо	DEPTH (m)	ELEV. (m)		onization Detectitile Organic Rdg. (pp	> -
	STRATA	TYPE	NUMBER	% RECOVERY	N VALUE or RQD				r Explosive Limi	% % % % % % % % % % % % % % % % % % %
GROUND SURFACE	\.· ^ ^ . /	x .		<u> </u>	-	0-	61.79	20	40 60 80	
Asphaltic concrete 0.10 FILL: Brown silty sand with gravel, clay, brick, mortar		AU SS	1 2		4		-60.79			3 · 3 · 3 · 3 · 3 · 3 · 3 · 3 · 3 · 3 ·
End of Borehole	\bowtie	<u>{\}</u> -								<u>: : : </u>
Practical refusal to augering at 1.40m depth								100 RKIE	200 300 400 Eagle Rdg. (ppm)	

SOIL PROFILE AND TEST DATA

FILE NO.

141 George Street

9 Auriga Drive, Ottawa, Ontario K2E 7T9

Geodetic

Phase II - Environmental Site Assessment Ottawa, Ontario

DATUM

REMARKS								PE2289	
BORINGS BY Portable Drill				D	ATE I	May 11, 2011		HOLE NO. BH 1-11	
	PLOT		SAN	/IPLE		DEPTH ELEV.	1	onization Detector	
SOIL DESCRIPTION		Ä	SER	ÆRY	VALUE r RQD	(m) (m)		tile Organic Rdg. (ppm)	Construction
	STRATA	TYPE	NUMBER	% RECOVERY	N VA or F			r Explosive Limit %	S
GROUND SURFACE		1.		K	-	0+61.30	20	40 60 80	
Interlocking brick 0.06		SS	1	58	9			74,014,010,010	
FILL: Brown silty sand with gravel		ss	2	8	4	1-60.30			
1.37	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	ss	3	33	46				
GLACIAL TILL: Brown silty sand with gravel and clay, trace cobbles		ss	4	53	50+	2-59.30			
2.26	^^^^^ ^^^^^	RC	1	32					
GLACIAL TILL: Grey silty clay with sand, gravel, cobbles and boulders	^^^^^ ^^^^^	_				3+58.30			
	^^^^^ ^^^^^ ^^^^	RC RC	2	18					
<u>3.9</u> 6	^^^^^ ^^^^^	RC 	4	64		4-57.30			
		RC	5	100		4 37.30			
BEDROCK: Grey limestone		- RC	6	71		5-56.30			
5.64		_ _RC	7						Y
End of Borehole									
(GWL @ 5.20m - May 16, 2011)							100	200 300 400 500	
							RKI E	Eagle Rdg. (ppm) as Resp. △ Methane Elim.	

9 Auriga Drive, Ottawa, Ontario K2E 7T9

Geodetic

SOIL PROFILE AND TEST DATA

FILE NO.

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM PE2709 REMARKS HOLE NO. BORINGS BY CME-55 Low Clearance Drill **BH 1-23** DATE February 24, 2023 **SAMPLE Photo Ionization Detector** PLOT DEPTH ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD STRATA NUMBER **Lower Explosive Limit %** N o v **GROUND SURFACE** 80 0+61.52Asphaltic concrete 0.05 1 FILL: Crushed stone 0.30 FILL: Dark brown silty sand, some 0.69 1 + 60.52SS 2 29 14 \gravel FILL: Brown silty sand with gravel SS 3 20 58 and crushed stone 2+59.52SS 4 23 2.74 54 3+58.52SS 5 86 50 +**GLACIAL TILL:** Dense to very 4+57.52 dense, brown silty sand to sandy silt RC 1 with gravel, cobbles and boulders 5+56.525.36 RC 2 100 87 6+55.527 + 54.52100 RC 3 100 8+53.52RC 4 100 100 9+52.52**BEDROCK:** Good to excellent 10+51.52 5 RC 100 100 quality, grey limestone 11 + 50.52RC 6 100 100 12+49.52 13+48.527 100 RC 100 14 + 47.52RC 8 100 100 15 + 46.5215.32 End of Borehole (GWL @ 4.49m - March 8, 2023) 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. BORINGS BY CME-55 Low Clearance Drill **BH 2-23** DATE February 24, 2023 **SAMPLE Photo Ionization Detector** PLOT DEPTH ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD STRATA NUMBER **Lower Explosive Limit %** N or v **GROUND SURFACE** 80 0+61.53Asphaltic concrete 0.06 1 0.69 **FILL:** Crushed stone with sand 1+60.53SS 2 42 5 FILL: Brown silty sand, some SS 3 gravel, wood, crushed stone, trace 58 15 2+59.53brick SS 4 62 38 3+58.53SS 5 100 50+ **GLACIAL TILL:** Dense to very 50+ ss 🛚 6 100 4+57.53 dense, brown silty sand to sandy silt with gravel, cobbles and boulders RC 1 100 5+56.535.64 RC 2 100 100 6+55.53**BEDROCK:** Excellent quality, grey limestone 7+54.53 RC 3 100 100 End of Borehole (GWL @ 4.98m - March 8, 2023) 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. **BH 3-23** BORINGS BY CME-55 Low Clearance Drill DATE February 27, 2023 **SAMPLE Photo Ionization Detector** PLOT DEPTH ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD STRATA NUMBER **Lower Explosive Limit %** N or v **GROUND SURFACE** 80 0+61.87Asphaltic concrete 0.06 2 ፟

尽 AU FILL: Dark brown sand with grave 0.69 and crushed stone 1+60.873 50 18 SS FILL: Brown silty sand to sandy silt, Isome gravel, cobbles, trace topsoil SS 4 33 58 2+59.87**GLACIAL TILL:** Compact to dense. 3+58.87brown silty sand to sandy silt with SS 5 42 27 gravel, cobbles and boulders 4 + 57.87SS 6 100 50 +4.95 5+56.87RC 1 100 100 6+55.87**BEDROCK:** Excellent quality, grey limestone RC 2 100 100 7 + 54.877.57 End of Borehole (GWL @ 3.48m - March 8, 2023) 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. BORINGS BY CME-55 Low Clearance Drill **BH 4-23** DATE February 27, 2023 **SAMPLE Photo Ionization Detector** PLOT DEPTH ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD STRATA NUMBER **Lower Explosive Limit %** N o v **GROUND SURFACE** 80 0+61.03Asphaltic concrete 0.05 2 ΑU FILL: Crushed stone with sand 0.25 FILL: Brown sand 1 + 60.03SS 3 25 18 - some crushed stone by 0.6m dept45 SS 4 46 24 2+59.03GLACIAL TILL: Compact to dense, brown silty sand with gravel, cobbles 3+58.03and boulders RC 1 35 4 + 57.03 4.65 5+56.03RC 2 100 93 6+55.03RC 3 100 100 7 + 54.038+53.03RC 4 100 100 9+52.03**BEDROCK:** Excellent quality, grey 5 RC 100 100 10+51.03 limestone 11 + 50.03RC 6 100 100 12+49.03 7 RC 100 100 13+48.0314 + 47.03RC 8 100 100 15 ± 46.03 End of Borehole 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. **BH 5-23** BORINGS BY CME-55 Low Clearance Drill DATE February 28, 2023 **SAMPLE Photo Ionization Detector** PLOT DEPTH ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD STRATA NUMBER **Lower Explosive Limit %** N or v **GROUND SURFACE** 80 0+62.09Asphaltic concrete 0.06 1 FILL: Dark brown silty sand, some0.25 1 + 61.092 50 50 FILL: Dark brown silty sand with Igravel and crushed stone SS 3 22 33 2+60.09**GLACIAL TILL:** Compact to very SS 4 92 50+ dense, brown silty sand to sandy silt 3+59.09with gravel, cobbles and boulders SS 5 67 37 4 + 58.09SS 6 96 50 +- grey by 4.1m depth 7 SS 50+ 100 4.93 5+57.09RC 1 100 76 **BEDROCK:** Good to excellent 6+56.09quality, grey limestone RC 2 100 100 End of Borehole (GWL @ 2.51m - March 8, 2023) 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. **BH 6-23** BORINGS BY CME-55 Low Clearance Drill DATE February 28, 2023 **SAMPLE Photo Ionization Detector** STRATA PLOT DEPTH ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY N VALUE or RQD NUMBER **Lower Explosive Limit % GROUND SURFACE** 80 0+62.08Asphaltic concrete 0.06 1 1+61.08 SS 2 54 18 FILL: Dark brown sand with crushed SS 3 26 stone 58 2+60.08SS 4 19 58 3+59.08SS 5 70 50+ 3.58 4+58.08RC 1 100 Reinforced concrete slab (former crane base) 5.18 5+57.08RC 2 100 95 6+56.08**BEDROCK:** Excellent quality, grey limestone RC 3 100 100 7+55.08 End of Borehole (GWL @ 2.77m - March 8, 2023) 100 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic									FILE NO.		
REMARKS BORINGS BY CME-55 Low Clearance	Drill			Б	ΔTF	August 1	4 2023		HOLE NO	Э.	
Borning BT OWL 30 LOW Olcaranoc			SAN	ИPLE	AIL .			Photo		n Detector	= -
SOIL DESCRIPTION	PLOT				_	DEPTH (m)	ELEV. (m)			Rdg. (ppm)	Monitoring Well Construction
	STRATA	TYPE	NUMBER	RECOVERY	VALUE			O Lowe	r Evnloe	ive Limit %	itorir nstru
GROUND SURFACE	STI	Ţ	Ŋ	RECO	N O N			20	-	60 80	Moo
Asphaltic concrete 0.08		AU	1			0-	62.03		Δ		
FILL: Brown silty sand wtih gravel and crushed stone, trace brick 1.45		√ ss	2	8	18	1-	61.03				
1.40		∯- ∏ss	3	50	14	2.	-60.03		Δ		
CLACIAL TILL. Compost to your		X ss	4	50	38	2	+60.03				
GLACIAL TILL: Compact to very dense, brown silty sand with gravel, cobbles and boulders		∑ SS	5	67	50+	3-	59.03	Δ			
cobbles and boulders		= SS	6	100	50+	4-	-58.03	4			
4. <u>6</u> 7	, , , , , , , , , , , , , , , , , , , ,	=-SS	7	100	50+			4			
DEDDOOK Overlage collect						5-	-57.03				
BEDROCK: Good to excellent quality, grey limestone						6-	56.03				
6.88											
End of Borehole											
								100	200 2	00 400 50	
									Eagle Rd	g. (ppm)	00
			1		1					Methane Elim.	

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. BORINGS BY CME-55 Low Clearance Drill **BH 8-23 DATE** August 14, 2023 **SAMPLE Photo Ionization Detector** PLOT **DEPTH** ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD STRATA NUMBER TYPE **Lower Explosive Limit %** N o v **GROUND SURFACE** 80 0+62.02Asphaltic concrete 0.10 1 FILL: Brown sifty sand, some gravel and crushed stone, trace brick and .07 1+61.022 25 7 wood SS 3 25 50 2+60.02**GLACIAL TILL:** Compact to very SS 4 25 58 dense, brown silty sand with gravel, 3+59.02cobbles and boulders SS 5 25 44 4+58.02 SS 6 50 50 +4.75 5+57.02RC 1 100 62 6+56.022 RC 100 100 7+55.02**BEDROCK:** Fair to excellent quality, grey limestone 8+54.02 RC 3 100 100 9+53.02RC 4 100 100 10+52.02 11 + 51.02RC 5 100 100 <u>12</u>.17 12 + 50.02End of Borehole 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic FILE NO. **PE2709 REMARKS** HOLE NO. BORINGS BY CME-55 Low Clearance Drill **BH 9-23 DATE** August 16, 2023 **SAMPLE Photo Ionization Detector** PLOT **DEPTH** ELEV. **SOIL DESCRIPTION** Volatile Organic Rdg. (ppm) (m) (m) RECOVERY VALUE r RQD STRATA NUMBER TYPE **Lower Explosive Limit %** N or v **GROUND SURFACE** 80 0+62.00Asphaltic concrete 0.08 ΑU 1 FILL: Brown silty sand, some grave 169 and crushed stone 1 + 61.00SS 2 67 8 SS 3 836 22 2+60.00SS 4 75 31 **GLACIAL TILL:** Compact to very dense, brown silty sand with gravel, 3+59.00SS 5 cobbles and boulders 67 46 4+58.00 SS 6 75 17 - some clay by 4.5m depth 7 SS 60 50+ 5+57.008 50+ 50 5.43 SS 1 RC 100 100 6+56.00RC 2 100 100 7 + 55.00**BEDROCK:** Excellent quality, grey limestone 8 + 54.00RC 3 100 100 9.07 9+53.00End of Borehole 200 300 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

9 Auriga Drive, Ottawa, Ontario K2E 7T9

SOIL PROFILE AND TEST DATA

Phase II - Environmental Site Assessment 137 & 141 George Street and 110 York Street Ottawa, Ontario

DATUM Geodetic									FILE NO.		
REMARKS									HOLE NO	D .	
BORINGS BY CME-55 Low Clearance I	Orill			D	ATE /	August 14	4, 2023	I	BH10-	-23	
SOIL DESCRIPTION	PLOT		SAN	/IPLE		DEPTH (m)	ELEV. (m)	1		Detecto c Rdg. (ppm	well Well
	STRATA	TYPE	NUMBER	% RECOVERY	N VALUE or RQD					ive Limit	Monitoring Well Construction
GROUND SURFACE				2	Z •	0-	62.04	20	40 6	60 80	2
							-61.04				
OVERBURDEN						2-	60.04				
OVERBURDEN							59.04				
							-58.04				
5.94						5-	-57.04				
End of Borehole									agle Rd	00 400 g. (ppm)	500

NORTHING:

SOIL PROFILE AND TEST DATA

Phase I - Environmental Site Assessment 116 York Street Ottawa, Ontario

9 Auriga Drive, Ottawa, Ontario K2E 7T9

368143.796

DATUM: Geodetic

EASTING:

REMARKS:

5032427.793 **ELEVATION**: 60.78

FILE NO.

PE6422

HOLE NO.

BORINGS BY: CME-55 Low Clearance [DATE:	May 8	2024	BH 1-24
SAMPLE DESCRIPTION	PLOT		SAN	/IPLE		DEPTH (m)	ELEV. (m)	Photo Ionization Detector Volatile Organic Rdg. (ppm)
GROUND SURFACE	STRATA	TYPE	NUMBER	% RECOVERY	N VALUE or RQD			Photo Ionization Detector Volatile Organic Rdg. (ppm) Current Lower Explosive Limit % 20 40 60 80
FILL: Compact brown silty sand with crushed concrete and stone, trace brick, metals, wood		ÃU	1			0-	-60.78	
		ss	2	21	10	1-	-59.78	
2.29		ss	3	42	4	2-	-58.78	
ILL: Stiff to very stiff brown silty lay with some sand, curshed stone nd concrete, trace cobbles, brick, rood, metals		ss	4	25	9	3-	-57.78	
3.81		ss	5	54	31			
SLACIAL TILL: Very dense grey andy silt with gravel, cobbles and oulders 4.42 EDROCK: Excellent quality grey	\^,^,^ \^,^,^ \ <u>^,</u> -	ss	6	29	+50	4-	-56.78	
mestone bedrock with interbedded hale seams		RC	1	100	97	5-	-55.78	
		_				6-	-54.78	
		RC	2	100	100	7-	-53.78	
		RC	3	100	100	8-	-52.78	
		_				9-	-51.78	
10.03 and of Borehole		RC 	4	100	100	10-	-50.78	
GWL @ 3.46m - May 22, 2024)								
								100 200 300 400 500 RKI Eagle Rdg. (ppm) ▲ Full Gas Resp. △ Methane Elim.

$paters on group {\tt Consulting Engineers}$

SOIL PROFILE AND TEST DATA

Phase I - Environmental Site Assessment 116 York Street Ottawa, Ontario

9 Auriga Drive, Ottawa, Ontario K2E 7T9

368156.266

NORTHING: Geodetic

5032440.634 **ELEVATION**: 60.38

FILE NO.

HOLE NO.

PE6422

REMARKS:

EASTING:

DATUM:

· · · · · · · · · · · · · · · · · · ·	Drill				DATE:	May 8	, 2024	1	BH 2-2	24
SAMPLE DESCRIPTION	PLOT		SAN	/IPLE		DEPTH		1	onization Detector Organic Rdg. (ppm)	TER
	STRATA PLOT	TYPE	NUMBER	% RECOVERY	N VALUE or RQD	(m)	(m)	O Lowe	r Explosive Limit %	PIEZOMETER
SPHALT 0.1				~	_	0-	60.38	20	40 60 80	-
SPHALT 0.1 ILL: Compact granualr with 0.2 rushed stone and gravel ILL: Compact brown silty sand ith some clay, trace graveland rushed stone		AU	1							
LACIAL TILL: Dense to very ense brown silty sand with clay,	7	SS	2	29	5	1-	-59.38	•		
ace gravel, occasional cobbles and oulders			6		4.5					
		ss	3	63	15	2-	-58.38	T		
		SS	4	42	16			•		
		ss	5		49	3-	-57.38	•		
3.9	(^^^^^ (^^^^^ (^^^^^)									
nd of Borehole										
								100	200 300 400 5	500

SOIL PROFILE AND TEST DATA

Phase I - Environmental Site Assessment 116 York Street Ottawa, Ontario

9 Auriga Drive, Ottawa, Ontario K2E 7T9

Geodetic

368157.789

NORTHING: 5032440.535 **ELEVATION**: FILE NO.

PE6422

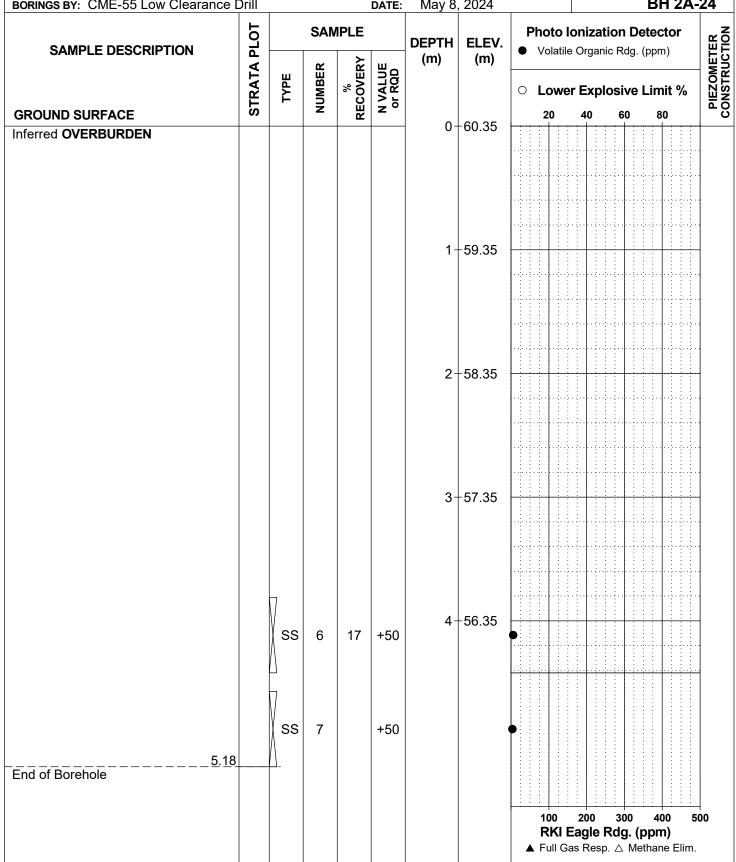
DATUM: **REMARKS:**

EASTING:

HOLE NO.

BH 2A-24 BORINGS BY: CME-55 Low Clearance Drill DATE: May 8, 2024

60.35



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NORTHING:

SOIL PROFILE AND TEST DATA

Phase I - Environmental Site Assessment 116 York Street Ottawa, Ontario

9 Auriga Drive, Ottawa, Ontario K2E 7T9

368178.7

EASTING:

DATUM: REMARKS:

Geodetic

5032404.654 **ELEVATION**: 60.93

FILE NO. PE6422

HOLE NO.

BORINGS BY: CME-55 Low Clearance	Drill	ı			DATE:	May 8	2024			E NO.	В	H 3-2	24
SAMPLE DESCRIPTION	PLOT			/IPLE		DEPTH (m)	ELEV. (m)	Photo Id					IG WELL
GROUND SURFACE	STRATA PLOT	TYPE	NUMBER	% RECOVERY	N VALUE or RQD	()	()	O Lower	Expl	losiv		nit %	MONITORING WEL
FILL: Brown silty sand with gravel and crushed stone, occasional		& AU	1			0-	-60.93	•					
concrete, brick and wood, trace clay		ss	2	17	2	1-	-59.93						
GLACIAL TILL: Dense to very		∆ V-ss	3	21	6		-58.93	•					
lense brown sandy silt with some gravel, occasional cobbles and coulders		∆ V ss	4	25	32	2-	-56.93	•					
Sandy pocket @ 2.29 m		∆ V ss	5	29	+50	3-	-57.93	•					
		RC	1	21	0	4-	-56.93						մարկանանանունական անանանանական անանանան անանանանան անանանան
5 11		ss	6		+50	5-	-55.93						
BEDROCK: Excellent quality grey mestone bedrock with interbedded hale seams	\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	⊔. RC	2	83	100	3-	-55.95						
		_	2		100	6-	-54.93						
		RC	3	100	100	7-	-53.93						
		_					50.00						
		RC	4	100	100	8-	-52.93						
		_				9-	-51.93						
10.26		RC	5	100	100	10-	-50.93						
ind of Borehole GWL @ 6.15m - May 22, 2024)	1 1 1 1	- -											1.1
C.1.2 & C.10111 May 22, 2027)								100 RKI E	200 agle	300 Rdg.			_ 500

SOIL PROFILE AND TEST DATA

Phase I - Environmental Site Assessment 116 York Street Ottawa, Ontario

9 Auriga Drive, Ottawa, Ontario K2E 7T9

368171.728 **EASTING:**

Geodetic

REMARKS:

DATUM:

5032397.855 **ELEVATION**: 61.33 NORTHING:

FILE NO. PE6422

HOLE NO.

BORINGS BY: CME-55 Low Clearance E	Orill				DATE:	May 9	, 2024		HOL	E NO		BH 4	1-24	1
SAMPLE DESCRIPTION	PLOT		SAN	/IPLE		DEPTH (m)	ELEV. (m)	Photo I					r	IG WELL
GROUND SURFACE	STRATA	TYPE	NUMBER	% RECOVERY	N VALUE or RQD	(,	(,	O Lowe	r Exp	olosi		imit '	%	MONITORING WEL
ASPHALT 0.03	$\times\!\!\times\!\!\times$	à AU	1			0-	61.33						: :	Í
FILL: Granular, light brown silty 0.20 sand with gravel and crushed stone FILL: Firm dark brown silty clay		AU I	ı											
with sand and gravel, some crush e ф2 stone, trace topsoil		ss	2	54	6	1-	-60.33	•						
GLACIAL TILL: Compact to very dense brown sandy silt, gravel, occasional cobbles and boulders		ss	3	38	18	2-	-59.33	•						
		ss	4	13	29			•						
		ss	5	50	12	3-	-58.33	•			- - -			
		∆ √ss	6	50	+50	4-	-57.33	•						
4.95		A ss	7	29	+50									
BEDROCK: Good to excellent quality grey limestone bedrock with nterbedded shale seams		RC	1	100	67	5-	-56.33							
norpeduca shale esame		_	'	100	07	6-	-55.33							
		RC	2	100	86	7-	-54.33							4
		_				8-	-53.33							
		RC	3	100	100	0-	-52.33							
		-		400		9-	52.33							
10.34		RC 	4	100	8	10-	-51.33							
End of Borehole														
(GWL @ 6.29m - May 22, 2024)								100 RKI I A Full G			j. (pj			00

$paters on group {\tt Consulting Engineers}$

SOIL PROFILE AND TEST DATA

Phase I - Environmental Site Assessment 116 York Street Ottawa, Ontario

9 Auriga Drive, Ottawa, Ontario K2E 7T9

368165.269 **EASTING:**

DATUM: Geodetic

REMARKS:

5032418.67 **ELEVATION**: 60.79 NORTHING:

FILE NO.

PE6422

HOLE NO.

BORINGS BY: CME-55 Low Clearance [Drill				DATE:	May 9	, 2024		BH 5-24	4
SAMPLE DESCRIPTION	PLOT		SAN	IPLE		DEPTH		1	onization Detector organic Rdg. (ppm)	TER
GROUND SURFACE	STRATAI	TYPE	NUMBER	% RECOVERY	N VALUE or RQD	(m)	(m)		er Explosive Limit %	PIEZOMETER
ASPHALT 0.05 FILL: Granular, brown silty sand 0.23 with gravel and crushed stone FILL: Firm brown silty sand with clay, gravel, crushed stone		AU	1	ш.		0-	-60.79	•		
GLACIAL TILL: Compact to dense brown silty sand with clay and gravel, occasional cobbles and boulders		-SS	2	54	5	1-	-59.79	•		
		ss	3	63	29	2-	-58.79	•		
		ss	4	38	+50	3-	-57.79			
3.76 End of Borehole		SS	5	54	20		,			
Practical refusal to augering @ 3.76 m										
									200 300 400 5 Eagle Rdg. (ppm) as Resp. △ Methane 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 Soft Firm Stiff Very Stiff Hard	<12 12-25 25-50 50-100 100-200 >200	<2 2-4 4-8 8-15 15-30 >30

SYMBOLS AND TERMS (continued)

SOIL DESCRIPTION (continued)

Cohesive soils can also be classified according to their "sensitivity". The sensitivity, S_t , 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:

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

Cc - Concavity coefficient = $(D30)^2 / (D10 \times D60)$

Cu - Uniformity coefficient = D60 / D10

Cc and Cu are used to assess the grading of sands and gravels:

Well-graded gravels have: 1 < Cc < 3 and Cu > 4 Well-graded sands have: 1 < Cc < 3 and Cu > 6

Sands 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 Ratio Initial sample void ratio = volume of voids / volume of solids

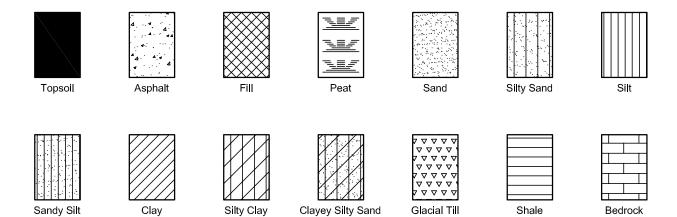
Wo - Initial water content (at start of consolidation test)

PERMEABILITY TEST

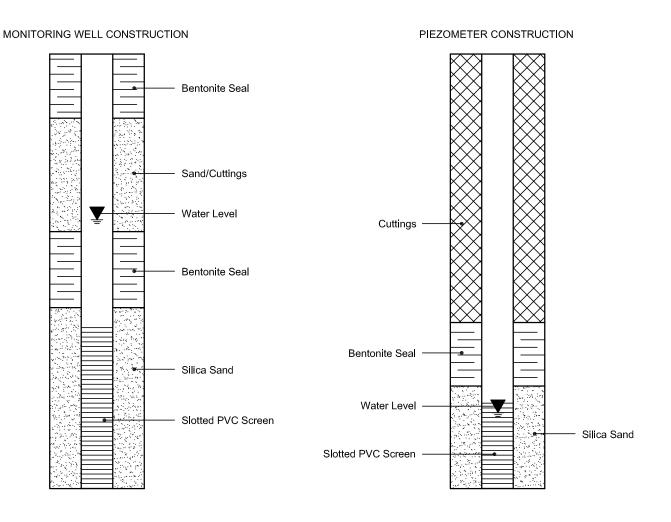
Coefficient of permeability or hydraulic conductivity is a measure of the ability of water to flow through the sample. The value of k is measured at a specified unit weight for (remoulded) cohesionless soil samples, because its value will vary with the unit weight or density of the sample during the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

SYMBOLS AND TERMS (continued)

STRATA PLOT



MONITORING WELL AND PIEZOMETER CONSTRUCTION





Head Office

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www.paracellabs.com

Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South

Nepean, ON K2E 7J5

Phone: (613) 226-7381

Fax: (613) 226-6344

Attn: Dan Arnott

 Client PO: 13160
 Report Date: 20-Jul-2012

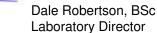
 Project: PE2709
 Order Date: 18-Jul-2012

 Custody: 94851
 Order #: 1229139

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

Paracel ID	Client II
1229139-01	BH1-SS7
1229139-02	BH2-SS8
1229139-03	BH3-SS6

Approved By:





Order #: 1229139

Certificate of Analysis

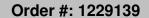
Client: Paterson Group Consulting Engineers

Client PO: 13160 Project Description: PE2709

Report Date: 20-Jul-2012 Order Date:18-Jul-2012

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date Analysis Date
CCME PHC F1	CWS Tier 1 - P&T GC-FID	18-Jul-12 20-Jul-12
CCME PHC F2 - F4	CWS Tier 1 - GC-FID, extraction	19-Jul-12 20-Jul-12
Solids, %	Gravimetric, calculation	19-Jul-12 19-Jul-12
VOCs	EPA 8260 - P&T GC-MS	18-Jul-12 20-Jul-12



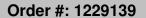


Client: Paterson Group Consulting Engineers

Client PO: 13160 Project Description: PE2709

Report Date: 20-Jul-2012 Order Date:18-Jul-2012

	Client ID: Sample Date:	BH1-SS7 17-Jul-12	BH2-SS8 17-Jul-12	BH3-SS6 17-Jul-12	- -
	Sample ID:	1229139-01	1229139-02	1229139-03	-
	MDL/Units	Soil	Soil	Soil	-
Physical Characteristics	0.4.0/ 1		1	T	
% Solids	0.1 % by Wt.	89.1	91.7	90.4	-
Volatiles	0 F0 ug/g dgy		T		
Acetone	0.50 ug/g dry	<0.50	<0.50	<0.50	-
Benzene	0.02 ug/g dry	<0.02	<0.02	<0.02	-
Bromodichloromethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Bromoform	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Bromomethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Carbon Tetrachloride	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Chlorobenzene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Chloroethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Chloroform	0.05 ug/g dry	< 0.05	<0.05	<0.05	-
Chloromethane	0.20 ug/g dry	<0.20	<0.20	<0.20	-
Dibromochloromethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Dichlorodifluoromethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,2-Dibromoethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,2-Dichlorobenzene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,3-Dichlorobenzene	0.05 ug/g dry	<0.05	< 0.05	<0.05	-
1,4-Dichlorobenzene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,1-Dichloroethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,2-Dichloroethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,1-Dichloroethylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,2-Dichloroethylene, total	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,2-Dichloropropane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,3-Dichloropropene, total	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Ethylbenzene	0.05 ug/g dry	<0.05	< 0.05	<0.05	-
Hexane	0.05 ug/g dry	<0.05	< 0.05	<0.05	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	<0.50	<0.50	<0.50	-
Methyl Butyl Ketone (2-Hexanone)	2.00 ug/g dry	<2.00	<2.00	<2.00	-
Methyl Isobutyl Ketone	0.50 ug/g dry	<0.50	<0.50	<0.50	-



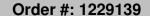


Client: Paterson Group Consulting Engineers

Client PO: 13160 Project Description: PE2709

Report Date: 20-Jul-2012 Order Date:18-Jul-2012

	Client ID:	BH1-SS7	BH2-SS8	BH3-SS6	-
	Sample Date: Sample ID:	17-Jul-12 1229139-01	17-Jul-12 1229139-02	17-Jul-12 1229139-03	-
	MDL/Units	Soil	Soil	Soil	-
Methyl tert-butyl ether	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Methylene Chloride	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Styrene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Tetrachloroethylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Toluene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,2,4-Trichlorobenzene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,1,1-Trichloroethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,1,2-Trichloroethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Trichloroethylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Trichlorofluoromethane	0.05 ug/g dry	<0.05	<0.05	<0.05	-
1,3,5-Trimethylbenzene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Vinyl chloride	0.02 ug/g dry	<0.02	<0.02	<0.02	-
m,p-Xylenes	0.05 ug/g dry	<0.05	<0.05	<0.05	-
o-Xylene	0.05 ug/g dry	<0.05	<0.05	<0.05	-
Xylenes, total	0.05 ug/g dry	<0.05	<0.05	<0.05	-
4-Bromofluorobenzene	Surrogate	102%	109%	115%	-
Dibromofluoromethane	Surrogate	92.6%	92.9%	91.6%	-
Toluene-d8	Surrogate	97.3%	100%	103%	-
Hydrocarbons			_		
F1 PHCs (C6-C10)	7 ug/g dry	182	20	<7	-
F2 PHCs (C10-C16)	4 ug/g dry	118	18	<4	-
F3 PHCs (C16-C34)	8 ug/g dry	<8	<8	<8	-
F4 PHCs (C34-C50)	6 ug/g dry	<6	<6	<6	-





Client: Paterson Group Consulting Engineers

Client PO: 13160 Project Description: PE2709

Report Date: 20-Jul-2012 Order Date:18-Jul-2012

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
•			0.1110	riodait	,				
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									
Acetone	ND	0.50	ug/g						
Benzene	ND	0.02	ug/g						
Bromodichloromethane	ND	0.05	ug/g						
Bromoform	ND	0.05	ug/g						
Bromomethane	ND	0.05	ug/g						
Carbon Tetrachloride	ND	0.05	ug/g						
Chlorobenzene	ND	0.05	ug/g						
Chloroethane	ND	0.05	ug/g						
Chloroform	ND	0.05	ug/g						
Chloromethane	ND	0.20	ug/g						
Dibromochloromethane	ND	0.05	ug/g						
Dichlorodifluoromethane	ND	0.05	ug/g						
1,2-Dibromoethane	ND	0.05	ug/g						
1,2-Dichlorobenzene	ND	0.05	ug/g						
1,3-Dichlorobenzene	ND	0.05	ug/g						
1,4-Dichlorobenzene	ND	0.05	ug/g						
1,1-Dichloroethane	ND	0.05	ug/g						
1,2-Dichloroethane	ND	0.05	ug/g						
1,1-Dichloroethylene	ND	0.05	ug/g						
cis-1,2-Dichloroethylene	ND	0.05	ug/g						
trans-1,2-Dichloroethylene	ND	0.05	ug/g						
1,2-Dichloroethylene, total 1,2-Dichloropropane	ND ND	0.05 0.05	ug/g						
cis-1,3-Dichloropropylene	ND ND	0.05	ug/g						
trans-1,3-Dichloropropylene	ND ND	0.05	ug/g						
1,3-Dichloropropene, total	ND ND	0.05	ug/g ug/g						
Ethylbenzene	ND	0.05	ug/g ug/g						
Hexane	ND	0.05	ug/g ug/g						
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g ug/g						
Methyl Butyl Ketone (2-Hexanone)	ND	2.00	ug/g						
Methyl Isobutyl Ketone	ND	0.50	ug/g						
Methyl tert-butyl ether	ND	0.05	ug/g						
Methylene Chloride	ND	0.05	ug/g						
Styrene	ND	0.05	ug/g						
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g						
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g						
Tetrachloroethylene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
1,2,4-Trichlorobenzene	ND	0.05	ug/g						
1,1,1-Trichloroethane	ND	0.05	ug/g						
1,1,2-Trichloroethane	ND	0.05	ug/g						
Trichloroethylene	ND	0.05	ug/g						
Trichlorofluoromethane	ND	0.05	ug/g						
1,3,5-Trimethylbenzene	ND	0.05	ug/g						
Vinyl chloride	ND	0.02	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g		4.4-	50 440			
Surrogate: 4-Bromofluorobenzene	9.36		ug/g		117	50-140			
Surrogate: Dibromofluoromethane	7.69		ug/g		96.1	50-140			
Surrogate: Toluene-d8	8.43		ug/g		105	50-140			



Client: Paterson Group Consulting Engineers

Client PO: 13160 Project Description: PE2709 Report Date: 20-Jul-2012 Order Date:18-Jul-2012

Amalista	F ,	Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g dry	ND				40	
F2 PHCs (C10-C16)	ND	4	ug/g dry	ND				30	
F3 PHCs (C16-C34)	37	8	ug/g dry ug/g dry	129			110.0	30	QR-01
F4 PHCs (C34-C50)	69	6	ug/g dry ug/g dry	136			65.1	30	QR-01
	09	O	ug/g ury	130			05.1	30	QI I-0 I
Physical Characteristics									
% Solids	87.3	0.1	% by Wt.	91.9			5.2	25	
Volatiles									
Acetone	ND	0.50	ua/a da	ND				ΕO	
	ND	0.50	ug/g dry	ND				50	
Benzene	ND	0.02	ug/g dry	ND				50	
Bromodichloromethane	ND	0.05	ug/g dry	ND				50	
Bromoform	ND	0.05	ug/g dry	ND				50	
Bromomethane	ND	0.05	ug/g dry	ND				50	
Carbon Tetrachloride	ND	0.05	ug/g dry	ND				50	
Chlorobenzene	ND	0.05	ug/g dry	ND				50	
Chloroethane	ND	0.05	ug/g dry	ND				50	
Chloroform	ND	0.05	ug/g dry	ND				50	
Chloromethane	ND	0.20	ug/g dry	ND				50	
Dibromochloromethane	ND	0.05	ug/g dry	ND				50	
Dichlorodifluoromethane	ND	0.05	ug/g dry	ND				50	
1,2-Dibromoethane	ND	0.05	ug/g dry	ND				50	
1,2-Dichlorobenzene	ND	0.05	ug/g dry	ND				50	
1.3-Dichlorobenzene	ND	0.05	ug/g dry	ND				50	
1,4-Dichlorobenzene	ND	0.05	ug/g dry	ND				50	
1,1-Dichloroethane	ND	0.05	ug/g dry	ND				50	
1,2-Dichloroethane	ND	0.05	ug/g dry	ND				50	
1,1-Dichloroethylene	ND	0.05	ug/g dry ug/g dry	ND				50	
cis-1,2-Dichloroethylene	ND	0.05		ND				50	
			ug/g dry						
trans-1,2-Dichloroethylene	ND	0.05	ug/g dry	ND				50 50	
1,2-Dichloropropane	ND	0.05	ug/g dry	ND				50	
cis-1,3-Dichloropropylene	ND	0.05	ug/g dry	ND				50	
trans-1,3-Dichloropropylene	ND	0.05	ug/g dry	ND				50	
Ethylbenzene	ND	0.05	ug/g dry	ND				50	
Hexane	ND	0.05	ug/g dry	ND				50	
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g dry	ND				50	
Methyl Butyl Ketone (2-Hexanone)	ND	2.00	ug/g dry	ND				50	
Methyl Isobutyl Ketone	ND	0.50	ug/g dry	ND				50	
Methyl tert-butyl ether	ND	0.05	ug/g dry	ND				50	
Methylene Chloride	ND	0.05	ug/g dry	ND				50	
Styrene	ND	0.05	ug/g dry	ND				50	
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g dry	ND				50	
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g dry	ND				50	
Tetrachloroethylene	ND	0.05	ug/g dry	ND				50	
Toluene	ND	0.05	ug/g dry	ND				50	
1,2,4-Trichlorobenzene	ND	0.05	ug/g dry	ND				50	
1,1,1-Trichloroethane	ND	0.05	ug/g dry	ND				50	
1,1,2-Trichloroethane	ND	0.05	ug/g dry	ND				50	
Trichloroethylene	ND	0.05	ug/g dry	ND				50	
Trichlorofluoromethane	ND	0.05	ug/g dry ug/g dry	ND				50	
1,3,5-Trimethylbenzene	ND ND	0.05		ND				50 50	
Vinyl chloride	ND ND	0.05	ug/g dry	ND				50 50	
,			ug/g dry						
m,p-Xylenes	ND	0.05	ug/g dry	ND				50	
o-Xylene	ND	0.05	ug/g dry	ND	440	50 440		50	
Surrogate: 4-Bromofluorobenzene	6.68		ug/g dry	ND	113	50-140			
Surrogate: Dibromofluoromethane	5.51		ug/g dry	ND	93.0	50-140			



Order #: 1229139

Certificate of Analysis

Client: Paterson Group Consulting Engineers

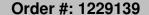
Client PO: 13160 Project Description: PE2709

Report Date: 20-Jul-2012 Order Date:18-Jul-2012

Method Quality Control: Duplicate

Analyte	Reporting Result Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Surrogate: Toluene-d8	5.93	ug/g dry	ND	100	50-140			

NIAGARA FALLS





Vinyl chloride

m,p-Xylenes

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 13160 Project Description: PE2709

Report Date: 20-Jul-2012 Order Date: 18-Jul-2012

Method Quality Control: Spike Reporting %REC **RPD** Source Result Units %REC **RPD** Notes Analyte Limit Limit Result Limit **Hydrocarbons** F1 PHCs (C6-C10) 218 7 ND 80-120 ug/g 109 F2 PHCs (C10-C16) 66 4 ug/g ND 82.5 80-120 8 F3 PHCs (C16-C34) 189 80-120 ug/g ND 94.5 F4 PHCs (C34-C50) 80-120 100 6 ug/g ND 83.6 Volatiles 0.50 ND 50-140 Acetone 11.7 ug/g 117 Benzene 4.10 0.02 ug/g ND 102 60-130 Bromodichloromethane 4.63 0.05 ND 116 60-130 ug/g 5.17 0.05 ND 129 60-130 Bromoform ug/g Bromomethane 3.46 0.05 ND 86.6 50-140 ug/g Carbon Tetrachloride 4.20 0.05 ND 105 60-130 ug/g Chlorobenzene 4.53 0.05 ug/g ND 113 60-130 Chloroethane 4.37 0.05 ND 109 50-140 ug/g Chloroform 4.60 0.05 ND 115 60-130 ug/g Chloromethane 2.70 0.20 ug/g ND 67.4 50-140 Dibromochloromethane 5.17 0.05 ug/g ND 129 60-130 Dichlorodifluoromethane 3.18 0.05 ug/g ND 79.6 50-140 1,2-Dibromoethane 4.94 0.05 ug/g ND 123 60-130 1,2-Dichlorobenzene 5.09 0.05 ug/g ND 127 60-130 1,3-Dichlorobenzene 4.75 0.05 ug/g ND 119 60-130 4.79 60-130 1,4-Dichlorobenzene 0.05 ug/g ND 120 1,1-Dichloroethane 4.51 0.05 ug/g ND 113 60-130 ND 109 60-130 1,2-Dichloroethane 4.37 0.05 ug/g 2.92 0.05 ND 73.1 60-130 1,1-Dichloroethylene ug/g cis-1,2-Dichloroethylene 4.47 0.05 ND 60-130 ug/g 112 86.3 trans-1,2-Dichloroethylene 3.45 0.05 ug/g ND 60-130 1,2-Dichloropropane 4.17 0.05 ug/g ND 104 60-130 0.05 ND 122 cis-1,3-Dichloropropylene 4.87 ug/g 60-130 trans-1,3-Dichloropropylene 0.05 ND 125 5.01 ug/g 60-130 Ethylbenzene 0.05 ND 97.4 60-130 3.90 ug/g Hexane 0.05 ND 110 60-130 4.38 ug/g Methyl Ethyl Ketone (2-Butanone) 10.8 0.50 ug/g ND 108 50-140 ND Methyl Butyl Ketone (2-Hexanone) 12.8 2.00 128 50-140 ug/g 0.50 ND Methyl Isobutyl Ketone 12.9 129 50-140 ug/g 0.05 ND 50-140 Methyl tert-butyl ether 12.1 121 ug/g Methylene Chloride 3.68 0.05 ug/g ND 92.0 60-130 Styrene 4.45 0.05 ug/g ND 111 60-130 4.66 0.05 ND 117 60-130 1,1,1,2-Tetrachloroethane ug/g 1,1,2,2-Tetrachloroethane 5.08 0.05 ND 127 60-130 ug/g Tetrachloroethylene 3.51 0.05 ND 87.7 60-130 ug/g Toluene 4.44 0.05 ND 111 60-130 ug/g 1,2,4-Trichlorobenzene 3.61 0.05 ug/g ND 90.3 60-130 1,1,1-Trichloroethane 3.95 0.05 ND 98.8 60-130 ug/g 1,1,2-Trichloroethane 5.15 0.05 ND 129 60-130 ug/g Trichloroethylene 3.49 0.05 ug/g ND 87.3 60-130 Trichlorofluoromethane 4.30 0.05 ug/g ND 107 50-140 1,3,5-Trimethylbenzene 3.53 0.05 ug/g ND 88.2 60-130

> P: 1-800-749-1947 E: paracel@paracellabs.com

3.16

0.02

0.05

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OTTAWA 300-2319 St. Laurent Blvd. Ottawa, ON K1G 4J8

ND

ND

ug/g

ug/g

MISSISSAUGA 6645 Kitimat Rd. Unit #27 Mississauga, ON L5N 6J3 NIAGARA FALLS 5415 Morning Glory Crt. Niagara Falls, ON L2J 0A3

50-140

60-130

SARNIA 123 Christina St. N. Sarnia, ON N7T 5T7

79.1

97.0



Order #: 1229139

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 13160 Project Description: PE2709

Report Date: 20-Jul-2012 Order Date: 18-Jul-2012

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
o-Xylene Surrogate: 4-Bromofluorobenzene	4.23 7.10	0.05	ug/g <i>ug/g</i>	ND	106 <i>88.8</i>	60-130 <i>50-140</i>			_



Order #: 1229139

Certificate of Analysis

Report Date: 20-Jul-2012 Client: Paterson Group Consulting Engineers Order Date:18-Jul-2012

Client PO: 13160 Project Description: PE2709

Qualifier Notes:

QC Qualifiers:

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

Sample Data Revisions

None

Work Order Revisions / Comments:

None

Other Report Notes:

n/a: not applicable

MDL: Method Detection Limit

Source Result: Data used as source for matrix and duplicate samples

%REC: Percent recovery.

RPD: Relative percent difference.

Soil results are reported on a dry weight basis when the units are denoted with 'dry'.

Where %Solids is reported, moisture loss includes the loss of volatile hydrocarbons.

CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.
- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.

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OTTAWA @ KINGSTON @ NIAGARA @ MISSIS	SAUGA	SA	RNIA		www.paracellabs.com								Pag	e of	f		
Client Name: Paterson Group Inc. Contact Name: Dan Arnott Address: 154 Colonnade Rd				Project Reference: PE 2709 Quote # PO# 12 / C						TAT:	TAT: Regular []3 Day						
Address: 154 Colonnade Rd Ottawa, ON KiE 75 Telephone: 613, 226, 7381		Email Address:	1	gra	p,	Ca					Date R	equired:			Ŧ		
Criteria: JO. Reg. 153/04 Table O. Reg. 153/11 (Curren	Table 3	RSC	Filing	1		4			B (St	orm) [JSUB	(Sanitary) Mu	micipality:] Other: _	
Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water)												Required .					
Paracel Order Number:	rix	Air Volume	of Containers	Sample	e Taken	FI-F4+BTEX	S		s by ICP/MS		WS)						
Sample ID/Location Name	Matrix	Air	Jo#	Date	Time	PHCs	VOCs	PAHs	Metals by	Hg	B (HWS)						
1 BHI-557	5		2	17-Jul-12		×	X										1
2 BH2-858	5					X	X										1
3 BH3-556	5		1	<u>V</u>		X	X										1
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Date/Time: [2-



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Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South

Nepean, ON K2E 7J5

Phone: (613) 226-7381

Fax: (613) 226-6344

Attn: Dan Arnott

 Client PO: 13161
 Report Date: 24-Jul-2012

 Project: PE2709
 Order Date: 19-Jul-2012

 Custody: 94801
 Order #: 1229217

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

Paracel ID Client ID 1229217-01 BH1-AU1 1229217-02 BH4-SS3

Approved By:

Mark Froto

Mark Foto, M.Sc. For Dale Robertson, BSc Laboratory Director



Order #: 1229217

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 13161 Project Description: PE2709

Report Date: 24-Jul-2012 Order Date:19-Jul-2012

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date Analysis	Date
Metals	EPA 6020 - Digestion - ICP-MS	20-Jul-12 21-	-Jul-12
PAHs by GC-MS, standard scan	EPA 8270 - GC-MS, extraction	20-Jul-12 24-	-Jul-12
Solids, %	Gravimetric, calculation	20-Jul-12 20-	-Jul-12



Report Date: 24-Jul-2012

Certificate of Analysis

Client: Paterson Group Consulting Engineers Order Date:19-Jul-2012

Client PO: 13161		Project Description	n: PE2709	.	uei Date.13-Jui-2012
	Client ID: Sample Date: Sample ID:	BH1-AU1 17-Jul-12 1229217-01	BH4-SS3 18-Jul-12 1229217-02	- - -	- - -
Dhariad Obarradariation	MDL/Units	Soil	Soil	-	-
Physical Characteristics	0.1 % by Wt.		1 000		
% Solids Metals	0.1 /8 by vvt.	98.6	89.6	-	-
Antimony	1 ug/g dry	<1	2		
Arsenic	1 ug/g dry	2	2		
Barium	1 ug/g dry	20	184		<u> </u>
Beryllium	0.5 ug/g dry	<0.5	<0.5		<u> </u>
Boron	5.0 ug/g dry	<5.0	<5.0		<u> </u>
Cadmium	0.5 ug/g dry	<0.5	0.5		
Chromium	5 ug/g dry	7	14		<u> </u>
Cobalt	1 ug/g dry	3	3	<u> </u>	<u> </u>
	5 ug/g dry	6	17	-	-
Copper Lead	1 ug/g dry	9	524	<u> </u>	-
	1 ug/g dry			<u>-</u>	-
Molybdenum Nickel	5 ug/g dry	2	1		-
Selenium	1 ug/g dry	11	8	<u>-</u>	-
Silver	0.3 ug/g dry	<1	<1		
Thallium	1 ug/g dry	1.8	1.5	-	-
Uranium	1 ug/g dry	<1	<1	-	-
	10 ug/g dry	<1	<1		-
Vanadium	20 ug/g dry	30	20	-	-
Zinc Semi-Volatiles	20 ug/g ury	<20	191	-	-
Acenaphthene	0.02 ug/g dry	<0.08 [1]			_
Acenaphthylene	0.02 ug/g dry	<0.08 [1]	-	<u> </u>	<u> </u>
Anthracene	0.02 ug/g dry	<0.08 [1]	-	-	
Benzo [a] anthracene	0.02 ug/g dry	0.09	-		<u>-</u>
Benzo [a] pyrene	0.02 ug/g dry		-		<u> </u>
Benzo [b] fluoranthene	0.02 ug/g dry	<0.08 [1]			-
Benzo [g,h,i] perylene	0.02 ug/g dry	<0.08 [1] <0.08 [1]	-		<u>-</u>
Benzo [k] fluoranthene	0.02 ug/g dry	<0.08 [1]	-	<u> </u>	
Biphenyl	0.02 ug/g dry	<0.08 [1]	-	<u> </u>	-
	0.02 ug/g dry	0.82	-		-
Chrysene Dibenzo [a,h] anthracene	0.02 ug/g dry			<u> </u>	
Fluoranthene	0.02 ug/g dry	<0.08 [1]	-	-	-
i iuoraninene	5.52 dg/g di j	<0.08 [1]	-	<u> </u>	-



Client PO: 13161

Order #: 1229217

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Project Description: PE2709

Report Date: 24-Jul-2012 Order Date:19-Jul-2012

	Client ID:	BH1-AU1	BH4-SS3	-	-
	Sample Date:	17-Jul-12	18-Jul-12	-	-
	Sample ID:	1229217-01	1229217-02	-	-
	MDL/Units	Soil	Soil	-	-
Fluorene	0.02 ug/g dry	<0.08 [1]	-	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	<0.08 [1]	-	•	-
1-Methylnaphthalene	0.02 ug/g dry	0.10	-	•	-
2-Methylnaphthalene	0.02 ug/g dry	0.13	-	-	-
Methylnaphthalene (1&2)	0.04 ug/g dry	0.23	-	-	-
Naphthalene	0.01 ug/g dry	0.05	-	•	-
Phenanthrene	0.02 ug/g dry	0.32	-	-	-
Pyrene	0.02 ug/g dry	0.22	-	-	-
2-Fluorobiphenyl	Surrogate	113%	-	-	-
Terphenyl-d14	Surrogate	108%	-	-	-



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 13161 Project Description: PE2709

Report Date: 24-Jul-2012 Order Date:19-Jul-2012

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Metals									
Antimony	ND	1	ug/g						
Arsenic	ND	1	ug/g						
Barium	ND	1	ug/g						
Beryllium	ND	0.5	ug/g						
Boron	ND	5.0	ug/g						
Cadmium	ND	0.5	ug/g						
Chromium	ND	5	ug/g						
Cobalt	ND	1	ug/g						
Copper	ND	5	ug/g						
Lead	ND	1	ug/g						
Molybdenum	ND	1	ug/g						
Nickel	ND	5	ug/g						
Selenium	ND	1	ug/g						
Silver	ND	0.3	ug/g						
Thallium	ND	1	ug/g						
Uranium	ND	1	ug/g						
Vanadium	ND	10	ug/g						
Zinc	ND	20	ug/g						



% Solids

Order #: 1229217

3.6

25

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 13161 Project Description: PE2709

84.8

0.1

Report Date: 24-Jul-2012 Order Date:19-Jul-2012

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Metals									
Antimony	ND	1	ug/g dry	ND			0.0	30	
Arsenic	ND	1	ug/g dry	ND			0.0	30	
Barium	14.2	1	ug/g dry	13.9			2.8	30	
Beryllium	ND	0.5	ug/g dry	ND			0.0	30	
Boron	ND	5.0	ug/g dry	ND			0.0	30	
Cadmium	ND	0.5	ug/g dry	ND			0.0	30	
Chromium	10.2	5	ug/g dry	10.0			2.3	30	
Cobalt	2.4	1	ug/g dry	2.3			2.8	30	
Copper	ND	5	ug/g dry	ND			0.0	30	
_ead	ND	1	ug/g dry	1.8			0.0	30	
Molybdenum	1.2	1	ug/g dry	ND			0.0	30	
Nickel	6.3	5	ug/g dry	6.1			4.0	30	
Selenium	ND	1	ug/g dry	ND			0.0	30	
Silver	ND	0.3	ug/g dry	ND			0.0	30	
Thallium	ND	1	ug/g dry	ND			0.0	30	
Jranium	ND	1	ug/g dry	ND			0.0	30	
Vanadium	14.8	10	ug/g dry	13.8			6.7	30	
Zinc	ND	20	ug/g dry	ND			0.0	30	
Physical Characteristics									
	010	• •							

% by Wt.

88.0



Vanadium

Zinc

Order #: 1229217

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 13161 Project Description: PE2709

57.7

53.2

Report Date: 24-Jul-2012 Order Date:19-Jul-2012

Method Quality Co	ntrol: Spike							
Analyte	Reportin Result Limit	g Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Metals								
Antimony	51.0	ug/L	0.01	102	70-130			
Arsenic	49.5	ug/L	0.1	98.7	70-130			
Barium	55.5	ug/L	5.5	100	70-130			
Beryllium	52.8	ug/L	0.08	105	70-130			
Boron	48.3	ug/L	0.5	95.6	70-130			
Cadmium	47.8	ug/L	0.007	95.5	70-130			
Chromium	55.4	ug/L	4.0	103	70-130			
Cobalt	50.9	ug/L	0.9	99.9	70-130			
Copper	50.9	ug/L	1.5	98.8	70-130			
Lead	51.4	ug/L	0.7	101	70-130			
Molybdenum	48.1	ug/L	0.07	96.1	70-130			
Nickel	52.8	ug/L	2.4	101	70-130			
Selenium	50.2	ug/L	ND	100	70-130			
Silver	41.4	ug/L	0.01	82.7	70-130			
Thallium	56.7	ug/L	ND	114	70-130			
Uranium	51.7	ug/L	0.1	103	70-130			

ug/L

ug/L

104

98.0

5.5 4.2 70-130

70-130



Certificate of Analysis

Client: Paterson Group Consulting Engineers Order Date:19-Jul-2012

Client PO: 13161 Project Description: PE2709

Qualifier Notes:

Sample Qualifiers:

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

Sample Data Revisions

None

Work Order Revisions / Comments:

None

Other Report Notes:

n/a: not applicable

MDL: Method Detection Limit

Source Result: Data used as source for matrix and duplicate samples

%REC: Percent recovery.

RPD: Relative percent difference.

Soil results are reported on a dry weight basis when the units are denoted with 'dry'. Where %Solids is reported, moisture loss includes the loss of volatile hydrocarbons. Report Date: 24-Jul-2012

	PARACEL LABORATORIES LTD. VA ® KINGSTON ® NIAGARA ® MISSISS.	RI	ELIAE	BLE.						300- Otta p: 1- e: pa	wa, C 800-7 arace	St. La Intario '49-19 @par	urent Blvd K1G 4J8 947 acellabs.cc		1	Λ. (Γ	of Cu ab Use On 948	(b) 1	
Client N Contact Address Telepho	Name: Paterson Group Inc. Name: Dan Arnott 154 Colonnade Rel, Ottawa,	0N 1	K267	J5	Project Reference Quote # PO # 131/2 Email Address: Jarn at 1 O. Reg. 558/00	bl @ pate	rsor	ng	500	φ.	Са				Date Re	Regula	ar []	3 Day 1 Day	
Matrix 7	Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) SS	(Storm/S	Sanitary S	ewer) P	(Paint) A (Air) O (Other)							Requir	ed An	alyses				
Parace	1 Order Number:	Matrix	Air Volume	of Containers	Sample	e Taken	Cs FI-F4+BTEX	Cs	Hs	Metals by ICP/MS		CrVI					8.1		
	Sample ID/Location Name	-	Ā	#	Date	Time	PHCs	VOCs	PAHS	Me	Hga	CrVI							
2	BH - MV	5			17-Jul-12		-		X	X		-		25	50 n	2			/
3	BH4-553	5		1	[8]VL12		\vdash		\vdash	X	Н	+	-	_	11				_/
4	_			_	-0.4.1.3		-		\vdash	_	\vdash	+	-	_					_
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Received at Lab:

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Temperature: 18 19 °C

Verified By:

Date/Time:

MC

pH Verified [| By: _

Received by Driver/Depot:

Temperature: _

Relinquished By (Print & Sign).

Date/Time: 1 22

ARNOTT



Head Office

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Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South

Nepean, ON K2E 7J5

Phone: (613) 226-7381

Fax: (613) 226-6344

Attn: Dan Arnott

 Client PO: 12278
 Report Date: 9-Aug-2012

 Project: PE2709
 Order Date: 8-Aug-2012

 Custody: 5106
 Order #: 1232102

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

Paracel ID	Client II
1232102-01	BH5-SS8
1232102-02	BH6-SS7
1232102-03	BH7-SS7

Approved By:

Mark Froto

Mark Foto, M.Sc. For Dale Robertson, BSc Laboratory Director



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 12278 Project Description: PE2709

Report Date: 09-Aug-2012 Order Date:8-Aug-2012

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date Analysis Date
BTEX	EPA 8260 - P&T GC-MS	8-Aug-12 9-Aug-12
CCME PHC F1	CWS Tier 1 - P&T GC-FID	8-Aug-12 9-Aug-12
CCME PHC F2 - F4	CWS Tier 1 - GC-FID, extraction	8-Aug-12 9-Aug-12
Solids, %	Gravimetric, calculation	9-Aug-12 9-Aug-12



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Report Date: 09-Aug-2012 Order Date:8-Aug-2012

Client PO: 12278 Project Description: PE2709

	Client ID:	BH5-SS8	BH6-SS7	BH7-SS7	-
	Sample Date:	08-Aug-12	08-Aug-12	08-Aug-12	-
	Sample ID:	1232102-01	1232102-02	1232102-03	-
	MDL/Units	Soil	Soil	Soil	-
Physical Characteristics					
% Solids	0.1 % by Wt.	74.0	96.4	96.5	-
Volatiles	-		-	-	
Benzene	0.02 ug/g dry	<0.02	<0.02	<0.02	-
Ethylbenzene	0.05 ug/g dry	< 0.05	<0.05	<0.05	-
Toluene	0.05 ug/g dry	< 0.05	<0.05	<0.05	-
m,p-Xylenes	0.05 ug/g dry	< 0.05	<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	90.2%	92.8%	93.3%	-
Hydrocarbons					
F1 PHCs (C6-C10)	7 ug/g dry	65	8	<7	-
F2 PHCs (C10-C16)	4 ug/g dry	74	<4	<4	-
F3 PHCs (C16-C34)	8 ug/g dry	<8	<8	62	-
F4 PHCs (C34-C50)	6 ug/g dry	<6	<6	112	-



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 12278 Project Description: PE2709 Report Date: 09-Aug-2012

Order Date:8-Aug-2012

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						
n,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g						
Surrogate: Toluene-d8	3.22		ug/g		101	50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 12278 Project Description: PE2709 Report Date: 09-Aug-2012

Order Date:8-Aug-2012

Method Quality Control: Duplicate Reporting Source %REC RPD Analyte Result Ĺimit RPD Units Result %REC Limit Limit Notes **Physical Characteristics** % Solids 68.7 0.1 % by Wt. 66.6 3.1 25



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 12278 Project Description: PE2709

Report Date: 09-Aug-2012 Order Date:8-Aug-2012

Method Quality Control: Spike									
Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	205	7	ug/g	ND	103	80-120			
Volatiles									
Benzene	3.89	0.02	ug/g	ND	97.1	60-130			
Ethylbenzene	2.78	0.05	ug/g	ND	69.4	60-130			
Toluene	3.99	0.05	ug/g	ND	99.7	60-130			
m,p-Xylenes	6.54	0.05	ug/g	ND	81.8	60-130			
o-Xylene	2.69	0.05	ug/g	ND	67.2	60-130			
Surrogate: Toluene-d8	2.72		ug/g		85.1	50-140			



Certificate of Analysis

Report Date: 09-Aug-2012 Client: Paterson Group Consulting Engineers Order Date:8-Aug-2012

Client PO: 12278 Project Description: PE2709

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.

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.



8-AU1,-12

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°C

Date/Time:

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Chain of Custody

(Lab Use Only)

e: paracel@paracellabs.com OTTAWA @ KINGSTON @ NIAGARA @ MISSISSAUGA @ SARNIA www.paracellabs.com Page of Project Reference: TAT: [] Regular []3 Day Contact Name: Quote #) Day Address: |5 []2 Day Email Address: Date Required: Telephone: Criteria: | | O. Reg. 153/04 Table ____ | O. Reg. 153/11 (Current) Table __ | | RSC Filing | | O. Reg. 558/00 | | PWQO | | CCME | | SUB (Storm) | | SUB (Sanitary) Municipality: [] Other: Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) SS (Storm/Sanitary Sewer) P (Paint) A (Air) O (Other) Required Analyses Paracel Order Number: Containers Air Volume Sample Taken Matrix JO # Sample ID/Location Name Date Time 1 8-Ax-12 190 m2 x 1 DAM 2 PM 3 2pm 4 5 6 7 8 9 10 should read SS8 per Day Comments: Method of Delivery Received by Driver/Depot: Received at Le Date/Time: Date/Time: Date/Time:

Chain of Custody (Blank) - Rev 0.0 December 2011

Temperature: 48.3 0

pH Verified [] By:



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

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive Ottawa, ON K2E 7T9 Attn: Karyn Munch

Client PO: 56968 Project: PE2709

Custody:

Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

Order #: 2310245

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

Paracel ID	Client ID
2310245-01	BH1-23-SS3
2310245-02	BH2-23-SS2
2310245-03	BH2-23-SS4
2310245-04	BH3-23-AU2
2310245-05	BH3-23-SS4
2310245-06	BH4-23-SS4
2310245-07	BH5-23-AU1
2310245-08	BH5-23-SS3
2310245-09	BH5-23-SS6
2310245-10	BH6-23-SS5
2310245-11	DUP
2310245-12	BH1-23-SS2
2310245-13	BH1-23-SS5
2310245-14	BH3-23-SS5
2310245-15	BH3-23-SS6

Approved By:

Mark Foto

Mark Foto, M.Sc. Lab Supervisor



Order #: 2310245

Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 8-Mar-2023

 Client PO:
 56968
 Project Description: PE2709

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	7-Mar-23	7-Mar-23
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	9-Mar-23	10-Mar-23
Conductivity	MOE E3138 - probe @25 °C, water ext	9-Mar-23	9-Mar-23
Mercury by CVAA	EPA 7471B - CVAA, digestion	10-Mar-23	10-Mar-23
pH, soil	EPA 150.1 - pH probe @ 25 °C, CaCl buffered ext.	10-Mar-23	10-Mar-23
PHC F1	CWS Tier 1 - P&T GC-FID	7-Mar-23	7-Mar-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	8-Mar-23	10-Mar-23
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	9-Mar-23	9-Mar-23
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	8-Mar-23	11-Mar-23
REG 153: VOCs by P&T GC/MS	EPA 8260 - P&T GC-MS	9-Mar-23	9-Mar-23
SAR	Calculated	9-Mar-23	10-Mar-23
Solids, %	CWS Tier 1 - Gravimetric	9-Mar-23	9-Mar-23



Report Date: 13-Mar-2023

Order Date: 8-Mar-2023 **Project Description: PE2709**

Client: Paterson Group Consulting Engineers

Client PO: 56968

Certificate of Analysis

BH2-23-SS2 Client ID: BH1-23-SS3 BH2-23-SS4 BH3-23-AU2 Sample Date: 24-Feb-23 09:00 27-Feb-23 09:00 27-Feb-23 09:00 27-Feb-23 09:00 2310245-01 2310245-02 2310245-03 2310245-04 Sample ID: MDL/Units Soil Soil Soil Soil **Physical Characteristics** 0.1 % by Wt. % Solids 93.2 88.4 92.7 90.6 General Inorganics 0.01 N/A SAR 0.80 1.61 1.77 2.16 5 uS/cm Conductivity 1360 693 442 469 0.05 pH Units рΗ 7.69 7.76 7.66 Metals 1.0 ug/g dry Antimony <1.0 2.2 <1.0 2.2 1.0 ug/g dry Arsenic 2.6 3.8 8.0 6.8 1.0 ug/g dry Barium 30.7 321 27.1 126 Beryllium 0.5 ug/g dry <0.5 0.6 <0.5 0.6 5.0 ug/g dry Boron 6.4 7.0 6.5 <5.0 0.5 ug/g dry Cadmium < 0.5 < 0.5 < 0.5 0.6 5.0 ug/g dry Chromium 12.2 29.2 10.1 16.7 0.2 ug/g dry Chromium (VI) < 0.2 < 0.2 <0.2 <0.2 1.0 ug/g dry Cobalt 3.8 7.1 3.6 5.7 5.0 ug/g dry Copper 9.3 120 5.7 36.6 1.0 ug/g dry Lead 3.6 432 32 217 0.1 ug/g dry Mercury < 0.1 2.7 < 0.1 0.3 1.0 ug/g dry Molybdenum <1.0 1.1 <1.0 1.3 5.0 ug/g dry Nickel 8.2 17.0 7.1 14.3 1.0 ug/g dry Selenium <1.0 1.3 <1.0 <1.0 0.3 ug/g dry Silver < 0.3 0.6 < 0.3 < 0.3 1.0 ug/g dry Thallium <1.0 <1.0 <1.0 <1.0 1.0 ug/g dry Uranium <1.0 <1.0 <1.0 <1.0 Vanadium 10.0 ug/g dry 15.7 17.5 26.1 22.1 Zinc 20.0 ug/g dry 429 <20.0 74.8 21.8 Volatiles 0.02 ug/g dry Benzene < 0.02 < 0.02 < 0.02 < 0.02 0.05 ug/g dry Ethylbenzene < 0.05 < 0.05 < 0.05 < 0.05 0.05 ug/g dry Toluene < 0.05 < 0.05 < 0.05 < 0.05 0.05 ug/g dry m,p-Xylenes < 0.05 < 0.05 < 0.05 < 0.05 0.05 ug/g dry o-Xylene < 0.05 < 0.05 < 0.05 < 0.05 Xylenes, total 0.05 ug/g dry < 0.05 < 0.05 < 0.05 < 0.05 Toluene-d8 Surrogate 123% 127% 124% 118% Hydrocarbons



Order #: 2310245

lanart Data: 12 Mar 20

Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

Client: Paterson Group Consulting Engineers
Client PO: 56968

Project Description: PE2709

	Client ID: Sample Date: Sample ID: MDL/Units	BH1-23-SS3 24-Feb-23 09:00 2310245-01 Soil	BH2-23-SS2 27-Feb-23 09:00 2310245-02 Soil	BH2-23-SS4 27-Feb-23 09:00 2310245-03 Soil	BH3-23-AU2 27-Feb-23 09:00 2310245-04 Soil
F1 PHCs (C6-C10)	7 ug/g dry	<7	<7	<7	<7
F2 PHCs (C10-C16)	4 ug/g dry	<4	<40 [1]	<4	<4
F3 PHCs (C16-C34)	8 ug/g dry	<8	312	<8	48 [2]
F4 PHCs (C34-C50)	6 ug/g dry	<6	298	<6	23 [2]
Semi-Volatiles	-				
Acenaphthene	0.02 ug/g dry	<0.02	0.76	-	0.07
Acenaphthylene	0.02 ug/g dry	<0.02	0.19	-	0.18
Anthracene	0.02 ug/g dry	<0.02	1.58	-	0.29
Benzo [a] anthracene	0.02 ug/g dry	<0.02	2.72	-	0.70
Benzo [a] pyrene	0.02 ug/g dry	<0.02	2.37	-	0.70
Benzo [b] fluoranthene	0.02 ug/g dry	<0.02	3.00	-	0.73
Benzo [g,h,i] perylene	0.02 ug/g dry	<0.02	1.45	-	0.38
Benzo [k] fluoranthene	0.02 ug/g dry	<0.02	1.68	-	0.42
Chrysene	0.02 ug/g dry	<0.02	2.68	-	0.71
Dibenzo [a,h] anthracene	0.02 ug/g dry	<0.02	0.38	-	0.11
Fluoranthene	0.02 ug/g dry	<0.02	6.09	-	1.47
Fluorene	0.02 ug/g dry	<0.02	0.53	-	0.08
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	<0.02	1.40	-	0.37
1-Methylnaphthalene	0.02 ug/g dry	<0.02	0.12	-	0.09
2-Methylnaphthalene	0.02 ug/g dry	<0.02	0.16	-	0.12
Methylnaphthalene (1&2)	0.04 ug/g dry	<0.04	0.28	-	0.21
Naphthalene	0.01 ug/g dry	<0.01	0.22	-	0.16
Phenanthrene	0.02 ug/g dry	<0.02	4.82	-	0.89
Pyrene	0.02 ug/g dry	<0.02	4.84	-	1.28
2-Fluorobiphenyl	Surrogate	106%	101%	-	108%
Terphenyl-d14	Surrogate	115%	110%	-	108%



Order #: 2310245

Report Date: 13-Mar-2023

Order Date: 8-Mar-2023

Client: Paterson Group Consulting Engineers Client PO: 56968 **Project Description: PE2709**

	Client ID: Sample Date: Sample ID: MDL/Units	BH3-23-SS4 27-Feb-23 09:00 2310245-05 Soil	BH4-23-SS4 28-Feb-23 09:00 2310245-06 Soil	BH5-23-AU1 28-Feb-23 09:00 2310245-07 Soil	BH5-23-SS3 28-Feb-23 09:00 2310245-08 Soil
Physical Characteristics	MIDE/OTHES			Con	Gon
% Solids	0.1 % by Wt.	96.2	92.5	92.6	98.4
General Inorganics	+		!	!	
SAR	0.01 N/A	1.19	4.44	0.90	1.32
Conductivity	5 uS/cm	209	1090	280	293
рН	0.05 pH Units	-	-	-	7.89
Metals					
Antimony	1.0 ug/g dry	<1.0	<1.0	<1.0	<1.0
Arsenic	1.0 ug/g dry	4.1	3.5	5.9	2.8
Barium	1.0 ug/g dry	38.6	27.1	24.1	38.1
Beryllium	0.5 ug/g dry	<0.5	<0.5	<0.5	<0.5
Boron	5.0 ug/g dry	5.6	7.2	6.8	5.6
Cadmium	0.5 ug/g dry	<0.5	<0.5	<0.5	<0.5
Chromium	5.0 ug/g dry	11.3	12.5	15.9	10.5
Chromium (VI)	0.2 ug/g dry	<0.2	<0.2	<0.2	<0.2
Cobalt	1.0 ug/g dry	4.2	3.5	6.8	3.7
Copper	5.0 ug/g dry	8.7	7.0	12.4	5.9
Lead	1.0 ug/g dry	5.3	4.5	32.5	4.5
Mercury	0.1 ug/g dry	<0.1	<0.1	<0.1	<0.1
Molybdenum	1.0 ug/g dry	<1.0	<1.0	6.7	<1.0
Nickel	5.0 ug/g dry	8.6	7.7	15.6	6.8
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	<1.0	<1.0	<1.0	<1.0
Vanadium	10.0 ug/g dry	19.2	19.5	25.8	14.3
Zinc	20.0 ug/g dry	20.5	23.8	26.3	<20.0
Volatiles					
Acetone	0.50 ug/g dry	<0.50	-	-	-
Benzene	0.02 ug/g dry	<0.02	-	-	-
Bromodichloromethane	0.05 ug/g dry	<0.05	-	-	-
Bromoform	0.05 ug/g dry	<0.05	-	-	-
Bromomethane	0.05 ug/g dry	<0.05	-	-	-
Carbon Tetrachloride	0.05 ug/g dry	<0.05	-	-	-
Chlorobenzene	0.05 ug/g dry	<0.05	-	-	-
Chloroform	0.05 ug/g dry	<0.05	-	-	-



Client: Paterson Group Consulting Engineers

Certificate of Analysis

Order #: 2310245

Report Date: 13-Mar-2023

Order Date: 8-Mar-2023

Client PO: 56968 Project Description: PE2709

	Client ID: Sample Date:	BH3-23-SS4 27-Feb-23 09:00 2310245-05	BH4-23-SS4 28-Feb-23 09:00 2310245-06	BH5-23-AU1 28-Feb-23 09:00 2310245-07	BH5-23-SS3 28-Feb-23 09:00 2310245-08
	Sample ID: MDL/Units	23 10245-05 Soil	2310245-06 Soil	2310245-07 Soil	23 10245-06 Soil
Dibromochloromethane	0.05 ug/g dry	<0.05	-	-	-
Dichlorodifluoromethane	0.05 ug/g dry	<0.05	-	-	-
1,2-Dichlorobenzene	0.05 ug/g dry	<0.05	-	-	-
1,3-Dichlorobenzene	0.05 ug/g dry	<0.05	-	-	-
1,4-Dichlorobenzene	0.05 ug/g dry	<0.05	-	-	-
1,1-Dichloroethane	0.05 ug/g dry	<0.05	-	-	-
1,2-Dichloroethane	0.05 ug/g dry	<0.05	-	-	-
1,1-Dichloroethylene	0.05 ug/g dry	<0.05	-	-	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	-	-	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	-	-	-
1,2-Dichloropropane	0.05 ug/g dry	<0.05	-	-	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	-	-	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	-	-	-
1,3-Dichloropropene, total	0.05 ug/g dry	<0.05	-	-	-
Ethylbenzene	0.05 ug/g dry	<0.05	-	-	-
Ethylene dibromide (dibromoethane, 1	0.05 ug/g dry	<0.05	-	-	-
Hexane	0.05 ug/g dry	<0.05	-	-	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	<0.50	-	-	-
Methyl Isobutyl Ketone	0.50 ug/g dry	<0.50	-	-	-
Methyl tert-butyl ether	0.05 ug/g dry	<0.05	-	-	-
Methylene Chloride	0.05 ug/g dry	<0.05	-	-	-
Styrene	0.05 ug/g dry	<0.05	-	-	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	<0.05	-	-	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	<0.05	-	-	-
Tetrachloroethylene	0.05 ug/g dry	<0.05	-	-	-
Toluene	0.05 ug/g dry	<0.05	-	-	-
1,1,1-Trichloroethane	0.05 ug/g dry	<0.05	-	-	-
1,1,2-Trichloroethane	0.05 ug/g dry	<0.05	-	-	-
Trichloroethylene	0.05 ug/g dry	<0.05	-	-	-
Trichlorofluoromethane	0.05 ug/g dry	<0.05	-	-	-
Vinyl chloride	0.02 ug/g dry	<0.02	-	-	-
m,p-Xylenes	0.05 ug/g dry	<0.05	-	-	-
o-Xylene	0.05 ug/g dry	<0.05	-	-	-
Xylenes, total	0.05 ug/g dry	<0.05	-	-	-
4-Bromofluorobenzene	Surrogate	98.1%	-	-	-



Order #: 2310245

Report Date: 13-Mar-2023

Order Date: 8-Mar-2023

Client: Paterson Group Consulting Engineers
Client PO: 56968

Project Description: PE2709

	Client ID:	BH3-23-SS4	BH4-23-SS4	BH5-23-AU1	BH5-23-SS3
	Sample Date:	27-Feb-23 09:00	28-Feb-23 09:00 2310245-06	28-Feb-23 09:00 2310245-07	28-Feb-23 09:00
	Sample ID:	2310245-05 Soil	2310245-06 Soil	2310245-07 Soil	2310245-08 Soil
Dibromofluoromethane	Surrogate	104%	-	-	-
Toluene-d8	Surrogate	121%	-	-	-
Benzene	0.02 ug/g dry	-	<0.02	<0.02	<0.02
Ethylbenzene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
Toluene	0.05 ug/g dry	-	<0.05	<0.05	<0.05
m,p-Xylenes	0.05 ug/g dry	-	<0.05	<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	-	126%	126%	120%
Hydrocarbons					
F1 PHCs (C6-C10)	7 ug/g dry	<7	<7	<7	<7
F2 PHCs (C10-C16)	4 ug/g dry	<4	<4	<80 [1]	<4
F3 PHCs (C16-C34)	8 ug/g dry	<8	<8	699	<8
F4 PHCs (C34-C50)	6 ug/g dry	<6	<6	1650	<6
Semi-Volatiles	1 1		T	Г	1
Acenaphthene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Acenaphthylene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Anthracene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Benzo [a] anthracene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Benzo [a] pyrene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Benzo [b] fluoranthene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Benzo [g,h,i] perylene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Benzo [k] fluoranthene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Chrysene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Dibenzo [a,h] anthracene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Fluoranthene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Fluorene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
1-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
2-Methylnaphthalene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Methylnaphthalene (1&2)	0.04 ug/g dry	<0.04	<0.04	<0.80 [1]	<0.04
Naphthalene	0.01 ug/g dry	<0.01	<0.01	<0.20 [1]	<0.01
Phenanthrene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
Pyrene	0.02 ug/g dry	<0.02	<0.02	<0.40 [1]	<0.02
2-Fluorobiphenyl	Surrogate	104%	96.1%	69.4%	95.0%



Order #: 2310245

Report Date: 13-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 8-Mar-2023

 Client PO:
 56968
 Project Description: PE2709

	Client ID:	BH3-23-SS4	BH4-23-SS4	BH5-23-AU1	BH5-23-SS3
	Sample Date:		28-Feb-23 09:00 28-Feb-23 09:00		28-Feb-23 09:00
	Sample ID:	2310245-05	2310245-06	2310245-07	2310245-08
	MDL/Units	Soil	Soil	Soil	Soil
Terphenyl-d14	Surrogate	109%	111%	101%	111%



Order #: 2310245

Report Date: 13-Mar-2023

Order Date: 8-Mar-2023

Client: Paterson Group Consulting Engineers
Client PO: 56968

Project Description: PE2709

	Client ID: Sample Date:	BH5-23-SS6 28-Feb-23 09:00	BH6-23-SS5 01-Mar-23 09:00	DUP 28-Feb-23 09:00	BH1-23-SS2 24-Feb-23 09:00
	Sample ID:	2310245-09 Soil	2310245-10 Soil	2310245-11 Soil	2310245-12 Soil
Physical Characteristics	MDL/Units	3011	3011	3011	3011
% Solids	0.1 % by Wt.	94.2	94.7	96.0	92.4
General Inorganics	- 	V	<u> </u>	1 00.0	02
SAR	0.01 N/A	1.25	1.88	1.43	1.90
Conductivity	5 uS/cm	226	3310	304	943
Metals	· · · · · · · · · · · · · · · · · · ·		•		•
Antimony	1.0 ug/g dry	<1.0	<1.0	<1.0	-
Arsenic	1.0 ug/g dry	4.1	10.2	4.6	-
Barium	1.0 ug/g dry	129	64.7	211	-
Beryllium	0.5 ug/g dry	<0.5	<0.5	0.5	-
Boron	5.0 ug/g dry	8.0	11.6	8.9	-
Cadmium	0.5 ug/g dry	<0.5	<0.5	<0.5	-
Chromium	5.0 ug/g dry	13.5	13.6	15.5	-
Chromium (VI)	0.2 ug/g dry	<0.2	<0.2	<0.2	-
Cobalt	1.0 ug/g dry	4.5	11.8	4.9	-
Copper	5.0 ug/g dry	<5.0	9.6	5.5	-
Lead	1.0 ug/g dry	5.0	19.7	5.1	-
Mercury	0.1 ug/g dry	<0.1	<0.1	<0.1	-
Molybdenum	1.0 ug/g dry	<1.0	5.4	<1.0	-
Nickel	5.0 ug/g dry	11.4	20.3	12.5	-
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	<1.0	<1.0	<1.0	-
Vanadium	10.0 ug/g dry	15.4	10.5	18.0	-
Zinc	20.0 ug/g dry	<20.0	<20.0	<20.0	-
Volatiles					
Acetone	0.50 ug/g dry	<0.50	-	<0.50	-
Benzene	0.02 ug/g dry	<0.02	-	<0.02	-
Bromodichloromethane	0.05 ug/g dry	<0.05	-	<0.05	-
Bromoform	0.05 ug/g dry	<0.05	-	<0.05	-
Bromomethane	0.05 ug/g dry	<0.05	-	<0.05	-
Carbon Tetrachloride	0.05 ug/g dry	<0.05	-	<0.05	-
Chlorobenzene	0.05 ug/g dry	<0.05	-	<0.05	-
Chloroform	0.05 ug/g dry	<0.05	-	<0.05	-
Dibromochloromethane	0.05 ug/g dry	<0.05	-	<0.05	-



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Client: Paterson Group Consulting Engineers

Client PO: 56968 Project Description: PE2709

	Client ID: Sample Date:	BH5-23-SS6 28-Feb-23 09:00	BH6-23-SS5 01-Mar-23 09:00	DUP 28-Feb-23 09:00	BH1-23-SS2 24-Feb-23 09:00
	Sample ID: MDL/Units	2310245-09 Soil	2310245-10 Soil	2310245-11 Soil	2310245-12 Soil
Dichlorodifluoromethane	0.05 ug/g dry	<0.05	-	<0.05	-
1,2-Dichlorobenzene	0.05 ug/g dry	<0.05	-	<0.05	-
1,3-Dichlorobenzene	0.05 ug/g dry	<0.05	-	<0.05	-
1,4-Dichlorobenzene	0.05 ug/g dry	<0.05	-	<0.05	-
1,1-Dichloroethane	0.05 ug/g dry	<0.05	-	<0.05	-
1,2-Dichloroethane	0.05 ug/g dry	<0.05	-	<0.05	-
1,1-Dichloroethylene	0.05 ug/g dry	<0.05	-	<0.05	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	-	<0.05	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	-	<0.05	-
1,2-Dichloropropane	0.05 ug/g dry	<0.05	-	<0.05	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	-	<0.05	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	-	<0.05	-
1,3-Dichloropropene, total	0.05 ug/g dry	<0.05	-	<0.05	-
Ethylbenzene	0.05 ug/g dry	0.10	-	0.08	-
Ethylene dibromide (dibromoethane, 1	0.05 ug/g dry	<0.05	-	<0.05	-
Hexane	0.05 ug/g dry	<0.05	-	<0.05	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	<0.50	-	<0.50	-
Methyl Isobutyl Ketone	0.50 ug/g dry	<0.50	-	<0.50	-
Methyl tert-butyl ether	0.05 ug/g dry	<0.05	-	<0.05	-
Methylene Chloride	0.05 ug/g dry	<0.05	-	<0.05	-
Styrene	0.05 ug/g dry	<0.05	-	<0.05	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	<0.05	-	<0.05	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	<0.05	-	<0.05	-
Tetrachloroethylene	0.05 ug/g dry	<0.05	-	<0.05	-
Toluene	0.05 ug/g dry	<0.05	-	<0.05	-
1,1,1-Trichloroethane	0.05 ug/g dry	<0.05	-	<0.05	-
1,1,2-Trichloroethane	0.05 ug/g dry	<0.05	-	<0.05	-
Trichloroethylene	0.05 ug/g dry	<0.05	-	<0.05	-
Trichlorofluoromethane	0.05 ug/g dry	<0.05	-	<0.05	-
Vinyl chloride	0.02 ug/g dry	<0.02	-	<0.02	-
m,p-Xylenes	0.05 ug/g dry	0.22	-	0.21	-
o-Xylene	0.05 ug/g dry	<0.05	-	<0.05	-
Xylenes, total	0.05 ug/g dry	0.22	-	0.21	-
4-Bromofluorobenzene	Surrogate	98.2%	-	98.4%	-
Dibromofluoromethane	Surrogate	108%	-	104%	-



Order #: 2310245

Report Date: 13-Mar-2023

Order Date: 8-Mar-2023

Client: Paterson Group Consulting Engineers

Client PO: 56968 **Project Description: PE2709**

	Client ID: Sample Date: Sample ID: MDL/Units	BH5-23-SS6 28-Feb-23 09:00 2310245-09 Soil	BH6-23-SS5 01-Mar-23 09:00 2310245-10 Soil	DUP 28-Feb-23 09:00 2310245-11 Soil	BH1-23-SS2 24-Feb-23 09:00 2310245-12 Soil
Toluene-d8	Surrogate	121%	-	120%	_
Benzene	0.02 ug/g dry	-	<0.02	_	<0.02
Ethylbenzene	0.05 ug/g dry	-	<0.05	_	<0.05
Toluene	0.05 ug/g dry	-	<0.05	_	<0.05
m,p-Xylenes	0.05 ug/g dry	-	<0.05	_	<0.05
o-Xylene	0.05 ug/g dry	-	<0.05	_	<0.05
Xylenes, total	0.05 ug/g dry	-	<0.05	_	<0.05
Toluene-d8	Surrogate		123%	_	122%
Hydrocarbons	Jan		12070	<u> </u>	12270
F1 PHCs (C6-C10)	7 ug/g dry	<7	<7	8	<7
F2 PHCs (C10-C16)	4 ug/g dry	<4	<4	28	17
F3 PHCs (C16-C34)	8 ug/g dry	<8	15	<8	107
F4 PHCs (C34-C50)	6 ug/g dry	<6	11	<6	70
Semi-Volatiles				<u> </u>	1
Acenaphthene	0.02 ug/g dry	-	<0.02	-	-
Acenaphthylene	0.02 ug/g dry	-	<0.02	-	-
Anthracene	0.02 ug/g dry	-	<0.02	-	-
Benzo [a] anthracene	0.02 ug/g dry	-	<0.02	-	-
Benzo [a] pyrene	0.02 ug/g dry	-	<0.02	-	-
Benzo [b] fluoranthene	0.02 ug/g dry	-	<0.02	-	-
Benzo [g,h,i] perylene	0.02 ug/g dry	-	<0.02	-	-
Benzo [k] fluoranthene	0.02 ug/g dry	-	<0.02	-	-
Chrysene	0.02 ug/g dry	-	<0.02	-	-
Dibenzo [a,h] anthracene	0.02 ug/g dry	-	<0.02	-	-
Fluoranthene	0.02 ug/g dry	-	<0.02	-	-
Fluorene	0.02 ug/g dry	-	<0.02	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	-	<0.02	-	-
1-Methylnaphthalene	0.02 ug/g dry	-	<0.02	-	-
2-Methylnaphthalene	0.02 ug/g dry	-	<0.02	-	-
Methylnaphthalene (1&2)	0.04 ug/g dry	-	<0.04	-	-
Naphthalene	0.01 ug/g dry	-	<0.01	-	-
Phenanthrene	0.02 ug/g dry	-	<0.02	-	-
Pyrene	0.02 ug/g dry	-	<0.02	-	-
2-Fluorobiphenyl	Surrogate	-	105%	-	-
Terphenyl-d14	Surrogate	-	116%	-	-



Chloroform

Order #: 2310245

Report Date: 13-Mar-2023

Order Date: 8-Mar-2023

Project Description: PE2709

Client: Paterson Group Consulting Engineers

Client PO: 56968

BH3-23-SS5 Client ID: BH1-23-SS5 BH3-23-SS6 Sample Date: 24-Feb-23 09:00 27-Feb-23 09:00 27-Feb-23 09:00 2310245-13 2310245-14 2310245-15 Sample ID: Soil Soil Soil MDL/Units **Physical Characteristics** % Solids 0.1 % by Wt. 92.2 90.5 91.1 General Inorganics 0.01 N/A SAR 1.54 2.07 Conductivity 5 uS/cm 260 317 0.05 pH Units 7.88 pΗ Metals Antimony 1.0 ug/g dry <1.0 <1.0 Arsenic 1.0 ug/g dry 2.9 2.5 Barium 1.0 ug/g dry 87.3 68.1 0.5 ug/g dry <0.5 <0.5 Beryllium Boron 5.0 ug/g dry 5.3 8.1 0.5 ug/g dry <0.5 <0.5 Cadmium Chromium 5.0 ug/g dry 9.9 10.8 Chromium (VI) 0.2 ug/g dry <0.2 <0.2 Cobalt 1.0 ug/g dry 3.2 2.8 < 5.0 5.0 ug/g dry <5.0 Copper 3.3 Lead 1.0 ug/g dry 2.8 Mercury 0.1 ug/g dry < 0.1 < 0.1 <1.0 Molybdenum 1.0 ug/g dry <1.0 6.8 7.2 Nickel 5.0 ug/g dry <1.0 Selenium 1.0 ug/g dry <1.0 Silver 0.3 ug/g dry < 0.3 < 0.3 _ 1.0 ug/g dry <1.0 Thallium <1.0 1.0 ug/g dry <1.0 <1.0 Uranium _ 15.7 10.0 ug/g dry 12.9 Vanadium 20.0 ug/g dry <20.0 <20.0 Zinc _ Volatiles Acetone 0.50 ug/g dry < 0.50 Benzene 0.02 ug/g dry _ _ < 0.02 _ 0.05 ug/g dry < 0.05 Bromodichloromethane Bromoform 0.05 ug/g dry _ _ < 0.05 _ 0.05 ug/g dry < 0.05 Bromomethane 0.05 ug/g dry < 0.05 Carbon Tetrachloride _ _ _ Chlorobenzene 0.05 ug/g dry < 0.05

OTTAWA - MISSISSAUGA - HAMILTON - KINGSTON - LONDON - NIAGARA - WINDSOR - RICHMOND HILL

_

_

0.05 ug/g dry

-

< 0.05



Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

Certificate of Analysis
Client: Paterson Group Consulting Engineers

Client PO: 56968 Project I

Project Description: PE2709

	Client ID:	BH1-23-SS5	BH3-23-SS5	BH3-23-SS6	-
	Sample Date: Sample ID:	24-Feb-23 09:00 2310245-13	27-Feb-23 09:00 2310245-14	27-Feb-23 09:00 2310245-15	-
	MDL/Units	Soil	Soil	Soil	-
Dibromochloromethane	0.05 ug/g dry	-	-	<0.05	-
Dichlorodifluoromethane	0.05 ug/g dry	-	-	<0.05	-
1,2-Dichlorobenzene	0.05 ug/g dry	-	-	<0.05	-
1,3-Dichlorobenzene	0.05 ug/g dry	-	-	<0.05	-
1,4-Dichlorobenzene	0.05 ug/g dry	-	-	<0.05	-
1,1-Dichloroethane	0.05 ug/g dry	-	-	<0.05	-
1,2-Dichloroethane	0.05 ug/g dry	-	-	<0.05	-
1,1-Dichloroethylene	0.05 ug/g dry	-	-	<0.05	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	-	-	<0.05	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	-	-	<0.05	-
1,2-Dichloropropane	0.05 ug/g dry	-	-	<0.05	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	-	-	<0.05	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	-	-	<0.05	-
1,3-Dichloropropene, total	0.05 ug/g dry	-	-	<0.05	-
Ethylbenzene	0.05 ug/g dry	-	-	<0.05	-
Ethylene dibromide (dibromoethane, 1	0.05 ug/g dry	-	-	<0.05	-
Hexane	0.05 ug/g dry	-	-	<0.05	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	-	-	<0.50	-
Methyl Isobutyl Ketone	0.50 ug/g dry	-	-	<0.50	-
Methyl tert-butyl ether	0.05 ug/g dry	-	-	<0.05	-
Methylene Chloride	0.05 ug/g dry	-	-	<0.05	-
Styrene	0.05 ug/g dry	-	-	<0.05	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	-	-	<0.05	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	-	-	<0.05	-
Tetrachloroethylene	0.05 ug/g dry	-	-	<0.05	-
Toluene	0.05 ug/g dry	-	-	<0.05	-
1,1,1-Trichloroethane	0.05 ug/g dry	-	-	<0.05	-
1,1,2-Trichloroethane	0.05 ug/g dry	-	-	<0.05	-
Trichloroethylene	0.05 ug/g dry	-	-	<0.05	-
Trichlorofluoromethane	0.05 ug/g dry	-	-	<0.05	-
Vinyl chloride	0.02 ug/g dry	-	-	<0.02	-
m,p-Xylenes	0.05 ug/g dry	-	-	<0.05	-
o-Xylene	0.05 ug/g dry	-	-	<0.05	-
Xylenes, total	0.05 ug/g dry	-	-	<0.05	-
4-Bromofluorobenzene	Surrogate	-	-	103%	-



Order #: 2310245

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Order Date: 8-Mar-2023

Client: Paterson Group Consulting Engineers

Client PO: 56968 Project Description: PE2709

	Client ID:	BH1-23-SS5	BH3-23-SS5	BH3-23-SS6	_
	Sample Date:	24-Feb-23 09:00	27-Feb-23 09:00	27-Feb-23 09:00	-
	Sample ID:	2310245-13	2310245-14	2310245-15	-
	MDL/Units	Soil	Soil	Soil	-
Dibromofluoromethane	Surrogate	-	-	106%	-
Toluene-d8	Surrogate	-	-	125%	-
Benzene	0.02 ug/g dry	<0.02	<0.02	-	-
Ethylbenzene	0.05 ug/g dry	<0.05	<0.05	-	-
Toluene	0.05 ug/g dry	<0.05	<0.05	-	-
m,p-Xylenes	0.05 ug/g dry	<0.05	<0.05	-	-
o-Xylene	0.05 ug/g dry	<0.05	<0.05	-	-
Xylenes, total	0.05 ug/g dry	<0.05	<0.05	-	-
Toluene-d8	Surrogate	125%	119%	-	-
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	-



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 Client:
 Paterson Group Consulting Engineers
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 Client PO:
 56968
 Project Description: PE2709

Method Quality Control: Blank

Analyta	5	Reporting		Source		%REC		RPD	.
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
General Inorganics									
Conductivity	ND	5	uS/cm						
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 Beryllium	ND ND	1.0 0.5	ug/g ug/g						
Boron	ND ND	5.0	ug/g 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 Mercury	ND ND	1.0 0.1	ug/g ug/g						
Molybdenum	ND ND	1.0	ug/g 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 Zinc	ND ND	10.0 20.0	ug/g						
Semi-Volatiles	ND	20.0	ug/g						
Acenaphthene	ND	0.02	ug/g						
Acenaphthylene	ND	0.02	ug/g 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 Chrysene	ND ND	0.02 0.02	ug/g						
Dibenzo [a,h] anthracene	ND ND	0.02	ug/g ug/g						
Fluoranthene	ND	0.02	ug/g						
Fluorene	ND	0.02	ug/g						
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g						
1-Methylnaphthalene	ND	0.02	ug/g						
2-Methylnaphthalene	ND	0.02	ug/g						
Methylnaphthalene (1&2) Naphthalene	ND	0.04	ug/g						
Naphthalene Phenanthrene	ND ND	0.01 0.02	ug/g						
Pyrene	ND ND	0.02	ug/g ug/g						
Surrogate: 2-Fluorobiphenyl	1.56	0.02	ug/g ug/g		117	50-140			
Surrogate: Terphenyl-d14	1.64		ug/g		123	50-140			
Volatiles			- -						
Acetone	ND	0.50	ug/g						
Benzene	ND	0.02	ug/g						
Bromodichloromethane	ND	0.05	ug/g						
Bromoform	ND	0.05	ug/g						
Bromomethane	ND	0.05	ug/g						
Carbon Tetrachloride	ND	0.05	ug/g						
Chlorobenzene	ND	0.05	ug/g						



Report Date: 13-Mar-2023

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Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 56968

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
-				Nesuit	701120	Liiiii		Liiiii	
Chloroform	ND	0.05	ug/g						
Dibromochloromethane	ND	0.05	ug/g						
Dichlorodifluoromethane	ND	0.05	ug/g						
1,2-Dichlorobenzene	ND	0.05	ug/g						
1,3-Dichlorobenzene	ND	0.05	ug/g						
1,4-Dichlorobenzene	ND	0.05	ug/g						
1,1-Dichloroethane	ND	0.05	ug/g						
1,2-Dichloroethane	ND	0.05	ug/g						
1,1-Dichloroethylene	ND	0.05	ug/g						
cis-1,2-Dichloroethylene	ND	0.05	ug/g						
trans-1,2-Dichloroethylene	ND	0.05	ug/g						
1,2-Dichloropropane	ND	0.05	ug/g						
cis-1,3-Dichloropropylene	ND	0.05	ug/g						
trans-1,3-Dichloropropylene	ND	0.05	ug/g						
1,3-Dichloropropene, total	ND	0.05	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Ethylene dibromide (dibromoethane, 1,2	ND	0.05	ug/g						
Hexane	ND	0.05	ug/g						
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g						
Methyl Isobutyl Ketone	ND	0.50	ug/g						
Methyl tert-butyl ether	ND	0.05	ug/g						
Methylene Chloride	ND	0.05	ug/g						
Styrene	ND	0.05	ug/g						
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g						
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g						
Tetrachloroethylene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
1,1,1-Trichloroethane	ND	0.05	ug/g						
1,1,2-Trichloroethane	ND	0.05	ug/g						
Trichloroethylene	ND	0.05	ug/g ug/g						
Trichlorofluoromethane	ND ND	0.05	ug/g ug/g						
Vinyl chloride	ND	0.02	ug/g ug/g						
m,p-Xylenes	ND	0.05	ug/g ug/g						
o-Xylene	ND ND	0.05	ug/g ug/g						
Xylenes, total	ND ND	0.05	ug/g ug/g						
Surrogate: 4-Bromofluorobenzene	9.44	0.03	ug/g ug/g		118	50-140			
Surrogate: Dibromofluoromethane	9.44 8.12				102	50-140 50-140			
-			ug/g						
Surrogate: Toluene-d8	9.93		ug/g		124	50-140			
Benzene	ND	0.02	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g						
Surrogate: Toluene-d8	9.93		ug/g		124	50-140			



Order #: 2310245

Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 8-Mar-2023

 Client PO:
 56968
 Project Description: PE2709

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
General Inorganics									
SAR	0.78	0.01	N/A	0.76			2.6	30	
Conductivity	214	5	uS/cm	216			0.9	5	
pH	7.18	0.05	pH Units	7.20			0.3	2.3	
Hydrocarbons	7.10	0.00	pri onito	7.20			0.0	2.0	
F1 PHCs (C6-C10)	ND	7	ug/g	ND			NC	40	
F2 PHCs (C10-C16)	ND	4	ug/g	ND			NC	30	
F3 PHCs (C16-C34)	9	8	ug/g	14			39.7	30	
F4 PHCs (C34-C50)	ND	6	ug/g	10			NC	30	
Metals			-9.9						
Antimony	ND	1.0	ug/g	ND			NC	30	
Arsenic	11.1	1.0	ug/g	9.8			12.6	30	
Barium	37.0	1.0	ug/g	37.7			1.7	30	
Beryllium	1.4	0.5	ug/g	1.3			4.2	30	
Boron	23.2	5.0	ug/g	20.6			11.8	30	
Cadmium	ND	0.5	ug/g	ND			NC	30	
Chromium (VI)	0.7	0.3	ug/g ug/g	1.0			25.5	35	
Chromium	29.4	5.0	ug/g ug/g	28.0			4.6	30	
Cobalt	18.9	1.0		18.1			4.0	30	
	58.3	5.0	ug/g	55.7			4.0 4.6	30	
Copper			ug/g						
Lead	7.8	1.0	ug/g	6.3			20.8	30	
Melyhdanum	ND	0.1	ug/g	ND			NC	30	
Molybdenum	ND	1.0	ug/g	ND			NC	30	
Nickel	36.6	5.0	ug/g	34.7			5.4	30	
Selenium	ND	1.0	ug/g	ND			NC	30	
Silver	ND	0.3	ug/g	ND			NC	30	
Thallium	ND	1.0	ug/g	ND			NC	30	
Uranium	ND	1.0	ug/g	ND			NC	30	
Vanadium	35.7	10.0	ug/g	33.7			5.8	30	
Zinc	78.9	20.0	ug/g	75.4			4.6	30	
Physical Characteristics									
% Solids	93.2	0.1	% by Wt.	93.2			0.1	25	
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g	ND			NC	40	
Acenaphthylene	ND	0.02	ug/g	ND			NC	40	
Anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] pyrene	ND	0.02	ug/g	ND			NC	40	
Benzo [b] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Benzo [g,h,i] perylene	ND	0.02	ug/g	ND			NC	40	
Benzo [k] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Chrysene	ND	0.02	ug/g	ND			NC	40	
Dibenzo [a,h] anthracene	ND	0.02	ug/g	ND			NC	40	
Fluoranthene	ND	0.02	ug/g	ND			NC	40	
Fluorene	ND	0.02	ug/g	ND			NC	40	
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g	ND			NC	40	
1-Methylnaphthalene	ND	0.02	ug/g	ND			NC	40	
2-Methylnaphthalene	ND	0.02	ug/g	ND			NC	40	
Naphthalene	ND	0.01	ug/g	ND			NC	40	
Phenanthrene	ND	0.02	ug/g	ND			NC	40	
Pyrene	ND	0.02	ug/g	ND			NC	40	
Surrogate: 2-Fluorobiphenyl	1.28	5.52	ug/g		83.9	50-140			
Surrogate: Terphenyl-d14	1.44				93.9	50-140			
/olatiles	1.77		ug/g		30.3	JU-140			
Acetone	ND	0.50	uala	ND			NC	50	
			ug/g					50 50	
Benzene	ND	0.02	ug/g	ND			NC	5 U	



Order #: 2310245

Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 8-Mar-2023

 Client PO:
 56968
 Project Description: PE2709

Method Quality Control: Duplicate

A mali da		Reporting				%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Bromodichloromethane	ND	0.05	ug/g	ND			NC	50	
Bromoform	ND	0.05	ug/g	ND			NC	50	
Bromomethane	ND	0.05	ug/g	ND			NC	50	
Carbon Tetrachloride	ND	0.05	ug/g	ND			NC	50	
Chlorobenzene	ND	0.05	ug/g	ND			NC	50	
Chloroform	ND	0.05	ug/g	ND			NC	50	
Dibromochloromethane	ND	0.05	ug/g	ND			NC	50	
Dichlorodifluoromethane	ND	0.05	ug/g	ND			NC	50	
1,2-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,3-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,4-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,1-Dichloroethane	ND	0.05	ug/g	ND			NC	50	
1,2-Dichloroethane	ND	0.05	ug/g	ND			NC	50	
1,1-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
cis-1,2-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
rans-1,2-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
1,2-Dichloropropane	ND	0.05	ug/g ug/g	ND			NC	50	
sis-1,3-Dichloropropylene	ND	0.05	ug/g ug/g	ND			NC	50	
rans-1,3-Dichloropropylene	ND	0.05	ug/g	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g ug/g	ND			NC	50	
Ethylene dibromide (dibromoethane, 1,2	ND	0.05	ug/g ug/g	ND			NC	50	
lexane	ND	0.05	ug/g ug/g	ND			NC	50	
Nethyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g ug/g	ND			NC	50	
Methyl Isobutyl Ketone	ND	0.50	ug/g ug/g	ND			NC	50 50	
Methyl tert-butyl ether	ND	0.05	ug/g ug/g	ND			NC	50 50	
Methylene Chloride	ND	0.05	ug/g ug/g	ND			NC	50 50	
Styrene	ND	0.05	ug/g ug/g	ND			NC	50 50	
,1,1,2-Tetrachloroethane	ND	0.05	ug/g ug/g	ND			NC	50 50	
,1,2-Tetrachioroethane	ND	0.05	ug/g ug/g	ND			NC	50 50	
etrachloroethylene	ND ND	0.05	ug/g ug/g	ND			NC	50 50	
oluene	ND ND	0.05	ug/g ug/g	ND ND			NC NC	50 50	
,1,1-Trichloroethane	ND ND	0.05	ug/g ug/g	ND ND			NC NC	50 50	
,1,2-Trichloroethane	ND ND	0.05	ug/g ug/g	ND ND			NC NC	50 50	
richloroethylene	ND ND	0.05	ug/g ug/g	ND ND			NC NC	50 50	
Trichloroethylerie Frichlorofluoromethane	ND ND	0.05	ug/g ug/g	ND ND			NC NC	50 50	
inchloronuoromethane /invl chloride	ND ND	0.05		ND ND			NC NC	50 50	
n,p-Xylenes	ND ND	0.02	ug/g ug/g	ND ND			NC NC	50 50	
i,p-Ayleries -Xylene	ND ND	0.05	ug/g ug/g	ND ND			NC NC	50 50	
Surrogate: 4-Bromofluorobenzene	8.72	0.03		שויי	102	50-140	140	50	
•	8.72 8.95		ug/g		102	50-140 50-140			
Surrogate: Dibromofluoromethane			ug/g						
Surrogate: Toluene-d8	10.4	0.00	ug/g		121	50-140		5 0	
Benzene	ND	0.02	ug/g	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g	ND			NC	50	
oluene	ND	0.05	ug/g	ND			NC	50	
n,p-Xylenes	ND	0.05	ug/g	ND			NC	50	
-Xylene	ND	0.05	ug/g	ND			NC	50	
Surrogate: Toluene-d8	10.4		ug/g		121	50-140			



Order #: 2310245

Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 8-Mar-2023

 Client PO:
 56968
 Project Description: PE2709

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	182	7	ug/g	ND	91.2	80-120			
F2 PHCs (C10-C16)	119	4	ug/g	ND	130	60-140			
F3 PHCs (C16-C34)	304	8	ug/g	14	129	60-140			
F4 PHCs (C34-C50)	171	6	ug/g	10	113	60-140			
Metals									
Antimony	35.6	1.0	ug/g	ND	70.9	70-130			
Arsenic	52.4	1.0	ug/g	3.9	97.0	70-130			
Barium	60.2	1.0	ug/g	15.1	90.4	70-130			
Beryllium	49.0	0.5	ug/g	0.5	96.9	70-130			
Boron	53.9	5.0	ug/g ug/g	8.3	91.3	70-130			
Cadmium	46.1	0.5	ug/g	ND	92.2	70-130			
Chromium (VI)	5.2	0.2	ug/g ug/g	1.0	69.5	70-130		(QM-05
Chromium	63.2	5.0		11.2	104	70-130		`	ZIVI-03
Cobalt	56.9	1.0	ug/g	7.3	99.3	70-130 70-130			
Copper	71.3	5.0	ug/g	7.3 22.3	99.3 98.0	70-130 70-130			
Lead	71.3 47.8	1.0	ug/g	22.5 2.5	90.5	70-130 70-130			
	1.42	0.1	ug/g	ND	90.5	70-130			
Melvindenum			ug/g						
Molybdenum	50.5	1.0	ug/g	ND	101	70-130			
Nickel	63.5	5.0	ug/g	13.9	99.1	70-130			
Selenium	45.7	1.0	ug/g	ND	91.1	70-130			
Silver	42.6	0.3	ug/g	ND	85.0	70-130			
Thallium	45.6	1.0	ug/g	ND	91.0	70-130			
Uranium	47.5	1.0	ug/g	ND	94.4	70-130			
Vanadium Zina	66.6	10.0	ug/g	13.5	106	70-130			
Zinc	75.7	20.0	ug/g	30.2	91.2	70-130			
Semi-Volatiles									
Acenaphthene	0.144	0.02	ug/g	ND	75.3	50-140			
Acenaphthylene	0.131	0.02	ug/g	ND	68.5	50-140			
Anthracene	0.139	0.02	ug/g	ND	72.8	50-140			
Benzo [a] anthracene	0.142	0.02	ug/g	ND	74.5	50-140			
Benzo [a] pyrene	0.149	0.02	ug/g	ND	77.8	50-140			
Benzo [b] fluoranthene	0.204	0.02	ug/g	ND	107	50-140			
Benzo [g,h,i] perylene	0.131	0.02	ug/g	ND	68.4	50-140			
Benzo [k] fluoranthene	0.155	0.02	ug/g	ND	80.9	50-140			
Chrysene	0.172	0.02	ug/g	ND	90.0	50-140			
Dibenzo [a,h] anthracene	0.133	0.02	ug/g	ND	69.8	50-140			
Fluoranthene	0.144	0.02	ug/g	ND	75.1	50-140			
Fluorene	0.141	0.02	ug/g	ND	73.7	50-140			
Indeno [1,2,3-cd] pyrene	0.141	0.02	ug/g	ND	74.0	50-140			
1-Methylnaphthalene	0.168	0.02	ug/g	ND	87.7	50-140			
2-Methylnaphthalene	0.171	0.02	ug/g	ND	89.7	50-140			
Naphthalene	0.164	0.01	ug/g	ND	85.6	50-140			
Phenanthrene	0.150	0.02	ug/g	ND	78.5	50-140			
Pyrene	0.140	0.02	ug/g	ND	73.4	50-140			
Surrogate: 2-Fluorobiphenyl	1.17		ug/g		76.4	50-140			
Surrogate: Terphenyl-d14	1.31		ug/g		85.7	50-140			
olatiles									
Acetone	11.3	0.50	ug/g	ND	113	50-140			



Report Date: 13-Mar-2023 Order Date: 8-Mar-2023

Project Description: PE2709

Certificate of Analysis Client: Paterson Group Consulting Engineers

Client PO: 56968

Method Quality Control: Spike

RPD Reporting Source %REC RPD Result Units %REC Notes Analyte Limit Limit Limit Result ND Benzene 3.57 0.02 ug/g 89.3 60-130 Bromodichloromethane 3.64 0.05 ND 90.9 60-130 ug/g 0.05 ND 101 Bromoform 4.02 ug/g 60-130 4.19 0.05 ND 105 50-140 Bromomethane ug/g 3.55 0.05 ND 88.8 60-130 Carbon Tetrachloride ug/g 98.7 Chlorobenzene 3.95 0.05 ug/g ND 60-130 Chloroform 3.69 0.05 ND 92.3 60-130 ug/g Dibromochloromethane 4.20 0.05 ug/g ND 105 60-130 Dichlorodifluoromethane 3.11 0.05 ND 77.7 50-140 ug/g 0.05 ND 88.3 60-130 1.2-Dichlorobenzene 3.53 ug/g ND 60-130 1,3-Dichlorobenzene 3.62 0.05 ug/g 90.5 1,4-Dichlorobenzene 3.45 0.05 ND 86.3 60-130 ug/g 1,1-Dichloroethane 3.97 0.05 ND 99.3 60-130 ug/g 1,2-Dichloroethane 3.63 0.05 ug/g ND 90.9 60-130 1,1-Dichloroethylene 3.78 0.05 ND 94.5 60-130 ug/g cis-1,2-Dichloroethylene 3.59 0.05 ND 89.8 60-130 ug/g trans-1,2-Dichloroethylene 3.59 0.05 ug/g ND 89.8 60-130 1,2-Dichloropropane 3.42 0.05 ug/g ND 85.5 60-130 cis-1,3-Dichloropropylene 3.41 0.05 ug/g ND 85.2 60-130 trans-1,3-Dichloropropylene 3.52 0.05 ND 88.0 60-130 ug/g 60-130 3.86 0.05 ND 96.5 Ethylbenzene ug/g Ethylene dibromide (dibromoethane, 1,2-3.94 0.05 ug/g ND 98.5 60-130 Hexane 4.02 0.05 ND 101 60-130 ug/g Methyl Ethyl Ketone (2-Butanone) 0.50 ND 97.0 50-140 9.70 ug/g Methyl Isobutyl Ketone 8.58 0.50 ug/g ND 85.8 50-140 0.05 ND 76.5 Methyl tert-butyl ether 7.65 50-140 ug/g ND Methylene Chloride 4.10 0.05 ug/g 103 60-130 Styrene 3.94 0.05 ug/g ND 98.5 60-130 1,1,1,2-Tetrachloroethane 4.00 0.05 ND 100 60-130 ug/g 1,1,2,2-Tetrachloroethane 4.22 0.05 ug/g ND 105 60-130 Tetrachloroethylene 4.07 0.05 ND 102 60-130 ug/g Toluene ND 100 4 00 0.05 ug/g 60-130 1,1,1-Trichloroethane 3.58 0.05 ug/g ND 89.6 60-130 1,1,2-Trichloroethane 3.45 0.05 ug/g ND 86.2 60-130 Trichloroethylene 3 37 0.05 ug/g ND 84 1 60-130 Trichlorofluoromethane 4.47 0.05 ug/g ND 112 50-140 Vinyl chloride 3.53 0.02 ug/g ND 88.2 50-140 m,p-Xylenes 8.06 0.05 ug/g ND 101 60-130 o-Xylene 4.11 0.05 ug/g ND 103 60-130 Surrogate: 4-Bromofluorobenzene 7.53 ug/g 94.2 50-140 Surrogate: Dibromofluoromethane 7.79 97.4 50-140 ug/g Surrogate: Toluene-d8 8.30 104 50-140 ug/g Benzene 3.57 0.02 ug/g ND 89.3 60-130 Ethylbenzene 3.86 0.05 ug/g ND 96.5 60-130 0.05 60-130 Toluene 4.00 ug/g ND 100 8.06 0.05 ND 101 60-130 m,p-Xylenes ug/g 0.05 ND 103 60-130 o-Xylene 4 11 ug/g Surrogate: Toluene-d8 8.30 104 50-140

ug/g



 Certificate of Analysis
 Report Date: 13-Mar-2023

 Client: Paterson Group Consulting Engineers
 Order Date: 8-Mar-2023

Project Description: PE2709

Qualifier Notes:

Client PO: 56968

Sample Qualifiers:

- 1: Elevated reporting limits due to the nature of the sample matrix.
- 2: Some peak(s) in the GC-FID Chromatogram are not typical of petroleum hydrocarbon distillates. May be the result of high concentrations of non-mineral based compounds not completely removed by the method cleanup. Results may be biased high.

QC Qualifiers:

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

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.



Laurent Blvd. rio K1G4J8 1947 racellabs.com

Chain Of Custody Paracel Order Number (Lab Use Only) (Lab Use Only)

Contact Name: Karyn Munch			_	ect Ref.	re	2709		73							Pa	ge]	of 4	
v.doi.622.		_	Quo												Turna	round	d Time	
9 Auriga Drive			PO#		5	968								1 day	,			3 day
Telephone: 4-13 226 7381		_	E-ma	il: }			(5004[0,0).							2 day			I	Regul
▼REG 153/04 □					Becs	be@pyter	rsongrap.ca songroup.ca	1					Date	e Requ	ired:			
Table 1 Regulat		-	Matrix				round Water)	193	19.50			435	SVT.					
☐ Table 2 ☐ Ind/Comm ☐ a	PWQO		SW (So	ırface	Water)	SS (Storm/Sa	nitary Sewer)					Re	quire	d Ana	lysis			
Table 3 Agri/Other	MISA		_	P(Paint)	A (Air) O (Oti	ner)	EX									T	T
Table Mun.	SU - Storm			ners				-F4+BTEX			۵							
For RSC: Yes No Cther:		v	lume	Containers		Sample	Taken	F1-F			by IC			·	SAR			
Sample ID/Location Name		Matrix	Air Volume	o.	-	Data		PHCs	VOCs	PAHs	Metals by ICP		5	B (HWS)	_	НΑ		
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Revsion 4.0





ead Office 0-2319 St. Laurent Blvd. tawa, Ontario K1G 4,8 1-800-749-1947

Paracel Order Number (Lab Use Only)

Chain Of Custody (Lab Use Only)

Client Name: Patison Contact Name: Karyn Munch	_	_	PE 2709									Page	2 of 4	
100		ote#:									1	Turnarou	ınd Time	9
9 Auriga Prive	PO	#:	56968 Munch@piter Berube @pater	5							1 day		[☐ 3 day
elephone: 613 226 738)	E-11	ail: ¥	Munch@pater	5019129.69							2 day		[Regula
▼ REG 153/04 □		5	Berube 6 bater	songrup. (9						Date	Requi	red:		
Table 1 Res/Park Classes Other Regulation			S (Soil/Sed.) GW (G		3)(9				310	ST.	9.7.		langer.	
Table 2 I Ind/Comm II	SW (S	urface	Water) SS (Storm/Sa	nitary Sewer)					Rec	quire	Anal	ysis		
Table 3 April/Only		P (Paint) A (Air) O (Oti	ner)	X								TT	Т
Table SU-Storm SU-Storm		673			F1-F4+BTEX			۵						
For RSC: Yes No Other:	a B	Containers	Sample	Taken	1-F4			oy 10				X X		
Sample ID/Location Name	Matrix Air Volume	of Co			PHCs F	VOCs	£	Metals by ICP		_	B (HWS)	EC/SAR		
pol	_	#	Date	Time	PH	8	PAHs	Me	Ÿ	S.	B (F	Εc		
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Chain Of Custody Paracel Order Number (Lab Use Only) (Lab Use Only)

lient Name: Pakison	_	Deal	net D-C		. /											
ontact Name: K4/Yn Munch		_	ect Ref:	PE 2709									Pa	ge 3 (of 4	
Add 622;		Quot									\vdash		-	round	-	-
9 Auriga Drive		PO#		56968							1.	1 day				3 day
lephone: 613 236 726		E-ma	il: K	Munch@pater	son goop. (a	1					-	2 day				Regul
FREG 153/04 DREG 406/19			5	Berube@pqte	SON GOUP. CO	9						Requ				vegui
Othor Pandat						· Interest				Lawrence	Date	nequ	irea.			
Table 1 Res/Park Med/Fine REG 558 PWQO	1 '	Matrix ' SW (Su	Type: urface \	S (Soil/Sed.) GW (G Water) SS (Storm/Sa	round Water)					Re	quire	d Anal	lysis			
Table 2 Ind/Comm Coarse CCME MISA		(00	P (F	Paint) A (Air) O (Oth	er)	V										_
Table 3 Agri/Other Su-Sani Su-Storm		T	50			⊢Ë.										
For Psc. of		e	Containers	Sample	Taken	PHCs F1-F4+BTEX			ICP				0.0			
- other	ž	Air Volume	Cont	Jumple	Idkell	<u>F</u>			s by			(3)	348			
Sample ID/Location Name BH3-23-555	Matrix	Air V	# of	Date	Time	유	Vocs	PAHs	Metals by ICP	Hg	CrVI	B (HWS)	EC/SAR	ЬH		
BH3-23-556	5		2	Feb 27, 2-27		X	>	ο.	χ	X	x	В		-	+	+
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sts:女 These samples are already at the lab. The	010	e/ 10	mbo	r is 230	9472				1	Metho	d of Del	ivery:			1	
hed By (Sign): Received By Drive				-									EL	Co	DUKIL	5
NG DV (Print): U -		1	1	A second	ecelved at Lab:	1	7	2	1	/erified	_			9780		
e: Date/Time: Page Lee	3/0	77	17	3 17010	ate/Time:08/	03/	22	12	40	Date/F	me:N	20	-	213	112	5
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Paracel ID: 2310245

LABORATORIES LTD

int Blvd.	Paracel Order Number	
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Chain Of Custody (Lab Use Only)

Cor	ntact Name: Patetson			Pro	ject Ref:	PE 2709					,	-				4	- 11	-
Ado	ntact Name: Karyn Munch Idress:			Que	te#:				-				\vdash		-	age 4	-	
	9 Avriga Prive			PO		56969	ζ						1.	1 day		around	Time □ 3	day
Tele	ephone: 6 B 226 7381			E-ma	ili: ķ /	Munch @ pater	en 9100p. (a						-	2 day				
F	V REG 153/04				5	Munch @ Pater Berube @ pater	san graup. Ca	ı						Requ			₫R	eguia
_	✓ REG 153/04 ☐ REG 406/19 Oth Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558	er Regulation						1850	0.80		100	Marine.	Date	nequ	illeu:			
П.	Table 2 Ind/Comm Coarse Come			SW (S	urface	S (Soil/Sed.) GW (G Water) SS (Storm/Sa	nitary Sewer)					Re	quire	d Ana	lysis			
9 1	Table 3 ☐ Agrl/Other ☐ SU-Sani	☐ MISA ☐ SU-Storm	_	_	P()	Paint) A (Air) O (Oti	ner)	X								П	1	Т
] 1	Table Mun;	□ 20 - 2form			ners			F1-F4+BTEX			Q.							
_	For RSC: Yes No Other:		L	Inme	Containers	Sample	Taken	1-F			oy IC				8			
	Sample ID/Location Name		Matrix	Air Volume	of	0		PHCs F	VOCs	s I	Metals by ICP		-	(HWS)	EC/SAR	h H		
1	BH1-23-552		5	∢	12	Date	Time	_	>	PAHs	Me	Ď	CrVI	B (F	Ec	۵.		
2	BH1-23-555		J	-	2	Feb 24, 2023		X							X	X		
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iquis	shed By (Sign): Blue	Received By Drive	≥r/Dep	ot:									1	92	ACL	2	Con	TEC
quis	shed By (Print): Bryce Lee		NOT THE	_	1	COUSE	eceived at Lab:	1	2		- V	erified	By:	1.	,		10	-0.0
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Revsion 4.0

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Chain of Custody (Blank).xlsx



ad Office +2319 St. Laurent Blvd. IMIG. Ontario K1G 4J8 +800-749-1947 araceleparacellabs.com Paracel Order Number

(Lab Use Only)

Chain Of Custody

(Lab Use Only)

LABORATORIES w.paracellabs.com Client Name: Project Ref: PF 2709 of Page Quote #: Turnaround Time Address: PO#: ☐ 1 day ☐ 3 day Kmunchalpatersong roup, ca Regular 2 day Date Required: REG 153/04 REG 406/19 Other Regulation Matrix Type: \$ (Soil/Sed.) GW (Ground Water) ☐ Table 1 ☐ Res/Park ☐ Med/Fine Required Analysis ☐ PWQO SW (Surface Water) SS (Storm/Sanitary Sewer) ☐ Table 2 ☐ Ind/Comm ☐ Coarse P (Paint) A (Air) O (Other) ☐ CCME ☐ MISA F1-F4+BTEX Table 3 Agri/Other ☐ SU - Sani ☐ SU - Storm # of Containers П ☐ Table Mun: Sample Taken Air Volume Metals by B (HWS) For RSC: Yes No Other: Matrix PHCs PAHs Cr≧ Sample ID/Location Name Date Β̈́I Time B41-23-552 S Feb. 24/2 2 3 4 5 6 7 8 9 10 Method of Delivery: Relinquished By (Sign) Relinquished By (Print); Date/Time Date/Time: pH Verified:

Revsion 4.0



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

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive Ottawa, ON K2E 7T9 Attn: Sam Berube

Client PO: 56893 Project: PE2709

Custody:

Report Date: 2-Mar-2023 Order Date: 27-Feb-2023

Order #: 2309081

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

 Paracel ID
 Client ID

 2309081-01
 BH1-23-SS2

 2309081-02
 BH1-23-SS4

Approved By:

Mark Froto

Mark Foto, M.Sc. Lab Supervisor



Report Date: 02-Mar-2023 Order Date: 27-Feb-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 56893

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	1-Mar-23	2-Mar-23
Mercury by CVAA	EPA 7471B - CVAA, digestion	2-Mar-23	2-Mar-23
PHC F1	CWS Tier 1 - P&T GC-FID	28-Feb-23	28-Feb-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	27-Feb-23	28-Feb-23
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	1-Mar-23	1-Mar-23
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	27-Feb-23	2-Mar-23
REG 153: VOCs by P&T GC/MS	EPA 8260 - P&T GC-MS	28-Feb-23	28-Feb-23
Solids, %	CWS Tier 1 - Gravimetric	28-Feb-23	28-Feb-23



Report Date: 02-Mar-2023 Order Date: 27-Feb-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 56893

					•
	Client ID:	BH1-23-SS2	BH1-23-SS4	_	
	Sample Date:	24-Feb-23 09:00	24-Feb-23 09:00	-	-
	Sample ID:	2309081-01	2309081-02	-	-
	MDL/Units	Soil	Soil	-	-
Physical Characteristics	1 2 2 2 1 1 1 1		_		1 1
% Solids	0.1 % by Wt.	90.7	92.5	-	-
Metals	1.00/				1
Antimony	1.0 ug/g dry	<1.0	-	-	-
Arsenic	1.0 ug/g dry	3.5	-	-	-
Barium	1.0 ug/g dry	62.5	-	-	-
Beryllium	0.5 ug/g dry	<0.5	-	-	-
Boron	5.0 ug/g dry	<5.0	-	-	-
Cadmium	0.5 ug/g dry	<0.5	-	-	-
Chromium	5.0 ug/g dry	14.9	-	-	-
Chromium (VI)	0.2 ug/g dry	<0.2	-	-	-
Cobalt	1.0 ug/g dry	3.9	-	-	-
Copper	5.0 ug/g dry	13.0	-	-	-
Lead	1.0 ug/g dry	61.9	-	-	-
Mercury	0.1 ug/g dry	0.2	-	-	-
Molybdenum	1.0 ug/g dry	<1.0	-	-	-
Nickel	5.0 ug/g dry	8.7	-	-	-
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	17.2	-	-	-
Zinc	20.0 ug/g dry	55.3	-	-	-
Volatiles	•		•		•
Acetone	0.50 ug/g dry	-	<0.50	•	-
Benzene	0.02 ug/g dry	-	<0.02	-	-
Bromodichloromethane	0.05 ug/g dry	-	<0.05	•	-
Bromoform	0.05 ug/g dry	-	<0.05	-	-
Bromomethane	0.05 ug/g dry	-	<0.05	-	-
Carbon Tetrachloride	0.05 ug/g dry	-	<0.05	-	-
Chlorobenzene	0.05 ug/g dry	-	<0.05	-	-
Chloroform	0.05 ug/g dry	-	<0.05	-	-
Dibromochloromethane	0.05 ug/g dry	-	<0.05	-	-
Dichlorodifluoromethane	0.05 ug/g dry	-	<0.05	-	-
1,2-Dichlorobenzene	0.05 ug/g dry	-	<0.05	-	-
1,3-Dichlorobenzene	0.05 ug/g dry	-	<0.05	-	-



Report Date: 02-Mar-2023 Order Date: 27-Feb-2023

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 56893 **Project Description: PE2709**

BH1-23-SS4 Client ID: BH1-23-SS2 Sample Date: 24-Feb-23 09:00 24-Feb-23 09:00 2309081-01 2309081-02 Sample ID: MDL/Units Soil Soil 1,4-Dichlorobenzene 0.05 ug/g dry < 0.05 1,1-Dichloroethane 0.05 ug/g dry < 0.05 1,2-Dichloroethane 0.05 ug/g dry < 0.05 1,1-Dichloroethylene 0.05 ug/g dry < 0.05 0.05 ug/g dry cis-1,2-Dichloroethylene < 0.05 trans-1,2-Dichloroethylene 0.05 ug/g dry < 0.05 0.05 ug/g dry 1,2-Dichloropropane < 0.05 0.05 ug/g dry cis-1,3-Dichloropropylene < 0.05 trans-1,3-Dichloropropylene 0.05 ug/g dry < 0.05 0.05 ug/g dry 1,3-Dichloropropene, total < 0.05 0.05 ug/g dry Ethylbenzene < 0.05 Ethylene dibromide (dibromoethane, 1,2-) 0.05 ug/g dry < 0.05 0.05 ug/g dry Hexane < 0.05 0.50 ug/g dry Methyl Ethyl Ketone (2-Butanone) < 0.50 0.50 ug/g dry Methyl Isobutyl Ketone < 0.50 0.05 ug/g dry Methyl tert-butyl ether < 0.05 0.05 ug/g dry Methylene Chloride < 0.05 0.05 ug/g dry Styrene < 0.05 0.05 ug/g dry 1,1,1,2-Tetrachloroethane < 0.05 0.05 ug/g dry 1,1,2,2-Tetrachloroethane < 0.05 0.05 ug/g dry Tetrachloroethylene < 0.05 0.05 ug/g dry Toluene < 0.05 0.05 ug/g dry 1,1,1-Trichloroethane < 0.05 1,1,2-Trichloroethane 0.05 ug/g dry < 0.05 0.05 ug/g dry Trichloroethylene < 0.05 0.05 ug/g dry Trichlorofluoromethane < 0.05 0.02 ug/g dry Vinyl chloride < 0.02 0.05 ug/g dry m,p-Xylenes < 0.05 0.05 ug/g dry o-Xylene < 0.05 0.05 ug/g dry Xylenes total < 0.05 4-Bromofluorobenzene Surrogate 105% Dibromofluoromethane Surrogate 94.4% Toluene-d8 Surrogate 115% -Hydrocarbons 7 ug/g dry F1 PHCs (C6-C10) <7 4 ug/g dry F2 PHCs (C10-C16) 43

OTTAWA - MISSISSAUGA - HAMILTON - KINGSTON - LONDON - NIAGARA - WINDSOR - RICHMOND HILL



Pyrene

2-Fluorobiphenyl

Terphenyl-d14

Order #: 2309081

Report Date: 02-Mar-2023

Order Date: 27-Feb-2023

Client: Paterson Group Consulting Engineers Client PO: 56893 **Project Description: PE2709**

	Client ID:	BH1-23-SS2	BH1-23-SS4	-	-
	Sample Date:	24-Feb-23 09:00	24-Feb-23 09:00	-	-
	Sample ID:	2309081-01	2309081-02	-	-
	MDL/Units	Soil	Soil	-	-
F3 PHCs (C16-C34)	8 ug/g dry	-	364	-	-
F4 PHCs (C34-C50)	6 ug/g dry	-	197	-	-
Semi-Volatiles	•			•	
Acenaphthene	0.02 ug/g dry	0.03	-	-	-
Acenaphthylene	0.02 ug/g dry	0.05	-	-	-
Anthracene	0.02 ug/g dry	0.09	-	-	-
Benzo [a] anthracene	0.02 ug/g dry	0.23	-	-	-
Benzo [a] pyrene	0.02 ug/g dry	0.26	-	-	-
Benzo [b] fluoranthene	0.02 ug/g dry	0.29	-	-	-
Benzo [g,h,i] perylene	0.02 ug/g dry	0.16	-	-	-
Benzo [k] fluoranthene	0.02 ug/g dry	0.14	-	-	-
Chrysene	0.02 ug/g dry	0.29	-	-	-
Dibenzo [a,h] anthracene	0.02 ug/g dry	0.04	-	-	-
Fluoranthene	0.02 ug/g dry	0.56	-	-	-
Fluorene	0.02 ug/g dry	0.02	-	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	0.14	-	-	-
1-Methylnaphthalene	0.02 ug/g dry	0.13	-	-	-
2-Methylnaphthalene	0.02 ug/g dry	0.18	-	-	-
Methylnaphthalene (1&2)	0.04 ug/g dry	0.31	-	-	-
Naphthalene	0.01 ug/g dry	0.14	-	-	-
Phenanthrene	0.02 ug/g dry	0.42	-	-	-

0.50

101%

115%

0.02 ug/g dry

Surrogate

Surrogate



Report Date: 02-Mar-2023

Order Date: 27-Feb-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 56893

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 Cobalt	ND ND	5.0 1.0	ug/g						
Copper	ND ND	5.0	ug/g						
Lead	ND ND	1.0	ug/g ug/g						
Mercury	ND ND	0.1	ug/g ug/g						
Molybdenum	ND	1.0	ug/g						
Nickel	ND	5.0	ug/g						
Selenium	ND	1.0	ug/g						
Silver	ND	0.3	ug/g						
Thallium	ND	1.0	ug/g						
Uranium	ND	1.0	ug/g						
Vanadium	ND	10.0	ug/g						
Zinc	ND	20.0	ug/g						
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g						
Acenaphthylene	ND	0.02	ug/g						
Anthracene	ND	0.02	ug/g						
Benzo [a] anthracene	ND	0.02	ug/g						
Benzo [a] pyrene	ND	0.02	ug/g						
Benzo [b] fluoranthene	ND	0.02	ug/g						
Benzo [g,h,i] perylene Benzo [k] fluoranthene	ND ND	0.02 0.02	ug/g						
Chrysene	ND ND	0.02	ug/g ug/g						
Dibenzo [a,h] anthracene	ND	0.02	ug/g ug/g						
Fluoranthene	ND	0.02	ug/g						
Fluorene	ND	0.02	ug/g						
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g						
1-Methylnaphthalene	ND	0.02	ug/g						
2-Methylnaphthalene	ND	0.02	ug/g						
Methylnaphthalene (1&2)	ND	0.04	ug/g						
Naphthalene	ND	0.01	ug/g						
Phenanthrene	ND	0.02	ug/g						
Pyrene	ND	0.02	ug/g						
Surrogate: 2-Fluorobiphenyl	1.61		ug/g		121	50-140			
Surrogate: Terphenyl-d14	1.63		ug/g		122	50-140			
V olatiles									
Acetone	ND	0.50	ug/g						
Benzene	ND	0.02	ug/g						
Bromodichloromethane	ND	0.05	ug/g						
Bromoform	ND	0.05	ug/g						
Bromomethane	ND	0.05	ug/g						
Carbon Tetrachloride	ND	0.05	ug/g						
Chlorobenzene	ND	0.05	ug/g						
Chloroform	ND	0.05	ug/g						
Dibromochloromethane	ND	0.05	ug/g						
Dichlorodifluoromethane	ND	0.05	ug/g						



Order #: 2309081

Report Date: 02-Mar-2023

Order Date: 27-Feb-2023

Client: Paterson Group Consulting Engineers Client PO: 56893 **Project Description: PE2709**

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
1,2-Dichlorobenzene	ND	0.05	ug/g						
1,3-Dichlorobenzene	ND	0.05	ug/g						
1,4-Dichlorobenzene	ND	0.05	ug/g						
1,1-Dichloroethane	ND	0.05	ug/g						
1,2-Dichloroethane	ND	0.05	ug/g						
1,1-Dichloroethylene	ND	0.05	ug/g						
cis-1,2-Dichloroethylene	ND	0.05	ug/g						
trans-1,2-Dichloroethylene	ND	0.05	ug/g						
1,2-Dichloropropane	ND	0.05	ug/g						
cis-1,3-Dichloropropylene	ND	0.05	ug/g						
trans-1,3-Dichloropropylene	ND	0.05	ug/g						
1,3-Dichloropropene, total	ND	0.05	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Ethylene dibromide (dibromoethane, 1,2-	ND	0.05	ug/g						
Hexane	ND	0.05	ug/g						
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g						
Methyl Isobutyl Ketone	ND	0.50	ug/g						
Methyl tert-butyl ether	ND	0.05	ug/g						
Methylene Chloride	ND	0.05	ug/g						
Styrene	ND	0.05	ug/g						
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g						
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g						
Tetrachloroethylene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
1,1,1-Trichloroethane	ND	0.05	ug/g						
1,1,2-Trichloroethane	ND	0.05	ug/g						
Trichloroethylene	ND	0.05	ug/g						
Trichlorofluoromethane	ND	0.05	ug/g						
Vinyl chloride	ND	0.02	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g						
Surrogate: 4-Bromofluorobenzene	8.49		ug/g		106	50-140			
Surrogate: Dibromofluoromethane	7.36		ug/g		92.0	50-140			
Surrogate: Toluene-d8	8.87		ug/g		111	50-140			



Client PO: 56893

Order #: 2309081

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Order Date: 27-Feb-2023 **Project Description: PE2709**

Report Date: 02-Mar-2023

Analyta		Reporting		Source		%REC	_	RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
lydrocarbons									
F1 PHCs (C6-C10)	ND	7	ug/g	ND			NC	40	
F2 PHCs (C10-C16)	ND	4	ug/g	ND			NC	30	
F3 PHCs (C16-C34)	ND	8	ug/g	ND			NC	30	
F4 PHCs (C34-C50)	ND	6	ug/g	ND			NC	30	
Metals			-9.9						
Antimony	ND	1.0	ug/g	ND			NC	30	
Arsenic	3.3	1.0	ug/g	3.5			6.1	30	
Barium	71.5	1.0	ug/g	62.5			13.5	30	
Beryllium	ND	0.5	ug/g	ND			NC	30	
Boron	ND	5.0	ug/g	ND			NC	30	
Cadmium	ND	0.5	ug/g	ND			NC	30	
Chromium (VI)	ND	0.2	ug/g	ND			NC	35	
Chromium	15.5	5.0	ug/g	14.9			4.4	30	
Cobalt	3.9	1.0	ug/g ug/g	3.9			0.9	30	
Copper	12.2	5.0	ug/g ug/g	13.0			6.7	30	
* *	72.9						16.3	30	
Lead		1.0	ug/g	61.9			4.7		
Melvhdonum	0.217 ND	0.1	ug/g	0.228				30	
Molybdenum	ND	1.0	ug/g	ND			NC	30	
Nickel Selenium	8.5	5.0	ug/g	8.7 ND			2.6	30	
Selenium	ND	1.0	ug/g	ND			NC	30	
Silver	ND	0.3	ug/g	ND			NC	30	
Thallium	ND	1.0	ug/g	ND			NC	30	
Uranium	ND	1.0	ug/g	ND			NC	30	
Vanadium	16.6	10.0	ug/g	17.2			3.6	30	
Zinc	54.7	20.0	ug/g	55.3			1.1	30	
Physical Characteristics									
% Solids	92.5	0.1	% by Wt.	92.3			0.2	25	
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g	ND			NC	40	
Acenaphthylene	ND	0.02	ug/g	ND			NC	40	
Anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] pyrene	ND	0.02	ug/g	ND			NC	40	
Benzo [b] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Benzo [g,h,i] perylene	ND	0.02	ug/g	ND			NC	40	
Benzo [k] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Chrysene	ND	0.02	ug/g	ND			NC	40	
Dibenzo [a,h] anthracene	ND	0.02	ug/g	ND			NC	40	
Fluoranthene	ND	0.02	ug/g	ND			NC	40	
Fluorene	ND	0.02	ug/g	ND			NC	40	
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g	ND			NC	40	
1-Methylnaphthalene	ND	0.02	ug/g ug/g	ND			NC	40	
2-Methylnaphthalene	ND ND	0.02	ug/g ug/g	ND			NC	40	
Naphthalene	ND ND	0.02	ug/g ug/g	ND			NC	40	
Phenanthrene	ND ND	0.01		ND ND			NC NC	40	
		0.02	ug/g	ND ND			NC NC	40 40	
Pyrene Surregate: 2 Fluorabinhanul	ND	0.02	ug/g	ND	00.0	E0 440	NC	40	
Surrogate: 2-Fluorobiphenyl	1.38		ug/g		98.3	50-140			
Surrogate: Terphenyl-d14	1.73		ug/g		124	50-140			
/olatiles									
Acetone	ND	0.50	ug/g	ND			NC	50	
Benzene	ND	0.02	ug/g	ND			NC	50	
Bromodichloromethane	ND	0.05	ug/g	ND			NC	50	
Bromoform	ND	0.05	ug/g	ND			NC	50	
Bromomethane	ND	0.05	ug/g	ND			NC	50	
Carbon Tetrachloride	ND	0.05	ug/g	ND			NC	50	
Chlorobenzene	ND	0.05	ug/g	ND			NC	50	



Report Date: 02-Mar-2023

Order Date: 27-Feb-2023

Project Description: PE2709

Certificate of Analysis

Client PO: 56893

Client: Paterson Group Consulting Engineers

Method Quality Control: Duplicate

		Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Chloroform	ND	0.05	ug/g	ND			NC	50	
Dibromochloromethane	ND	0.05	ug/g	ND			NC	50	
Dichlorodifluoromethane	ND	0.05	ug/g	ND			NC	50	
1,2-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,3-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,4-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,1-Dichloroethane	ND	0.05	ug/g	ND			NC	50	
1,2-Dichloroethane	ND	0.05	ug/g	ND			NC	50	
1,1-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
cis-1,2-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
trans-1,2-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
1,2-Dichloropropane	ND	0.05	ug/g	ND			NC	50	
cis-1,3-Dichloropropylene	ND	0.05	ug/g	ND			NC	50	
trans-1,3-Dichloropropylene	ND	0.05	ug/g	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g	ND			NC	50	
Ethylene dibromide (dibromoethane, 1,2	ND	0.05	ug/g	ND			NC	50	
Hexane	ND	0.05	ug/g	ND			NC	50	
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g	ND			NC	50	
Methyl Isobutyl Ketone	ND	0.50	ug/g	ND			NC	50	
Methyl tert-butyl ether	ND	0.05	ug/g	ND			NC	50	
Methylene Chloride	ND	0.05	ug/g	ND			NC	50	
Styrene	ND	0.05	ug/g	ND			NC	50	
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g	ND			NC	50	
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g	ND			NC	50	
Tetrachloroethylene	ND	0.05	ug/g	ND			NC	50	
Toluene	ND	0.05	ug/g	ND			NC	50	
1,1,1-Trichloroethane	ND	0.05	ug/g	ND			NC	50	
1,1,2-Trichloroethane	ND	0.05	ug/g	ND			NC	50	
Trichloroethylene	ND	0.05	ug/g	ND			NC	50	
Trichlorofluoromethane	ND	0.05	ug/g	ND			NC	50	
Vinyl chloride	ND	0.02	ug/g	ND			NC	50	
m,p-Xylenes	ND	0.05	ug/g	ND			NC	50	
o-Xylene	ND	0.05	ug/g	ND			NC	50	
Surrogate: 4-Bromofluorobenzene	10.1		ug/g		111	50-140	-		
Surrogate: Dibromofluoromethane	8.71		ug/g		95.3	50-140			
Surrogate: Toluene-d8	10.8		ug/g		118	50-140			



Client: Paterson Group Consulting Engineers

Order Date: 27-Feb-2023 Client PO: 56893 **Project Description: PE2709**

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	81.3	80-120			
F2 PHCs (C10-C16)	89	4	ug/g	ND	102	60-140			
F3 PHCs (C16-C34)	228	8	ug/g	ND	107	60-140			
F4 PHCs (C34-C50)	166	6	ug/g	ND	123	60-140			
letals									
Arsenic	48.0	1.0	ug/g	1.4	93.2	70-130			
Barium	69.9	1.0	ug/g	25.0	89.9	70-130			
Beryllium	47.5	0.5	ug/g	ND	94.7	70-130			
Boron	44.1	5.0	ug/g ug/g	ND	85.0	70-130			
Cadmium	43.9	0.5	ug/g ug/g	ND	87.6	70-130			
Chromium (VI)	0.07	0.2		ND	35.0	70-130			QM-05
Chromium	54.7	5.0	ug/g	6.0	97.5	70-130			QIVI-03
Cobalt	48.5	1.0	ug/g	1.6	93.8	70-130			
	48.5 49.8	5.0	ug/g	5.2	93.8 89.3	70-130 70-130			
Copper Lead	49.8 71.4	5.0 1.0	ug/g	5.2 24.8	93.3	70-130 70-130			
			ug/g						
Mercury Molybdenum	1.42 45.8	0.1 1.0	ug/g	0.228 ND	79.3 91.2	70-130 70-130			
<u>-</u>			ug/g						
Nickel	50.4	5.0	ug/g	ND	93.8	70-130			
Selenium	45.0	1.0	ug/g	ND	89.6	70-130			
Silver	40.8	0.3	ug/g	ND	81.4	70-130			
Thallium 	43.4	1.0	ug/g	ND	86.6	70-130			
Uranium	47.6	1.0	ug/g	ND	94.9	70-130			
Vanadium 	54.9	10.0	ug/g	ND	96.0	70-130			
Zinc	64.1	20.0	ug/g	22.1	83.9	70-130			
emi-Volatiles									
Acenaphthene	0.208	0.02	ug/g	ND	119	50-140			
Acenaphthylene	0.172	0.02	ug/g	ND	98.3	50-140			
Anthracene	0.168	0.02	ug/g	ND	95.7	50-140			
Benzo [a] anthracene	0.136	0.02	ug/g	ND	77.9	50-140			
Benzo [a] pyrene	0.148	0.02	ug/g	ND	84.8	50-140			
Benzo [b] fluoranthene	0.186	0.02	ug/g	ND	106	50-140			
Benzo [g,h,i] perylene	0.155	0.02	ug/g	ND	88.8	50-140			
Benzo [k] fluoranthene	0.157	0.02	ug/g	ND	89.8	50-140			
Chrysene	0.209	0.02	ug/g	ND	120	50-140			
Dibenzo [a,h] anthracene	0.156	0.02	ug/g	ND	89.3	50-140			
Fluoranthene	0.178	0.02	ug/g	ND	102	50-140			
Fluorene	0.204	0.02	ug/g	ND	117	50-140			
Indeno [1,2,3-cd] pyrene	0.161	0.02	ug/g	ND	92.1	50-140			
1-Methylnaphthalene	0.231	0.02	ug/g	ND	132	50-140			
2-Methylnaphthalene	0.235	0.02	ug/g	ND	134	50-140			
Naphthalene	0.240	0.01	ug/g	ND	137	50-140			
Phenanthrene	0.209	0.02	ug/g	ND	119	50-140			
Pyrene	0.177	0.02	ug/g	ND	101	50-140			
Surrogate: 2-Fluorobiphenyl	1.61		ug/g		115	50-140			
Surrogate: Terphenyl-d14	1.95		ug/g		139	50-140			
olatiles of the state of the st									
Acetone	12.9	0.50	ug/g	ND	129	50-140			
Benzene	4.55	0.02	ug/g	ND	114	60-130			

Report Date: 02-Mar-2023



Order #: 2309081

Report Date: 02-Mar-2023

Order Date: 27-Feb-2023

Client: Paterson Group Consulting Engineers

Client PO: 56893 **Project Description: PE2709**

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Bromodichloromethane	3.68	0.05	ug/g	ND	92.1	60-130			
Bromoform	3.39	0.05	ug/g	ND	84.7	60-130			
Bromomethane	4.35	0.05	ug/g	ND	109	50-140			
Carbon Tetrachloride	3.04	0.05	ug/g	ND	75.9	60-130			
Chlorobenzene	4.11	0.05	ug/g	ND	103	60-130			
Chloroform	4.06	0.05	ug/g	ND	101	60-130			
Dibromochloromethane	3.45	0.05	ug/g	ND	86.2	60-130			
Dichlorodifluoromethane	4.12	0.05	ug/g	ND	103	50-140			
1,2-Dichlorobenzene	3.67	0.05	ug/g	ND	91.7	60-130			
1,3-Dichlorobenzene	3.53	0.05	ug/g	ND	88.2	60-130			
1,4-Dichlorobenzene	3.40	0.05	ug/g	ND	85.0	60-130			
1,1-Dichloroethane	4.12	0.05	ug/g	ND	103	60-130			
1,2-Dichloroethane	4.56	0.05	ug/g	ND	114	60-130			
1,1-Dichloroethylene	4.09	0.05	ug/g	ND	102	60-130			
cis-1,2-Dichloroethylene	3.87	0.05	ug/g	ND	96.6	60-130			
trans-1,2-Dichloroethylene	3.73	0.05	ug/g	ND	93.2	60-130			
1,2-Dichloropropane	4.53	0.05	ug/g	ND	113	60-130			
cis-1,3-Dichloropropylene	2.91	0.05	ug/g	ND	72.7	60-130			
trans-1,3-Dichloropropylene	2.62	0.05	ug/g	ND	65.5	60-130			
Ethylbenzene	4.38	0.05	ug/g	ND	110	60-130			
Ethylene dibromide (dibromoethane, 1,2	3.45	0.05	ug/g	ND	86.2	60-130			
Hexane	3.68	0.05	ug/g	ND	92.0	60-130			
Methyl Ethyl Ketone (2-Butanone)	12.7	0.50	ug/g	ND	127	50-140			
Methyl Isobutyl Ketone	12.1	0.50	ug/g	ND	121	50-140			
Methyl tert-butyl ether	12.2	0.05	ug/g	ND	122	50-140			
Methylene Chloride	4.11	0.05	ug/g	ND	103	60-130			
Styrene	3.42	0.05	ug/g	ND	85.6	60-130			
1,1,1,2-Tetrachloroethane	3.35	0.05	ug/g	ND	83.8	60-130			
1,1,2,2-Tetrachloroethane	3.34	0.05	ug/g	ND	83.4	60-130			
Tetrachloroethylene	3.73	0.05	ug/g	ND	93.2	60-130			
Toluene	4.42	0.05	ug/g	ND	110	60-130			
1,1,1-Trichloroethane	3.44	0.05	ug/g	ND	86.1	60-130			
1,1,2-Trichloroethane	3.93	0.05	ug/g	ND	98.3	60-130			
Trichloroethylene	3.83	0.05	ug/g	ND	95.6	60-130			
Trichlorofluoromethane	3.82	0.05	ug/g	ND	95.4	50-140			
Vinyl chloride	2.89	0.02	ug/g	ND	72.3	50-140			
m,p-Xylenes	8.25	0.05	ug/g	ND	103	60-130			
o-Xylene	4.24	0.05	ug/g	ND	106	60-130			
Surrogate: 4-Bromofluorobenzene	8.35		ug/g		104	50-140			
Surrogate: Dibromofluoromethane	8.33		ug/g		104	50-140			
Surrogate: Toluene-d8	8.58		ug/g		107	50-140			



Client: Paterson Group Consulting Engineers

Order #: 2309081

Report Date: 02-Mar-2023 Order Date: 27-Feb-2023

Client PO: 56893 Project Description: PE2709

Qualifier Notes:

QC Qualifiers:

Certificate of Analysis

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

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.

Chain of Custody (Rlank) visy



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Paracel Order Number (Lab Use Only)

Chain Of Custody (Lab Use Only)

7269081

LABORATORIES LTD.					com	Y 2	090	01						
Contact Name: Paterson Group Force Contact Name: Samuel Borache Karyn Mu Address: 9 Auriza Drive, ottawa, On Telephone: (12 days 7200)		Project	Ref:	PEQ	109							Page	1_of _	1
Contact Name: Samuel Borathe Karun Mu	ndo	Quote	#:								•	urnaro	und Tim	e
Address:	100	PO#:	V	56893							□ 1 day			☐ 3 day
9 Aurisa Drive attawa OA	/	E-mail:	Sh	embe op	vtersonsr	eup	٠٥٤				□ 2 day			Regular
Telephone: 613 - 226 - 7381			tm	enise ope unch opate	exanavou) · ca				Da	te Requ	ired:		
REG 153/04 REG 406/19 Other Regulation				(Soil/Sed.) GW (Gr						Requi	ed Ana	veie		
☐ Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558 ☐ PWQO			face W	/ater) SS (Storm/San	itary Sewer)	included the state of the state								
☐ Table 2 ☐ Ind/Comm ☐ Coarse ☐ CCME ☐ MISA			P (P	aint) A (Air) O (Oth	er)	JEX								
Table 3 ☐ Agri/Other ☐ SU-Sani ☐ SU-Storm			ers			F1-F4+BTEX			G G					
☐ Table Mun:		ıme	Containers	Sample	Taken	F1-F			by ICP		(S)			
For RSC: Yes No Other:	Matrix	۸ir Volume	of Co			PHCs	VOCs	PAHs	Metals	BH 3	(HWS)			
Sample ID/Location Name	_	Air	#	Date	Time	<u>a</u>	š	1	2/	P 3	<u></u>		+	
1 BH1-23-SS2	5			Feb. 24123		 	1/	V	+	-	+		+	
2 BH1-23-554			2			V	V		_	_			+	
3 BH1-23-555 (hold)	1		2	1					_	_				
4										_				
5												Ш	\perp	
6														
7														
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Comments: * Please place BH1-23-555 on	ho	ıd			Q ₀		2			8	Delivery	P	+	
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Date/Time: Temperature:				°C	Temperature:	13	6		1	H Verifie	d: 🗆	By:		



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

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive Ottawa, ON K2E 7T9 Attn: Sam Berube

Client PO: 56931 Project: PE2709

Custody:

Report Date: 8-Mar-2023 Order Date: 3-Mar-2023

Order #: 2309472

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

Paracel ID	Client ID
2309472-01	BH2-23-SS5
2309472-03	BH3-23-SS5
2309472-05	BH4-23-SS3
2309472-06	BH6-23-SS2
2309472-07	DUP1-23

Approved By:

Mark Froto

Mark Foto, M.Sc. Lab Supervisor



Order #: 2309472

Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

Client PO: 56931 Project Description: PE2709

Analysis Summary Table

Client: Paterson Group Consulting Engineers

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	6-Mar-23	7-Mar-23
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	7-Mar-23	7-Mar-23
Conductivity	MOE E3138 - probe @25 °C, water ext	7-Mar-23	7-Mar-23
Mercury by CVAA	EPA 7471B - CVAA, digestion	7-Mar-23	7-Mar-23
pH, soil	EPA 150.1 - pH probe @ 25 °C, CaCl buffered ext.	6-Mar-23	7-Mar-23
PHC F1	CWS Tier 1 - P&T GC-FID	6-Mar-23	7-Mar-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	6-Mar-23	7-Mar-23
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	7-Mar-23	7-Mar-23
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	3-Mar-23	6-Mar-23
REG 153: VOCs by P&T GC/MS	EPA 8260 - P&T GC-MS	6-Mar-23	7-Mar-23
SAR	Calculated	7-Mar-23	7-Mar-23
Solids, %	CWS Tier 1 - Gravimetric	6-Mar-23	6-Mar-23



Order #: 2309472

Report Date: 08-Mar-2023

Order Date: 3-Mar-2023

Client PO: 56931

Client: Paterson Group Consulting Engineers

Project Description: PE2709

	Client ID: Sample Date: Sample ID: MDL/Units	BH2-23-SS5 27-Feb-23 00:00 2309472-01 Soil	BH3-23-SS5 27-Feb-23 00:00 2309472-03 Soil	BH4-23-SS3 28-Feb-23 00:00 2309472-05 Soil	BH6-23-SS2 01-Mar-23 00:00 2309472-06 Soil
Physical Characteristics	MDL/OIIIIS				00
% Solids	0.1 % by Wt.	92.6	92.7	92.5	96.5
General Inorganics	+		-		
SAR	0.01 N/A	-	-	4.39	4.65
Conductivity	5 uS/cm	-	-	1160	1770
рН	0.05 pH Units	-	-	7.59	7.89
Metals			•		
Antimony	1.0 ug/g dry	-	-	<1.0	<1.0
Arsenic	1.0 ug/g dry	-	-	3.4	2.2
Barium	1.0 ug/g dry	-	-	31.3	211
Beryllium	0.5 ug/g dry	-	-	<0.5	<0.5
Boron	5.0 ug/g dry	-	-	6.8	16.2
Cadmium	0.5 ug/g dry	-	-	<0.5	<0.5
Chromium	5.0 ug/g dry	-	-	13.0	11.0
Chromium (VI)	0.2 ug/g dry	-	-	<0.2	<0.2
Cobalt	1.0 ug/g dry	-	_	3.3	5.5
Copper	5.0 ug/g dry	-	-	6.1	7.2
Lead	1.0 ug/g dry	-	-	12.3	5.8
Mercury	0.1 ug/g dry	-	_	<0.1	<0.1
Molybdenum	1.0 ug/g dry	-	-	<1.0	1.0
Nickel	5.0 ug/g dry	-	-	7.5	10.1
Selenium	1.0 ug/g dry	-	-	<1.0	<1.0
Silver	0.3 ug/g dry	-	-	<0.3	<0.3
Thallium	1.0 ug/g dry	-	-	<1.0	<1.0
Uranium	1.0 ug/g dry	-	-	<1.0	<1.0
Vanadium	10.0 ug/g dry	-	-	18.9	<10.0
Zinc	20.0 ug/g dry	-	-	32.1	<20.0
Volatiles			1		
Acetone	0.50 ug/g dry	<0.50	<0.50	-	-
Benzene	0.02 ug/g dry	<0.02	<0.02	-	-
Bromodichloromethane	0.05 ug/g dry	<0.05	<0.05	-	-
Bromoform	0.05 ug/g dry	<0.05	<0.05	-	-
Bromomethane	0.05 ug/g dry	<0.05	<0.05	-	-
Carbon Tetrachloride	0.05 ug/g dry	<0.05	<0.05	-	-
Chlorobenzene	0.05 ug/g dry	<0.05	<0.05	-	-
Chloroform	0.05 ug/g dry	<0.05	<0.05	-	-
	- +		+	ļ	Į



Client: Paterson Group Consulting Engineers

Certificate of Analysis

Order #: 2309472

Report Date: 08-Mar-2023

Order Date: 3-Mar-2023

Client PO: 56931 Project Description: PE2709

ŗ	Client ID: Sample Date: Sample ID:	BH2-23-SS5 27-Feb-23 00:00 2309472-01	BH3-23-SS5 27-Feb-23 00:00 2309472-03	BH4-23-SS3 28-Feb-23 00:00 2309472-05	BH6-23-SS2 01-Mar-23 00:00 2309472-06
Dibromochloromethane	MDL/Units 0.05 ug/g dry	Soil	Soil	Soil	Soil
Dichlorodifluoromethane	0.05 ug/g dry	<0.05	<0.05	-	-
	0.05 ug/g dry	<0.05	<0.05	-	-
1,2-Dichlorobenzene		<0.05	<0.05	-	-
1,3-Dichlorobenzene	0.05 ug/g dry	<0.05	<0.05	-	-
1,4-Dichlorobenzene	0.05 ug/g dry	<0.05	<0.05	-	-
1,1-Dichloroethane	0.05 ug/g dry	<0.05	<0.05	-	-
1,2-Dichloroethane	0.05 ug/g dry	<0.05	<0.05	-	-
1,1-Dichloroethylene	0.05 ug/g dry	<0.05	<0.05	-	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	<0.05	-	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	<0.05	-	-
1,2-Dichloropropane	0.05 ug/g dry	<0.05	<0.05	-	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	<0.05	-	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	<0.05	-	-
1,3-Dichloropropene, total	0.05 ug/g dry	<0.05	<0.05	-	-
Ethylbenzene	0.05 ug/g dry	<0.05	<0.05	-	-
Ethylene dibromide (dibromoethane, 1,2-)	0.05 ug/g dry	<0.05	<0.05	-	-
Hexane	0.05 ug/g dry	<0.05	<0.05	-	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	<0.50	<0.50	-	-
Methyl Isobutyl Ketone	0.50 ug/g dry	<0.50	<0.50	-	-
Methyl tert-butyl ether	0.05 ug/g dry	<0.05	<0.05	-	-
Methylene Chloride	0.05 ug/g dry	<0.05	<0.05	-	-
Styrene	0.05 ug/g dry	<0.05	<0.05	-	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	<0.05	<0.05	-	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	<0.05	<0.05	-	-
Tetrachloroethylene	0.05 ug/g dry	<0.05	<0.05	-	-
Toluene	0.05 ug/g dry	<0.05	<0.05	-	-
1,1,1-Trichloroethane	0.05 ug/g dry	<0.05	<0.05	-	-
1,1,2-Trichloroethane	0.05 ug/g dry	<0.05	<0.05	-	-
Trichloroethylene	0.05 ug/g dry	<0.05	<0.05	-	-
Trichlorofluoromethane	0.05 ug/g dry	<0.05	<0.05	-	-
Vinyl chloride	0.02 ug/g dry	<0.02	<0.02	-	-
m,p-Xylenes	0.05 ug/g dry	<0.05	<0.05	_	-
o-Xylene	0.05 ug/g dry	<0.05	<0.05	-	_
Xylenes, total	0.05 ug/g dry	<0.05	<0.05	_	-
4-Bromofluorobenzene	Surrogate	104%	103%	-	-
Dibromofluoromethane	Surrogate	113%	113%	-	-



Order #: 2309472

Report Date: 08-Mar-2023

Order Date: 3-Mar-2023

Client: Paterson Group Consulting Engineers

Client PO: 56931 **Project Description: PE2709**

	Client ID: Sample Date: Sample ID:	BH2-23-SS5 27-Feb-23 00:00 2309472-01	BH3-23-SS5 27-Feb-23 00:00 2309472-03	BH4-23-SS3 28-Feb-23 00:00 2309472-05	BH6-23-SS2 01-Mar-23 00:00 2309472-06
Toluene-d8	MDL/Units Surrogate	Soil 119%	Soil 119%	Soil -	Soil -
Benzene	0.02 ug/g dry	-	-	<0.02	<0.02
Ethylbenzene	0.05 ug/g dry	<u> </u>	_	<0.05	<0.05
Toluene	0.05 ug/g dry		_	<0.05	<0.05
m,p-Xylenes	0.05 ug/g dry	<u>-</u>		<0.05	<0.05
o-Xylene	0.05 ug/g dry		-	<0.05	<0.05
Xylenes, total	0.05 ug/g dry	-	-	<0.05	<0.05
Toluene-d8	Surrogate	-	-	120%	115%
Hydrocarbons			ļ	12070	1.070
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	50
F4 PHCs (C34-C50)	6 ug/g dry	<6	-	<6	42
Semi-Volatiles		-			
Acenaphthene	0.02 ug/g dry	-	-	<0.02	<0.02
Acenaphthylene	0.02 ug/g dry	-	-	<0.02	<0.02
Anthracene	0.02 ug/g dry	-	-	<0.02	<0.02
Benzo [a] anthracene	0.02 ug/g dry	-	-	<0.02	<0.02
Benzo [a] pyrene	0.02 ug/g dry	-	-	<0.02	<0.02
Benzo [b] fluoranthene	0.02 ug/g dry	-	-	<0.02	<0.02
Benzo [g,h,i] perylene	0.02 ug/g dry	-	-	<0.02	<0.02
Benzo [k] fluoranthene	0.02 ug/g dry	-	-	<0.02	<0.02
Chrysene	0.02 ug/g dry	-	-	<0.02	<0.02
Dibenzo [a,h] anthracene	0.02 ug/g dry	-	-	<0.02	<0.02
Fluoranthene	0.02 ug/g dry	-	-	<0.02	<0.02
Fluorene	0.02 ug/g dry	-	-	<0.02	<0.02
Indeno [1,2,3-cd] pyrene	0.02 ug/g dry	-	-	<0.02	<0.02
1-Methylnaphthalene	0.02 ug/g dry	-	-	<0.02	<0.02
2-Methylnaphthalene	0.02 ug/g dry	-	-	<0.02	<0.02
Methylnaphthalene (1&2)	0.04 ug/g dry	-	-	<0.04	<0.04
Naphthalene	0.01 ug/g dry	-	-	<0.01	<0.01
Phenanthrene	0.02 ug/g dry	-	-	<0.02	<0.02
Pyrene	0.02 ug/g dry	-	_	<0.02	<0.02
2-Fluorobiphenyl	Surrogate	-	-	81.7%	105%
Terphenyl-d14	Surrogate	-	-	119%	132%



Order #: 2309472

Report Date: 08-Mar-2023

Order Date: 3-Mar-2023

Client: Paterson Group Consulting Engineers Client PO: 56931

Project Description: PE2709

	Client ID	DUP1-23			
	Client ID: Sample Date:	27-Feb-23 00:00	-	-	-
	Sample ID:	2309472-07	-	-	-
	MDL/Units	Soil	-	-	-
Physical Characteristics					1
% Solids	0.1 % by Wt.	93.4	-	-	-
Volatiles	0.50 / 1		1		1
Acetone	0.50 ug/g dry	<0.50	-	-	-
Benzene	0.02 ug/g dry	<0.02	-	-	-
Bromodichloromethane	0.05 ug/g dry	<0.05	-	-	-
Bromoform	0.05 ug/g dry	<0.05	-	-	-
Bromomethane	0.05 ug/g dry	<0.05	-	-	-
Carbon Tetrachloride	0.05 ug/g dry	<0.05	-	-	-
Chlorobenzene	0.05 ug/g dry	<0.05	-	-	-
Chloroform	0.05 ug/g dry	<0.05	-	-	-
Dibromochloromethane	0.05 ug/g dry	<0.05	-	-	-
Dichlorodifluoromethane	0.05 ug/g dry	<0.05	-	-	-
1,2-Dichlorobenzene	0.05 ug/g dry	<0.05	-	-	-
1,3-Dichlorobenzene	0.05 ug/g dry	<0.05	-	-	-
1,4-Dichlorobenzene	0.05 ug/g dry	<0.05	-	-	-
1,1-Dichloroethane	0.05 ug/g dry	<0.05	-	-	-
1,2-Dichloroethane	0.05 ug/g dry	<0.05	-	-	-
1,1-Dichloroethylene	0.05 ug/g dry	<0.05	-	-	-
cis-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	-	-	-
trans-1,2-Dichloroethylene	0.05 ug/g dry	<0.05	-	-	-
1,2-Dichloropropane	0.05 ug/g dry	<0.05	-	-	-
cis-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	-	-	-
trans-1,3-Dichloropropylene	0.05 ug/g dry	<0.05	-	-	-
1,3-Dichloropropene, total	0.05 ug/g dry	<0.05	-	-	-
Ethylbenzene	0.05 ug/g dry	<0.05	-	-	-
Ethylene dibromide (dibromoethane, 1	0.05 ug/g dry	<0.05	-	-	-
Hexane	0.05 ug/g dry	<0.05	-	-	-
Methyl Ethyl Ketone (2-Butanone)	0.50 ug/g dry	<0.50	-	-	-
Methyl Isobutyl Ketone	0.50 ug/g dry	<0.50	-	-	-
Methyl tert-butyl ether	0.05 ug/g dry	<0.05	-	-	-
Methylene Chloride	0.05 ug/g dry	<0.05	-	-	-
Styrene	0.05 ug/g dry	<0.05	-	-	-
1,1,1,2-Tetrachloroethane	0.05 ug/g dry	<0.05	-	-	-
1,1,2,2-Tetrachloroethane	0.05 ug/g dry	<0.05	-	-	-



Client: Paterson Group Consulting Engineers

Certificate of Analysis

Order #: 2309472

Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

Client PO: 56931 Project Description: PE2709

	Client ID:	DUP1-23	-	-	-
	Sample Date:	27-Feb-23 00:00	-	-	-
	Sample ID:	2309472-07	-	-	-
	MDL/Units	Soil	-	-	-
Tetrachloroethylene	0.05 ug/g dry	<0.05	-	-	-
Toluene	0.05 ug/g dry	<0.05	-	-	-
1,1,1-Trichloroethane	0.05 ug/g dry	<0.05	-	-	-
1,1,2-Trichloroethane	0.05 ug/g dry	<0.05	-	-	-
Trichloroethylene	0.05 ug/g dry	<0.05	-	-	-
Trichlorofluoromethane	0.05 ug/g dry	<0.05	-	-	-
Vinyl chloride	0.02 ug/g dry	<0.02	-	-	-
m,p-Xylenes	0.05 ug/g dry	<0.05	-	-	-
o-Xylene	0.05 ug/g dry	<0.05	-	-	-
Xylenes, total	0.05 ug/g dry	<0.05	-	-	-
4-Bromofluorobenzene	Surrogate	99.6%	-	-	-
Dibromofluoromethane	Surrogate	111%	-	-	-
Toluene-d8	Surrogate	114%	-	-	-



Order #: 2309472

Report Date: 08-Mar-2023

Order Date: 3-Mar-2023

Client: Paterson Group Consulting Engineers Client PO: 56931 **Project Description: PE2709**

Method Quality Control: Blank

Result	Reporting Limit	Units	Source Result	%RFC	%REC	RPD	RPD Limit	Notes
, , , , ,	Liiill	Office	rtesuit	701 NEO	Liillit	111 D	Liitiit	
	_	0.4						
ND	5	uS/cm						
	7	ug/g						
ND	O	ug/g						
ND	4.0							
ND	0.5							
ND	5.0							
ND	0.5	ug/g						
ND	0.2	ug/g						
ND	5.0	ug/g						
ND	5.0							
ND	1.0	ug/g						
ND	0.3	ug/g						
ND	1.0	ug/g						
ND	1.0	ug/g						
ND	∠0.0	ug/g						
NB	0.00							
ND	0.02							
ND	0.02	ug/g						
ND	0.02	ug/g						
ND	0.02	ug/g						
ND	0.02	ug/g						
		ug/g						
ND	0.04	ug/g						
ND	0.01	ug/g						
ND	0.02	ug/g						
	0.02	ug/g			# * * * * *			
1.54		ug/g		115	50-140			
ND	0.50	ug/g						
ND		ug/g						
ND ND	0.05	ug/g ug/g						
שמו	0.00	ug/g						
	ND N	ND 5	ND 5	ND S US/cm	ND 5	Result Limit Units Result %REC Limit	ND 5	Result Limit Units Result %REC Limit RPD Limit



Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

Project Description: PE2709

Certificate of Analysis

Client PO: 56931

Client: Paterson Group Consulting Engineers

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Chloroform	ND	0.05	ug/g						
Dibromochloromethane	ND	0.05	ug/g						
Dichlorodifluoromethane	ND	0.05	ug/g						
1,2-Dichlorobenzene	ND	0.05	ug/g						
1,3-Dichlorobenzene	ND	0.05	ug/g						
1,4-Dichlorobenzene	ND	0.05	ug/g						
1,1-Dichloroethane	ND	0.05	ug/g						
1,2-Dichloroethane	ND	0.05	ug/g						
1,1-Dichloroethylene	ND	0.05	ug/g						
cis-1,2-Dichloroethylene	ND	0.05	ug/g						
trans-1,2-Dichloroethylene	ND	0.05	ug/g						
1,2-Dichloropropane	ND	0.05	ug/g						
cis-1,3-Dichloropropylene	ND	0.05	ug/g						
trans-1,3-Dichloropropylene	ND	0.05	ug/g						
1,3-Dichloropropene, total	ND	0.05	ug/g						
Ethylbenzene	ND	0.05	ug/g						
Ethylene dibromide (dibromoethane, 1,2	ND	0.05	ug/g						
Hexane	ND	0.05	ug/g						
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g						
Methyl Isobutyl Ketone	ND	0.50	ug/g						
Methyl tert-butyl ether	ND	0.05	ug/g						
Methylene Chloride	ND	0.05	ug/g						
Styrene	ND	0.05	ug/g						
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g						
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g						
Tetrachloroethylene	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
1,1,1-Trichloroethane	ND	0.05	ug/g						
1,1,2-Trichloroethane	ND	0.05	ug/g						
Trichloroethylene	ND	0.05	ug/g						
Trichlorofluoromethane	ND	0.05	ug/g						
Vinyl chloride	ND	0.02	ug/g						
m,p-Xylenes	ND	0.02	ug/g ug/g						
o-Xylene	ND	0.05	ug/g ug/g						
Xylenes, total	ND ND	0.05	ug/g ug/g						
Surrogate: 4-Bromofluorobenzene	8.90	0.00	ug/g ug/g		111	50-140			
Surrogate: 4-biomonuorobenzene Surrogate: Dibromofluoromethane	8.64				108	50-140 50-140			
			ug/g						
Surrogate: Toluene-d8	9.44	0.05	ug/g		118	50-140			
Benzene	ND	0.02	ug/g						
Ethylbenzene 	ND	0.05	ug/g						
Toluene	ND	0.05	ug/g						
m,p-Xylenes	ND	0.05	ug/g						
o-Xylene	ND	0.05	ug/g						
Xylenes, total	ND	0.05	ug/g						
Surrogate: Toluene-d8	9.44		ug/g		118	50-140			



Order #: 2309472

Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 3-Mar-2023

 Client PO:
 56931
 Project Description: PE2709

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
General Inorganics									
SAR	0.79	0.01	N/A	0.78			1.3	30	
Conductivity	479	5	uS/cm	476			0.6	5	
pH	7.26	0.05	pH Units	7.30			0.5	2.3	
lydrocarbons	0	0.00	pri onno				0.0	2.0	
F1 PHCs (C6-C10)	ND	7	ug/g	ND			NC	40	
F2 PHCs (C10-C16)	321	4	ug/g	317			1.2	30	
F3 PHCs (C16-C34)	7740	8	ug/g	7540			2.5	30	
F4 PHCs (C34-C50)	1310	6	ug/g	1280			2.3	30	
letals									
Antimony	ND	1.0	ug/g	ND			NC	30	
Arsenic	3.6	1.0	ug/g	3.4			6.2	30	
Barium	28.5	1.0	ug/g	31.3			9.6	30	
Beryllium	ND	0.5	ug/g	ND			NC	30	
Boron	5.9	5.0	ug/g	6.8			15.5	30	
Cadmium	ND	0.5	ug/g	ND			NC	30	
Chromium (VI)	ND	0.2	ug/g	ND			NC	35	
Chromium	11.8	5.0	ug/g	13.0			9.8	30	
Cobalt	3.3	1.0	ug/g ug/g	3.3			1.6	30	
Copper	6.0	5.0	ug/g ug/g	6.1			1.8	30	
Lead	12.2	1.0		12.3			1.0	30	
Mercury	12.2 ND	0.1	ug/g ug/g	12.3 ND			NC	30	
-	ND ND	1.0		ND			NC	30	
Molybdenum Nickel	7.2	5.0	ug/g				3.2	30	
Selenium	7.2 ND		ug/g	7.5 ND			NC	30	
		1.0	ug/g						
Silver	ND	0.3	ug/g	ND			NC	30	
Thallium	ND	1.0	ug/g	ND			NC	30	
Uranium	ND	1.0	ug/g	ND			NC	30	
Vanadium	16.7	10.0	ug/g	18.9			12.4	30	
Zinc	30.8	20.0	ug/g	32.1			4.1	30	
Physical Characteristics	00.0	0.4	0/ 5. 344	00.5			0.5	05	
% Solids Semi-Volatiles	92.0	0.1	% by Wt.	92.5			0.5	25	
	ND	0.00		ND			NO	40	
Acenaphthene	ND	0.02	ug/g	ND			NC	40	
Acenaphthylene	ND	0.02	ug/g	ND			NC	40	
Anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] pyrene	ND	0.02	ug/g	ND			NC	40	
Benzo [b] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Benzo [g,h,i] perylene	ND	0.02	ug/g	ND			NC	40	
Benzo [k] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Chrysene	ND	0.02	ug/g	ND			NC	40	
Dibenzo [a,h] anthracene	ND	0.02	ug/g	ND			NC	40	
Fluoranthene	ND	0.02	ug/g	ND			NC	40	
Fluorene	ND	0.02	ug/g	ND			NC	40	
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g	ND			NC	40	
1-Methylnaphthalene	ND	0.02	ug/g	ND			NC	40	
2-Methylnaphthalene	ND	0.02	ug/g	ND			NC	40	
Naphthalene	ND	0.01	ug/g	ND			NC	40	
Phenanthrene	ND	0.02	ug/g	ND			NC	40	
Pyrene	ND	0.02	ug/g	ND			NC	40	
Surrogate: 2-Fluorobiphenyl	0.885		ug/g		51.7	50-140			
Surrogate: Terphenyl-d14	1.16		ug/g		68.0	50-140			
/olatiles									
Acetone	ND	0.50	ug/g	ND			NC	50	
Benzene	ND	0.02	ug/g	ND			NC	50	



Order #: 2309472

Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

Client PO: 56931 Project Description: PE2709

Method Quality Control: Duplicate

Client: Paterson Group Consulting Engineers

Analyte	D	Reporting		Source	0/5==	%REC	000	RPD	NI. 4.
nialyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Bromodichloromethane	ND	0.05	ug/g	ND			NC	50	
Bromoform	ND	0.05	ug/g	ND			NC	50	
Bromomethane	ND	0.05	ug/g	ND			NC	50	
Carbon Tetrachloride	ND	0.05	ug/g	ND			NC	50	
Chlorobenzene	ND	0.05	ug/g	ND			NC	50	
Chloroform	ND	0.05	ug/g	ND			NC	50	
Dibromochloromethane	ND	0.05	ug/g	ND			NC	50	
Dichlorodifluoromethane	ND	0.05	ug/g	ND			NC	50	
1,2-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,3-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,4-Dichlorobenzene	ND	0.05	ug/g	ND			NC	50	
1,1-Dichloroethane	ND	0.05	ug/g	ND			NC	50	
1,2-Dichloroethane	ND	0.05	ug/g	ND			NC	50	
1,1-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
cis-1,2-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
trans-1,2-Dichloroethylene	ND	0.05	ug/g	ND			NC	50	
1,2-Dichloropropane	ND	0.05	ug/g	ND			NC	50	
cis-1,3-Dichloropropylene	ND	0.05	ug/g	ND			NC	50	
trans-1,3-Dichloropropylene	ND	0.05	ug/g	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g	ND			NC	50	
Ethylene dibromide (dibromoethane, 1,2	ND	0.05	ug/g	ND			NC	50	
Hexane	ND	0.05	ug/g	ND			NC	50	
Methyl Ethyl Ketone (2-Butanone)	ND	0.50	ug/g	ND			NC	50	
Methyl Isobutyl Ketone	ND	0.50	ug/g	ND			NC	50	
Methyl tert-butyl ether	ND	0.05	ug/g	ND			NC	50	
Methylene Chloride	ND	0.05	ug/g	ND			NC	50	
Styrene	ND	0.05	ug/g	ND			NC	50	
1,1,1,2-Tetrachloroethane	ND	0.05	ug/g	ND			NC	50	
1,1,2,2-Tetrachloroethane	ND	0.05	ug/g	ND			NC	50	
Tetrachloroethylene	ND	0.05	ug/g	ND			NC	50	
Toluene	ND	0.05	ug/g	ND			NC	50	
1,1,1-Trichloroethane	ND	0.05	ug/g	ND			NC	50	
1,1,2-Trichloroethane	ND	0.05	ug/g	ND			NC	50	
Trichloroethylene	ND	0.05	ug/g	ND			NC	50	
Trichlorofluoromethane	ND	0.05	ug/g	ND			NC	50	
Vinyl chloride	ND	0.02	ug/g	ND			NC	50	
m,p-Xylenes	ND	0.05	ug/g	ND			NC	50	
o-Xylene	ND	0.05	ug/g	ND			NC	50	
Surrogate: 4-Bromofluorobenzene	9.94		ug/g		106	50-140			
Surrogate: Dibromofluoromethane	10.9		ug/g		115	50-140			
Surrogate: Toluene-d8	11.3		ug/g		120	50-140			
Benzene	ND	0.02	ug/g	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g	ND			NC	50	
Toluene	ND	0.05	ug/g	ND			NC	50	
m,p-Xylenes	ND	0.05	ug/g	ND			NC	50	
o-Xylene	ND	0.05	ug/g	ND			NC	50	
Surrogate: Toluene-d8	11.3	0.00	ug/g		120	50-140			



Order #: 2309472

Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 3-Mar-2023

 Client PO:
 56931
 Project Description: PE2709

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	160	7	ug/g	ND	80.1	80-120			
F2 PHCs (C10-C16)	433	4	ug/g	317	123	60-140			
F3 PHCs (C16-C34)	233	8	ug/g	ND	119	80-120			
F4 PHCs (C34-C50)	121	6	ug/g	ND	97.6	80-120			
Metals									
Antimony	35.2	1.0	ug/g	ND	70.1	70-130			
Arsenic	47.7	1.0	ug/g	1.3	92.7	70-130			
Barium	55.0	1.0	ug/g ug/g	12.5	85.0	70-130			
Beryllium	45.7	0.5	ug/g ug/g	ND	91.0	70-130			
Boron	44.3	5.0	ug/g ug/g	ND	83.1	70-130			
Cadmium	45.2	0.5	ug/g ug/g	ND	90.3	70-130			
Chromium (VI)	4.4	0.2	ug/g ug/g	ND	81.5	70-130			
Chromium	54.1	5.0	ug/g ug/g	5.2	97.7	70-130			
Cobalt	48.0	1.0	ug/g ug/g	1.3	93.3	70-130			
Copper	46.3	5.0	ug/g ug/g	ND	93.3 87.7	70-130			
Lead	40.5 47.5	1.0	ug/g ug/g	4.9	85.1	70-130			
Mercury	47.5 1.15	0.1	ug/g ug/g	4.9 ND	76.4	70-130 70-130			
Molybdenum	49.3	1.0		ND	98.2	70-130			
Nickel	49.3	5.0	ug/g	ND	92.1	70-130			
Selenium	49.0		ug/g			70-130			
	42.3 42.2	1.0	ug/g	ND	84.3	70-130 70-130			
Silver		0.3	ug/g	ND	84.3				
Thallium	43.4	1.0	ug/g	ND	86.7	70-130			
Uranium	44.1	1.0	ug/g	ND	87.8	70-130			
Vanadium Zinc	55.5 53.3	10.0 20.0	ug/g	ND ND	95.8 81.0	70-130 70-130			
	55.5	20.0	ug/g	ND	01.0	70-130			
Semi-Volatiles									
Acenaphthene	0.142	0.02	ug/g	ND	85.4	50-140			
Acenaphthylene	0.119	0.02	ug/g	ND	71.5	50-140			
Anthracene	0.120	0.02	ug/g	ND	72.3	50-140			
Benzo [a] anthracene	0.100	0.02	ug/g	ND	60.2	50-140			
Benzo [a] pyrene	0.118	0.02	ug/g	ND	70.9	50-140			
Benzo [b] fluoranthene	0.144	0.02	ug/g	ND	86.6	50-140			
Benzo [g,h,i] perylene	0.124	0.02	ug/g	ND	74.5	50-140			
Benzo [k] fluoranthene	0.137	0.02	ug/g	ND	82.2	50-140			
Chrysene	0.130	0.02	ug/g	ND	78.0	50-140			
Dibenzo [a,h] anthracene	0.121	0.02	ug/g	ND	72.5	50-140			
Fluoranthene	0.113	0.02	ug/g	ND	67.9	50-140			
Fluorene	0.126	0.02	ug/g	ND	75.5	50-140			
Indeno [1,2,3-cd] pyrene	0.121	0.02	ug/g	ND	72.6	50-140			
1-Methylnaphthalene	0.133	0.02	ug/g	ND	79.8	50-140			
2-Methylnaphthalene	0.146	0.02	ug/g	ND	87.4	50-140			
Naphthalene	0.158	0.01	ug/g	ND	94.9	50-140			
Phenanthrene	0.124	0.02	ug/g	ND	74.4	50-140			
Pyrene	0.113	0.02	ug/g	ND	68.1	50-140			
Surrogate: 2-Fluorobiphenyl	1.14		ug/g		85.8	50-140			
Surrogate: Terphenyl-d14	1.58		ug/g		118	50-140			
/olatiles									
Acetone	13.0	0.50	ug/g	ND	130	50-140			



Order #: 2309472

Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 3-Mar-2023

 Client PO:
 56931
 Project Description: PE2709

Method Quality Control: Spike

nalyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Benzene	4.86	0.02	ug/g	ND	121	60-130			
Bromodichloromethane	4.94	0.05	ug/g	ND	124	60-130			
Bromoform	4.92	0.05	ug/g	ND	123	60-130			
Bromomethane	4.10	0.05	ug/g	ND	102	50-140			
Carbon Tetrachloride	4.62	0.05	ug/g	ND	115	60-130			
Chlorobenzene	4.61	0.05	ug/g	ND	115	60-130			
Chloroform	4.83	0.05	ug/g	ND	121	60-130			
Dibromochloromethane	4.93	0.05	ug/g	ND	123	60-130			
Dichlorodifluoromethane	3.55	0.05	ug/g	ND	88.7	50-140			
l,2-Dichlorobenzene	4.36	0.05	ug/g	ND	109	60-130			
,3-Dichlorobenzene	4.31	0.05	ug/g	ND	108	60-130			
l,4-Dichlorobenzene	4.14	0.05	ug/g	ND	103	60-130			
l,1-Dichloroethane	4.90	0.05	ug/g	ND	122	60-130			
l,2-Dichloroethane	4.82	0.05	ug/g	ND	121	60-130			
, 1,1-Dichloroethylene	4.86	0.05	ug/g	ND	122	60-130			
sis-1,2-Dichloroethylene	4.79	0.05	ug/g	ND	120	60-130			
rans-1,2-Dichloroethylene	4.79	0.05	ug/g	ND	120	60-130			
I,2-Dichloropropane	4.75	0.05	ug/g	ND	119	60-130			
sis-1,3-Dichloropropylene	4.76	0.05	ug/g	ND	119	60-130			
rans-1,3-Dichloropropylene	4.25	0.05	ug/g	ND	106	60-130			
Ethylbenzene	4.64	0.05	ug/g	ND	116	60-130			
Ethylene dibromide (dibromoethane, 1,2	4.88	0.05	ug/g	ND	122	60-130			
Hexane	4.47	0.05	ug/g	ND	112	60-130			
Methyl Ethyl Ketone (2-Butanone)	13.0	0.50	ug/g	ND	130	50-140			
Methyl Isobutyl Ketone	12.7	0.50	ug/g	ND	127	50-140			
Methyl tert-butyl ether	9.87	0.05	ug/g	ND	98.7	50-140			
Methylene Chloride	4.68	0.05	ug/g	ND	117	60-130			
Styrene	4.90	0.05	ug/g ug/g	ND	122	60-130			
I,1,1,2-Tetrachloroethane	4.57	0.05	ug/g ug/g	ND	114	60-130			
1,1,2,2-Tetrachloroethane	4.65	0.05	ug/g ug/g	ND	116	60-130			
Fetrachloroethylene	4.74	0.05	ug/g ug/g	ND	118	60-130			
Foluene	4.60	0.05	ug/g ug/g	ND	115	60-130			
,1,1-Trichloroethane	4.71	0.05	ug/g ug/g	ND	118	60-130			
1,1,2-Trichloroethane	4.31	0.05		ND	108	60-130			
Frichloroethylene	4.57	0.05	ug/g ug/g	ND	114	60-130			
Frichlorofluoromethane	4.18	0.05		ND	105	50-130			
	3.87	0.03	ug/g	ND					
/inyl chloride			ug/g		96.8	50-140 60 130			
n,p-Xylenes o-Xylene	9.41 4.88	0.05 0.05	ug/g	ND ND	118 122	60-130 60-130			
-	4.88 7.86	0.05	ug/g	ND	98.2	50-130 50-140			
Surrogate: 4-Bromofluorobenzene Surrogate: Dibromofluoromethane	7.86 8.34		ug/g ug/g		98.2 104	50-140 50-140			
Surrogate: Dibromonuorometriane Surrogate: Toluene-d8	7.72		ug/g ug/g		96.5	50-140 50-140			
Benzene	4.86	0.02	ug/g ug/g	ND	121	60-130			
Ethylbenzene	4.64	0.02	ug/g ug/g	ND	116	60-130			
Foluene	4.60	0.05		ND	115	60-130			
n,p-Xylenes	4.60 9.41	0.05	ug/g	ND	118	60-130			
•			ug/g						
o-Xylene Surrogate: Toluene-d8	4.88 7.72	0.05	ug/g <i>ug/g</i>	ND	122 96.5	60-130 <i>50-140</i>			



Report Date: 08-Mar-2023 Order Date: 3-Mar-2023

 Client: Paterson Group Consulting Engineers
 Order Date: 3-Mar-2023

 Client PO: 56931
 Project Description: PE2709

Qualifier Notes:

Sample Data Revisions

Certificate of Analysis

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 III								Paracel Order Number (Lab Use Only)				Chain Of Custody (Lab Use Only)					
Client Name: Sauvet Benke / Lary Contact Name:	Much	ach	Projec	t Ref:	PEZF	ल्य								Pa	ge 📗	of _	
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Telephone: 613.226-7381					Lmunc	happa	Hev	en	CNI	110.	ca	Date	Requ	ired:			
REG 153/04 REG 406/19 Other Regulat	tion	M	atriv T	may	S [Soil/Sed.) GW (Gr	-	888		J.					8/10			
☐ Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558 ☐	PWQO				Vater) SS (Storm/San						Re	quire	d Ana	lysis			
□ Table 2 □ Ind/Comm □ Coarse □ CCME □	MISA			P (P	aint) A (Air) O (Oth	er)	X		Π						П	T	T
☐ Table 3 ☐ Agri/Other ☐ SU-Sani ☐	SU - Storm			23			F1-F4+BTEX			۵					9		
Mun:			ne	Containers	Sample '	Taken	1-F4			by ICP				_	5		
For RSC: Yes No Cther:		Sample Taken O U U U U U U U U U U U U U U U U U U					l «	s			_	(HWS)	#	13			
Sample ID/Location Name		Matrix	Aĭr	# of	Date	Time	PHCs	VOCs	PAHs	Metals	Β̈́	CrVI	B (H	- All	1249		
1 BH2- 23-555		Ş		2	Feb. 2723		1	V									
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300 - 2319 St. Laurent Blvd Ottawa, ON, K1G 4J8 1-800-749-1947 www.paracellabs.com

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive

Ottawa, ON K2E 7T9

Attn: Sam Berube

Client PO: 58143

Project: PE2709

Custody:

Report Date: 18-Aug-2023

Order Date: 15-Aug-2023

Order #: 2333190

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

submitted:

Paracel ID Client ID

2333190-01 BH7-23-SS3

2333190-02 BH8-23-SS6

Das



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58143

Report Date: 18-Aug-2023

Order Date: 15-Aug-2023

Project Description: PE2709

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	16-Aug-23	17-Aug-23
PHC F1	CWS Tier 1 - P&T GC-FID	16-Aug-23	17-Aug-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	15-Aug-23	18-Aug-23
Solids, %	CWS Tier 1 - Gravimetric	17-Aug-23	17-Aug-23

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58143

Report Date: 18-Aug-2023

Order Date: 15-Aug-2023

Project Description: PE2709

	Client ID:	BH7-23-SS3	BH8-23-SS6	-	-		
	Sample Date:	14-Aug-23 09:00	14-Aug-23 09:00	-	-	-	-
	Sample ID:	2333190-01	2333190-02	-	-		
	Matrix:	Soil	Soil	-	-		
	MDL/Units						
Physical Characteristics	•						
% Solids	0.1 % by Wt.	91.8	93.6	-	-	-	-
Volatiles	•					•	
Benzene	0.02 ug/g	<0.02	<0.02	-	=	-	-
Ethylbenzene	0.05 ug/g	<0.05	0.28	-	-	-	-
Toluene	0.05 ug/g	<0.05	0.48	-	-	-	-
m,p-Xylenes	0.05 ug/g	<0.05	0.36	-	-	-	-
o-Xylene	0.05 ug/g	<0.05	<0.05	-	-	-	-
Xylenes, total	0.05 ug/g	<0.05	0.36	-	-	-	-
Toluene-d8	Surrogate	128%	126%	-	-	-	-
Hydrocarbons						-	
F1 PHCs (C6-C10)	7 ug/g	<7	355	-	-	-	-
F2 PHCs (C10-C16)	4 ug/g	<4	377	-	-	-	-
F3 PHCs (C16-C34)	8 ug/g	<8	74	-	-	-	-
F4 PHCs (C34-C50)	6 ug/g	<6	10	-	-	-	-

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58143

Report Date: 18-Aug-2023

Order Date: 15-Aug-2023

Project Description: PE2709

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	%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	9.61		%	120	50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58143

Report Date: 18-Aug-2023

Order Date: 15-Aug-2023

Project Description: PE2709

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									<u> </u>
F1 PHCs (C6-C10)	ND	7	ug/g	ND			NC	40	
F2 PHCs (C10-C16)	55	4	ug/g	48			13.5	30	
F3 PHCs (C16-C34)	76	8	ug/g	54			NC	30	
F4 PHCs (C34-C50)	12	6	ug/g	6			NC	30	
Physical Characteristics									
% Solids	92.2	0.1	% by Wt.	91.8			0.4	25	
Volatiles									
Benzene	ND	0.02	ug/g	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g	ND			NC	50	
Toluene	ND	0.05	ug/g	ND			NC	50	
m,p-Xylenes	ND	0.05	ug/g	ND			NC	50	
o-Xylene	ND	0.05	ug/g	ND			NC	50	
Surrogate: Toluene-d8	11.8		%		129	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58143

Report Date: 18-Aug-2023

Order Date: 15-Aug-2023

Project Description: PE2709

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	176	7	ug/g	ND	87.8	85-115			
F2 PHCs (C10-C16)	163	4	ug/g	48	133	60-140			
F3 PHCs (C16-C34)	310	8	ug/g	54	121	60-140			
F4 PHCs (C34-C50)	173	6	ug/g	6	124	60-140			
Volatiles									
Benzene	3.34	0.02	ug/g	ND	83.4	60-130			
Ethylbenzene	4.70	0.05	ug/g	ND	118	60-130			
Toluene	4.90	0.05	ug/g	ND	123	60-130			
m,p-Xylenes	7.42	0.05	ug/g	ND	92.7	60-130			
o-Xylene	4.50	0.05	ug/g	ND	113	60-130			
Surrogate: Toluene-d8	7.82		%		97.8	50-140			



Client: Paterson Group Consulting Engineers

Order #: 2333190

Report Date: 18-Aug-2023

Order Date: 15-Aug-2023

Project Description: PE2709

Certificate of Analysis

Client PO: 58143

Qualifier Notes:

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 unlesss otherwise noted.

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

Any use of these results implies your agreement that our total liabilty 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.





Paracel Order Number (Lab Use Only)

Chain Of Custody (Lab Use Only)

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Client Name: Paterson Group		Projec	t Ref:	PE 2709									Page	Lof	l and the same of
contact Name: Sam Berube		Quote	#:								Turnaround Time				
Address:		PO #:	58	143									iiiaiou		
9 Auriga Dr Ottawa		E-mail		142								1 day			☐ 3 day
Telephone: 613 - 226 - 7381		Sberube@patersongroup.ca						☐ 2 day			☑ Regular				
REG 153/04 ☐ REG 406/19 Other Regulation		_				- 34	v.)/135	9-35-66	7 300	15 U.C.	SAL	require			
☐ Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558 ☐ PWQO	1 1	Aatrix T SW (Su	ype: : rface V	S (Soil/Sed.) GW (G Vater) SS (Storm/Sa	round Water)					Red	quire	d Analys	is		
☐ Table 2 ☐ Ind/Comm ☐ Coarse ☐ CCME ☐ MISA		, (Sa		aint) A (Air) O (Oth		×		4197	316		Tag	92.5		_	
☐ Table 3 ☐ Agri/Other ☐ SU - Sani ☐ SU - Storm						F1-F4+BTEX									
☐ Table Mun:			iner	Sample	Takon	44			CP						
For RSC: Yes No Other:	×	lum.	Containers	Sample	raken	F1-F			by			(S)			
Sample ID/Location Name	Matrix	Air Volume	# of C	Date	Time	PHCs	VOCs	PAHs	Metals by ICP	_	CrVI	(HWS)			
1 BH7-23-553	5	1	2	Aug 14/23	ilme	_	>	-A	Ž	Ϋ́	ပ်	ω		\vdash	
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Aug 15 2023			23)	°C	Temperature: C	1,90	C	rapply :	100	pH Veri	fied: [17	By:	10	
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300 - 2319 St. Laurent Blvd Ottawa, ON, K1G 4J8 1-800-749-1947 www.paracellabs.com

Certificate of Analysis

Paterson Group Consulting Engineers (Ottawa)

9 Auriga Drive

Ottawa, ON K2E 7T9

Attn: Jeremy Camposarcone

Client PO: 60196

Project: PE6422

Custody:

Report Date: 31-May-2024

Order Date: 14-May-2024

Revised Report Order #: 2420225

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

Paracel ID	Client ID	Paracel ID	Client ID
2420225-01	BH1-24-AU1		
2420225-02	BH1-24-SS3		
2420225-03	BH1-24-SS5		
2420225-04	BH2-24-AU1		
2420225-05	BH2-24-SS3		
2420225-06	BH3-24-AU1		
2420225-07	BH3-24-SS5		
2420225-08	BH4-24-AU1		
2420225-09	BH4-24-SS2		
2420225-10	BH4-24-SS3		
2420225-11	BH4-24-SS6		
2420225-12	BH5-24-AU1		
2420225-13	BH5-24-SS2(BOTTOM)		
2420225-14	DUP1		
2420225-15	DUP2		

Approved By:

Mark Froto

Mark Foto, M.Sc.

Lab Supervisor



Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024 Order Date: 14-May-2024

Project Description: PE6422

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 8260 - P&T GC-MS	16-May-24	16-May-24
Chromium, hexavalent - soil	MOE E3056 - Extraction, colourimetric	21-May-24	21-May-24
Conductivity	MOE E3138 - probe @25 °C, water ext	16-May-24	16-May-24
Mercury by CVAA	EPA 7471B - CVAA, digestion	16-May-24	16-May-24
pH, soil	EPA 150.1 - pH probe @ 25 °C, CaCl buffered ext.	17-May-24	17-May-24
PHC F1	CWS Tier 1 - P&T GC-FID	16-May-24	16-May-24
PHC F4G (gravimetric)	CWS Tier 1 - Extraction Gravimetric	22-May-24	23-May-24
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	16-May-24	21-May-24
REG 153: Metals by ICP/MS, soil	EPA 6020 - Digestion - ICP-MS	16-May-24	16-May-24
REG 153: PAHs by GC-MS	EPA 8270 - GC-MS, extraction	15-May-24	17-May-24
SAR	Calculated	16-May-24	16-May-24
Solids, %	CWS Tier 1 - Gravimetric	17-May-24	21-May-24

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH1-24-AU1	BH1-24-SS3	BH1-24-SS5	BH2-24-AU1		
	Sample Date:	08-May-24 09:00	08-May-24 09:00	08-May-24 09:00	08-May-24 09:00	-	-
	Sample ID:	2420225-01	2420225-02	2420225-03	2420225-04		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Physical Characteristics			•	•	•		•
% Solids	0.1 % by Wt.	91.5	91.5	90.4	96.4	-	-
General Inorganics	•					•	
SAR	0.01 N/A	2.19	1.60	1.87	4.11	-	-
Conductivity	5 uS/cm	2330	583	402	544	-	-
рН	0.05 pH Units	-	-	8.79	7.28	-	-
Metals					•	•	
Antimony	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Arsenic	1.0 ug/g	2.6	2.0	2.1	2.7	-	-
Barium	1.0 ug/g	99.5	29.9	35.8	87.8	-	-
Beryllium	0.5 ug/g	<0.5	<0.5	<0.5	<0.5	-	-
Boron	5.0 ug/g	11.0	6.2	<5.0	11.1	-	-
Cadmium	0.5 ug/g	<0.5	<0.5	<0.5	<0.5	-	-
Chromium	5.0 ug/g	20.5	10.5	11.8	14.8	-	-
Chromium (VI)	0.2 ug/g	0.8	0.3	<0.2	<0.2	-	-
Cobalt	1.0 ug/g	5.7	4.9	4.6	5.4	-	-
Copper	5.0 ug/g	16.1	9.9	6.8	12.8	-	-
Lead	1.0 ug/g	32.9	10.5	4.0	32.1	-	-
Mercury	0.1 ug/g	<0.1	<0.1	<0.1	<0.1	-	-
Molybdenum	1.0 ug/g	1.5	<1.0	<1.0	<1.0	-	-
Nickel	5.0 ug/g	13.2	10.8	8.3	11.5	-	-
Selenium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Silver	0.3 ug/g	<0.3	<0.3	<0.3	<0.3	-	-
Thallium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Uranium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Vanadium	10.0 ug/g	27.5	14.9	18.7	25.1	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH1-24-AU1	BH1-24-SS3	BH1-24-SS5	BH2-24-AU1		
	Sample Date:	08-May-24 09:00	08-May-24 09:00	08-May-24 09:00	08-May-24 09:00	_	_
	Sample ID:	2420225-01	2420225-02	2420225-03	2420225-04		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Metals	L			ļ.	Į l		
Zinc	20.0 ug/g	35.1	<20.0	<20.0	34.7	-	-
Volatiles							
Benzene	0.02 ug/g	<0.02	<0.02	<0.02	<0.02	-	-
Ethylbenzene	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
Toluene	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
m,p-Xylenes	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
o-Xylene	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
Xylenes, total	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
Toluene-d8	Surrogate	82.2%	111%	115%	97.6%	-	-
Hydrocarbons	•					-	
F1 PHCs (C6-C10)	7 ug/g	<7	<7	<7	<7	-	-
F2 PHCs (C10-C16)	4 ug/g	<4	<4	<4	<40	-	-
F3 PHCs (C16-C34)	8 ug/g	105	44	<8	164	-	-
F4 PHCs (C34-C50)	6 ug/g	263 [3]	74	<6	1040 [3]	-	-
F4G PHCs (gravimetric)	50 ug/g	382	-	-	1680	-	-
Semi-Volatiles							
Acenaphthene	0.02 ug/g	<0.02	0.08	0.02	<0.40 [1]	-	-
Acenaphthylene	0.02 ug/g	<0.02	<0.02	<0.02	<0.40 [1]	-	-
Anthracene	0.02 ug/g	0.03	0.20	0.06	0.54	-	-
Benzo [a] anthracene	0.02 ug/g	0.05	0.22	0.05	1.59	-	-
Benzo [a] pyrene	0.02 ug/g	0.04	0.15	0.04	1.41	-	-
Benzo [b] fluoranthene	0.02 ug/g	0.05	0.16	0.04	1.51	-	-
Benzo [g,h,i] perylene	0.02 ug/g	0.03	0.08	0.02	0.87	-	-
Benzo [k] fluoranthene	0.02 ug/g	0.03	0.12	0.03	0.79	-	-
Chrysene	0.02 ug/g	0.04	0.20	0.05	1.92	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH1-24-AU1	BH1-24-SS3	BH1-24-SS5	BH2-24-AU1		
	Sample Date:	08-May-24 09:00	08-May-24 09:00	08-May-24 09:00	08-May-24 09:00	-	-
	Sample ID:	2420225-01	2420225-02	2420225-03	2420225-04		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Semi-Volatiles	•						
Dibenzo [a,h] anthracene	0.02 ug/g	<0.02	0.02	<0.02	<0.40 [1]	-	-
Fluoranthene	0.02 ug/g	0.11	0.77	0.20	4.61	-	-
Fluorene	0.02 ug/g	<0.02	0.09	0.03	<0.40	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g	0.02	0.08	<0.02	0.73	-	-
1-Methylnaphthalene	0.02 ug/g	<0.02	<0.02	<0.02	<0.40 [1]	-	-
2-Methylnaphthalene	0.02 ug/g	<0.02	<0.02	<0.02	<0.40 [1]	-	-
Methylnaphthalene (1&2)	0.04 ug/g	<0.04	<0.04	<0.04	<0.80 [1]	-	-
Naphthalene	0.01 ug/g	0.02	0.06	0.04	<0.20 [1]	-	-
Phenanthrene	0.02 ug/g	0.06	0.67	0.19	2.40	-	-
Pyrene	0.02 ug/g	0.09	0.54	0.14	4.59	-	-
2-Fluorobiphenyl	Surrogate	70.3%	66.6%	62.2%	68.0%	-	-
Terphenyl-d14	Surrogate	79.6%	83.4%	79.8%	87.4%	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH2-24-SS3	BH3-24-AU1	BH3-24-SS5	BH4-24-AU1		
	Sample Date:	08-May-24 09:00	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	-	-
	Sample ID:	2420225-05	2420225-06	2420225-07	2420225-08		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Physical Characteristics			•	•	•		•
% Solids	0.1 % by Wt.	91.8	91.6	92.1	93.6	-	-
General Inorganics						•	
SAR	0.01 N/A	6.83	2.29	2.08	0.28	-	-
Conductivity	5 uS/cm	944	406	459	156	-	-
рН	0.05 pH Units	-	-	7.79	-	-	-
Metals					•	•	
Antimony	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Arsenic	1.0 ug/g	2.3	2.8	2.9	4.0	-	-
Barium	1.0 ug/g	41.1	59.5	21.3	98.2	-	-
Beryllium	0.5 ug/g	<0.5	<0.5	<0.5	<0.5	-	-
Boron	5.0 ug/g	<5.0	6.5	7.4	7.3	-	-
Cadmium	0.5 ug/g	<0.5	<0.5	<0.5	<0.5	-	-
Chromium (VI)	0.2 ug/g	<0.2	0.2	-	<0.2	-	-
Chromium	5.0 ug/g	11.7	13.0	11.7	10.4	-	-
Cobalt	1.0 ug/g	4.6	4.7	3.6	5.1	-	-
Copper	5.0 ug/g	8.3	12.4	5.5	11.8	-	-
Lead	1.0 ug/g	3.7	43.4	5.2	50.7	-	-
Mercury	0.1 ug/g	<0.1	0.1	-	<0.1	-	-
Molybdenum	1.0 ug/g	<1.0	<1.0	<1.0	1.0	-	-
Nickel	5.0 ug/g	8.4	9.5	8.3	11.3	-	-
Selenium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Silver	0.3 ug/g	<0.3	<0.3	<0.3	<0.3	-	-
Thallium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Uranium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Vanadium	10.0 ug/g	20.6	19.8	16.7	27.6	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Order Date: 14-May-2024

Project Description: PE6422

Report Date: 31-May-2024

	Client ID:	BH2-24-SS3	BH3-24-AU1	BH3-24-SS5	BH4-24-AU1		
	Sample Date:	08-May-24 09:00	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	-	-
	Sample ID:	2420225-05	2420225-06	2420225-07	2420225-08		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Metals			•				
Zinc	20.0 ug/g	<20.0	45.1	<20.0	26.6	-	-
Volatiles							•
Benzene	0.02 ug/g	<0.02	<0.02	<0.02	<0.02	-	-
Ethylbenzene	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
Toluene	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
m,p-Xylenes	0.05 ug/g	<0.05	<0.05	<0.05	<0.05	-	-
o-Xylene	0.05 ug/g	<0.05	0.07	<0.05	<0.05	-	-
Xylenes, total	0.05 ug/g	<0.05	0.07	<0.05	<0.05	-	•
Toluene-d8	Surrogate	103%	102%	114%	111%	-	-
Hydrocarbons							
F1 PHCs (C6-C10)	7 ug/g	<7	<7	<7	<7	-	-
F2 PHCs (C10-C16)	4 ug/g	39	<4	<4	<80 [2]	-	-
F3 PHCs (C16-C34)	8 ug/g	58	169	<8	420	-	-
F4 PHCs (C34-C50)	6 ug/g	80	135 [3]	<6	1800 [3]	-	-
F4G PHCs (gravimetric)	50 ug/g	-	131	-	2540	-	-
Semi-Volatiles			•				
Acenaphthene	0.02 ug/g	<0.02	<0.40 [1]	<0.02	<0.02	-	-
Acenaphthylene	0.02 ug/g	<0.02	<0.40 [1]	<0.02	0.20	-	-
Anthracene	0.02 ug/g	<0.02	1.20	<0.02	0.15	-	-
Benzo [a] anthracene	0.02 ug/g	<0.02	1.76	<0.02	0.27	-	-
Benzo [a] pyrene	0.02 ug/g	<0.02	1.31	<0.02	0.20	-	-
Benzo [b] fluoranthene	0.02 ug/g	<0.02	1.27	<0.02	0.19	-	-
Benzo [g,h,i] perylene	0.02 ug/g	<0.02	0.65	<0.02	0.14	-	-
Benzo [k] fluoranthene	0.02 ug/g	<0.02	0.78	<0.02	0.12	-	-
Chrysene	0.02 ug/g	<0.02	1.78	<0.02	0.24	-	-

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024

Order Date: 14-May-2024

Project Description: PE6422

	Client ID:	BH2-24-SS3	BH3-24-AU1	BH3-24-SS5	BH4-24-AU1		
	Sample Date:	08-May-24 09:00	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	-	-
	Sample ID:	2420225-05	2420225-06	2420225-07	2420225-08		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Semi-Volatiles			-	•			
Dibenzo [a,h] anthracene	0.02 ug/g	<0.02	<0.40 [1]	<0.02	0.03	-	-
Fluoranthene	0.02 ug/g	0.02	5.49	<0.02	0.51	-	-
Fluorene	0.02 ug/g	<0.02	0.43 [1]	<0.02	<0.02	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g	<0.02	0.61	<0.02	0.07	-	-
1-Methylnaphthalene	0.02 ug/g	<0.02	<0.40 [1]	<0.02	<0.02	-	-
2-Methylnaphthalene	0.02 ug/g	<0.02	<0.40 [1]	<0.02	0.02	-	-
Methylnaphthalene (1&2)	0.04 ug/g	<0.04	<0.80 [1]	<0.04	<0.04	-	-
Naphthalene	0.01 ug/g	<0.01	0.38	<0.01	0.01	-	-
Phenanthrene	0.02 ug/g	<0.02	4.17	<0.02	0.19	-	-
Pyrene	0.02 ug/g	0.02	4.56	<0.02	0.45	-	-
2-Fluorobiphenyl	Surrogate	70.5%	63.0%	66.1%	61.6%	-	-
Terphenyl-d14	Surrogate	87.8%	94.9%	82.5%	68.9%	-	-

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH4-24-SS2	BH4-24-SS3	BH4-24-SS6	BH5-24-AU1		
	Sample Date:	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	-	-
	Sample ID:	2420225-09	2420225-10	2420225-11	2420225-12		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Physical Characteristics							
% Solids	0.1 % by Wt.	80.8	93.6	94.5	93.4	-	-
General Inorganics							
SAR	0.01 N/A	2.30	-	2.21	2.00	-	-
Conductivity	5 uS/cm	272	-	357	363	-	-
рН	0.05 pH Units	-	-	-	7.40	-	-
Metals							
Antimony	1.0 ug/g	<1.0	<1.0	<1.0	1.7	-	-
Arsenic	1.0 ug/g	5.2	4.8	3.0	4.2	-	-
Barium	1.0 ug/g	177	36.2	36.7	137	-	-
Beryllium	0.5 ug/g	<0.5	<0.5	<0.5	<0.5	-	-
Boron	5.0 ug/g	5.3	6.7	7.4	5.2	-	-
Cadmium	0.5 ug/g	<0.5	<0.5	<0.5	0.5	-	-
Chromium (VI)	0.2 ug/g	<0.2	-	-	0.4	-	-
Chromium	5.0 ug/g	18.5	13.9	9.5	14.2	-	-
Cobalt	1.0 ug/g	4.6	6.4	2.9	4.8	-	-
Copper	5.0 ug/g	22.7	10.2	<5.0	53.8	-	-
Lead	1.0 ug/g	349	8.5	5.6	224	-	-
Mercury	0.1 ug/g	0.9	<0.1	-	0.2	-	
Molybdenum	1.0 ug/g	<1.0	1.1	<1.0	1.1	-	-
Nickel	5.0 ug/g	10.2	10.4	6.5	14.9	-	
Selenium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Silver	0.3 ug/g	<0.3	<0.3	<0.3	<0.3	-	-
Thallium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Uranium	1.0 ug/g	<1.0	<1.0	<1.0	<1.0	-	-
Vanadium	10.0 ug/g	21.1	21.1	13.6	28.8	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH4-24-SS2	BH4-24-SS3	BH4-24-SS6	BH5-24-AU1		
	Sample Date:	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	-	-
	Sample ID:	2420225-09	2420225-10	2420225-11	2420225-12		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Metals							
Zinc	20.0 ug/g	94.5	75.7	<20.0	228	-	-
Volatiles	·			·			·
Benzene	0.02 ug/g	<0.02	-	<0.02	<0.02	-	-
Ethylbenzene	0.05 ug/g	<0.05	-	<0.05	<0.05	-	-
Toluene	0.05 ug/g	<0.05	-	<0.05	<0.05	-	-
m,p-Xylenes	0.05 ug/g	<0.05	-	<0.05	<0.05	•	-
o-Xylene	0.05 ug/g	<0.05	-	<0.05	<0.05	-	-
Xylenes, total	0.05 ug/g	<0.05	-	<0.05	<0.05	-	-
Toluene-d8	Surrogate	85.9%	-	126%	98.4%	-	-
Hydrocarbons				-			
F1 PHCs (C6-C10)	7 ug/g	<7	-	<7	<7	-	-
F2 PHCs (C10-C16)	4 ug/g	6	-	9	<40 [2]	-	-
F3 PHCs (C16-C34)	8 ug/g	54	-	14	269	-	-
F4 PHCs (C34-C50)	6 ug/g	67 [3]	-	25	777 [3]	-	-
F4G PHCs (gravimetric)	50 ug/g	87	-	-	707	-	-
Semi-Volatiles							·
Acenaphthene	0.02 ug/g	0.03	<0.02	-	0.10	-	-
Acenaphthylene	0.02 ug/g	0.06	<0.02	-	0.14	•	-
Anthracene	0.02 ug/g	0.13	<0.02	-	0.26	-	-
Benzo [a] anthracene	0.02 ug/g	0.31	<0.02	-	0.83	-	-
Benzo [a] pyrene	0.02 ug/g	0.26	<0.02	-	0.60	-	-
Benzo [b] fluoranthene	0.02 ug/g	0.23	<0.02	-	0.74	-	-
Benzo [g,h,i] perylene	0.02 ug/g	0.14	<0.02	-	0.26	-	-
Benzo [k] fluoranthene	0.02 ug/g	0.16	<0.02	-	0.42	-	-
Chrysene	0.02 ug/g	0.25	<0.02	-	0.76	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH4-24-SS2	BH4-24-SS3	BH4-24-SS6	BH5-24-AU1		
	Sample Date:	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	09-May-24 09:00	-	-
	Sample ID:	2420225-09	2420225-10	2420225-11	2420225-12		
	Matrix:	Soil	Soil	Soil	Soil		
	MDL/Units						
Semi-Volatiles					•	-	
Dibenzo [a,h] anthracene	0.02 ug/g	0.02	<0.02	-	0.08	-	-
Fluoranthene	0.02 ug/g	0.77	<0.02	-	2.00	-	-
Fluorene	0.02 ug/g	0.02	<0.02	-	0.09	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g	0.13	<0.02	-	0.25	-	-
1-Methylnaphthalene	0.02 ug/g	<0.02	<0.02	-	0.02	-	-
2-Methylnaphthalene	0.02 ug/g	<0.02	<0.02	-	0.02	-	-
Methylnaphthalene (1&2)	0.04 ug/g	<0.04	<0.04	-	0.04	-	-
Naphthalene	0.01 ug/g	0.04	<0.01	-	0.03	-	-
Phenanthrene	0.02 ug/g	0.37	<0.02	-	1.02	-	-
Pyrene	0.02 ug/g	0.73	<0.02	-	1.67	-	-
2-Fluorobiphenyl	Surrogate	54.7%	64.5%	-	64.4%	-	-
Terphenyl-d14	Surrogate	73.9%	88.6%	-	73.5%	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH5-24-SS2(BOTTO M)	DUP1	DUP2			
	Sample Date:	09-May-24 09:00	08-May-24 09:00	09-May-24 09:00		_	-
	Sample ID:	2420225-13	2420225-14	2420225-15			
	Matrix:	Soil	Soil	Soil			
	MDL/Units	•					
Physical Characteristics							
% Solids	0.1 % by Wt.	92.9	92.0	93.7	-	-	-
General Inorganics	•				•		
SAR	0.01 N/A	1.54	2.18	1.75	-	-	-
Conductivity	5 uS/cm	353	433	347	-	-	-
pН	0.05 pH Units	-	9.60	-	-	-	-
Metals	•				•		,
Antimony	1.0 ug/g	<1.0	<1.0	<1.0	-	-	-
Arsenic	1.0 ug/g	2.1	2.2	2.7	-	-	-
Barium	1.0 ug/g	30.2	35.9	43.0	-	-	-
Beryllium	0.5 ug/g	<0.5	<0.5	<0.5	-	-	-
Boron	5.0 ug/g	<5.0	5.2	7.7	-	-	-
Cadmium	0.5 ug/g	<0.5	<0.5	<0.5	-	-	-
Chromium (VI)	0.2 ug/g	-	<0.2	-	-	-	-
Chromium	5.0 ug/g	10.1	12.2	10.5	-	-	-
Cobalt	1.0 ug/g	3.6	4.8	3.3	-	-	-
Copper	5.0 ug/g	6.6	7.8	<5.0	-	-	-
Lead	1.0 ug/g	3.7	4.2	5.6	-	-	-
Mercury	0.1 ug/g	-	<0.1	-	-	-	-
Molybdenum	1.0 ug/g	<1.0	<1.0	<1.0	-	-	-
Nickel	5.0 ug/g	6.9	8.7	7.1	-	-	-
Selenium	1.0 ug/g	<1.0	<1.0	<1.0	-	-	-
Silver	0.3 ug/g	<0.3	<0.3	<0.3	-	-	-
Thallium	1.0 ug/g	<1.0	<1.0	<1.0	-	-	-
Uranium	1.0 ug/g	<1.0	<1.0	<1.0	-	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH5-24-SS2(BOTTO M)	DUP1	DUP2			
	Sample Date:	09-May-24 09:00	08-May-24 09:00	09-May-24 09:00		-	_
	Sample ID:	2420225-13	2420225-14	2420225-15			
	Matrix:	Soil	Soil	Soil			
	MDL/Units						
Metals							•
Vanadium	10.0 ug/g	16.2	19.1	14.9	-	-	-
Zinc	20.0 ug/g	<20.0	<20.0	<20.0	-	-	-
Volatiles							•
Benzene	0.02 ug/g	<0.02	<0.02	<0.02	-	-	-
Ethylbenzene	0.05 ug/g	<0.05	<0.05	<0.05	-	-	-
Toluene	0.05 ug/g	<0.05	<0.05	<0.05	-	-	-
m,p-Xylenes	0.05 ug/g	<0.05	<0.05	<0.05	-	-	-
o-Xylene	0.05 ug/g	<0.05	<0.05	<0.05	-	-	-
Xylenes, total	0.05 ug/g	<0.05	<0.05	<0.05	-	-	-
Toluene-d8	Surrogate	126%	107%	106%	-	-	-
Hydrocarbons				-			
F1 PHCs (C6-C10)	7 ug/g	<7	<7	<7	-	-	-
F2 PHCs (C10-C16)	4 ug/g	<4	<4	<4	-	-	-
F3 PHCs (C16-C34)	8 ug/g	<8	15	<8	-	-	-
F4 PHCs (C34-C50)	6 ug/g	<6	13	<6	-	-	-
Semi-Volatiles							
Acenaphthene	0.02 ug/g	<0.02	0.02	-	-	-	-
Acenaphthylene	0.02 ug/g	<0.02	<0.02	-	-	-	-
Anthracene	0.02 ug/g	<0.02	0.06	-	-	-	-
Benzo [a] anthracene	0.02 ug/g	<0.02	0.05	-	-	-	-
Benzo [a] pyrene	0.02 ug/g	<0.02	0.04	-	-	-	-
Benzo [b] fluoranthene	0.02 ug/g	<0.02	0.03	-	-	-	-
Benzo [g,h,i] perylene	0.02 ug/g	<0.02	<0.02	-	-	-	-
Benzo [k] fluoranthene	0.02 ug/g	<0.02	0.03	-	-	-	-
Chrysene	0.02 ug/g	<0.02	0.05	-	-	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

	Client ID:	BH5-24-SS2(BOTTO M)	DUP1	DUP2			
	Sample Date:	,	08-May-24 09:00	09-May-24 09:00		_	-
	Sample ID:	· ·	2420225-14	2420225-15			
	Matrix:	Soil	Soil	Soil			
	MDL/Units						
Semi-Volatiles							
Dibenzo [a,h] anthracene	0.02 ug/g	<0.02	<0.02	-	-	-	-
Fluoranthene	0.02 ug/g	<0.02	0.18	-	-	-	•
Fluorene	0.02 ug/g	<0.02	0.02	-	-	-	-
Indeno [1,2,3-cd] pyrene	0.02 ug/g	<0.02	<0.02	-	-	-	-
1-Methylnaphthalene	0.02 ug/g	<0.02	<0.02	-	-	-	-
2-Methylnaphthalene	0.02 ug/g	<0.02	<0.02	-	-	-	-
Methylnaphthalene (1&2)	0.04 ug/g	<0.04	<0.04	-	-	-	-
Naphthalene	0.01 ug/g	<0.01	0.02	-	-	-	-
Phenanthrene	0.02 ug/g	<0.02	0.18	-	-	-	-
Pyrene	0.02 ug/g	<0.02	0.14	-	-	-	-
2-Fluorobiphenyl	Surrogate	72.6%	54.1%	-	-	-	-
Terphenyl-d14	Surrogate	105%	62.6%	-	-	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024

Order Date: 14-May-2024

Project Description: PE6422

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	%REC	%REC Limit	RPD	RPD Limit	Notes
General Inorganics								
Conductivity	ND	5	uS/cm					
Hydrocarbons								
F1 PHCs (C6-C10)	ND	7	ug/g					
F2 PHCs (C10-C16)	ND	4	ug/g					
F3 PHCs (C16-C34)	ND	8	ug/g					
F4 PHCs (C34-C50)	ND	6	ug/g					
F4G PHCs (gravimetric)	ND	50	ug/g					
Metals								
Antimony	ND	1.0	ug/g					
Arsenic	ND	1.0	ug/g					
Barium	ND	1.0	ug/g					
Beryllium	ND	0.5	ug/g					
Boron	ND	5.0	ug/g					
Cadmium	ND	0.5	ug/g					
Chromium (VI)	ND	0.2	ug/g					
Chromium	ND	5.0	ug/g					
Cobalt	ND	1.0	ug/g					
Copper	ND	5.0	ug/g					
Lead	ND	1.0	ug/g					
Mercury	ND	0.1	ug/g					
Molybdenum	ND	1.0	ug/g					
Nickel	ND	5.0	ug/g					
Selenium	ND	1.0	ug/g					
Silver	ND	0.3	ug/g					
Thallium	ND	1.0	ug/g					
Uranium	ND	1.0	ug/g					
Vanadium	ND	10.0	ug/g					
Zinc	ND	20.0	ug/g					
Semi-Volatiles								
Acenaphthene	ND	0.02	ug/g					
Acenaphthylene	ND	0.02	ug/g					
Anthracene	ND	0.02	ug/g					

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024 Order Date: 14-May-2024 Project Description: PE6422

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	%REC	%REC Limit	RPD	RPD Limit	Notes
Benzo [a] anthracene	ND	0.02	ug/g					
Benzo [a] pyrene	ND	0.02	ug/g					
Benzo [b] fluoranthene	ND	0.02	ug/g					
Benzo [g,h,i] perylene	ND	0.02	ug/g					
Benzo [k] fluoranthene	ND	0.02	ug/g					
Chrysene	ND	0.02	ug/g					
Dibenzo [a,h] anthracene	ND	0.02	ug/g					
Fluoranthene	ND	0.02	ug/g					
Fluorene	ND	0.02	ug/g					
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g					
1-Methylnaphthalene	ND	0.02	ug/g					
2-Methylnaphthalene	ND	0.02	ug/g					
Methylnaphthalene (1&2)	ND	0.04	ug/g					
Naphthalene	ND	0.01	ug/g					
Phenanthrene	ND	0.02	ug/g					
Pyrene	ND	0.02	ug/g					
Surrogate: 2-Fluorobiphenyl	0.688		%	51.6	50-140			
Surrogate: Terphenyl-d14	1.11		%	83.4	50-140			
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	8.74		%	109	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024

Order Date: 14-May-2024

Project Description: PE6422

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
General Inorganics									
SAR	2.09	0.01	N/A	2.19			4.7	30	
Conductivity	2300	5	uS/cm	2330			1.3	5	
pH	6.22	0.05	pH Units	6.24			0.3	2.3	
Hydrocarbons F1 PHCs (C6-C10)	ND	7	ug/g	ND			NC	40	
F2 PHCs (C10-C16)	ND	4	ug/g	ND			NC	30	
F3 PHCs (C16-C34)	79	8	ug/g	105			28.5	30	
F4 PHCs (C34-C50)	194	6	ug/g	263			30.4	30	QR-05
Metals									
Antimony	ND	1.0	ug/g	ND			NC	30	
Arsenic	3.4	1.0	ug/g	3.4			0.6	30	
Barium	314	1.0	ug/g	333			6.0	30	
Beryllium	1.1	0.5	ug/g	1.0			9.1	30	
Boron	13.6	5.0	ug/g	13.3			2.9	30	
Cadmium	ND	0.5	ug/g	ND			NC	30	
Chromium (VI)	0.2	0.2	ug/g	0.2			0.0	35	
Chromium	48.8	5.0	ug/g	50.0			2.6	30	
Cobalt	15.5	1.0	ug/g	16.1			3.6	30	
Copper	31.1	5.0	ug/g	32.3			3.8	30	
Lead	23.7	1.0	ug/g	24.3			2.3	30	
Mercury	ND	0.1	ug/g	ND			NC	30	
Molybdenum	ND	1.0	ug/g	ND			NC	30	
Nickel	32.8	5.0	ug/g	33.5			2.1	30	
Selenium	ND	1.0	ug/g	ND			NC	30	
Silver	ND	0.3	ug/g	ND			NC	30	
Thallium	ND	1.0	ug/g	ND			NC	30	
Uranium	ND	1.0	ug/g	ND			NC	30	
Vanadium	66.6	10.0	ug/g	68.8			3.2	30	
Zinc	86.5	20.0	ug/g	90.3			4.3	30	
Physical Characteristics									

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024

Order Date: 14-May-2024

Project Description: PE6422

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
% Solids	92.5	0.1	% by Wt.	91.8			0.8	25	
Semi-Volatiles									
Acenaphthene	ND	0.02	ug/g	ND			NC	40	
Acenaphthylene	ND	0.02	ug/g	ND			NC	40	
Anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] anthracene	ND	0.02	ug/g	ND			NC	40	
Benzo [a] pyrene	ND	0.02	ug/g	ND			NC	40	
Benzo [b] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Benzo [g,h,i] perylene	ND	0.02	ug/g	ND			NC	40	
Benzo [k] fluoranthene	ND	0.02	ug/g	ND			NC	40	
Chrysene	ND	0.02	ug/g	ND			NC	40	
Dibenzo [a,h] anthracene	ND	0.02	ug/g	ND			NC	40	
Fluoranthene	ND	0.02	ug/g	ND			NC	40	
Fluorene	ND	0.02	ug/g	ND			NC	40	
Indeno [1,2,3-cd] pyrene	ND	0.02	ug/g	ND			NC	40	
1-Methylnaphthalene	ND	0.02	ug/g	ND			NC	40	
2-Methylnaphthalene	ND	0.02	ug/g	ND			NC	40	
Naphthalene	ND	0.01	ug/g	ND			NC	40	
Phenanthrene	ND	0.02	ug/g	ND			NC	40	
Pyrene	ND	0.02	ug/g	ND			NC	40	
Surrogate: 2-Fluorobiphenyl	1.44		%		66.7	50-140			
Surrogate: Terphenyl-d14	1.92		%		88.7	50-140			
Volatiles									
Benzene	ND	0.02	ug/g	ND			NC	50	
Ethylbenzene	ND	0.05	ug/g	ND			NC	50	
Toluene	ND	0.05	ug/g	ND			NC	50	
m,p-Xylenes	ND	0.05	ug/g	ND			NC	50	
o-Xylene	ND	0.05	ug/g	ND			NC	50	
Surrogate: Toluene-d8	9.46		%		108	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024

Order Date: 14-May-2024

Project Description: PE6422

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	186	7	ug/g	ND	108	85-115			
F2 PHCs (C10-C16)	102	4	ug/g	ND	117	60-140			
F3 PHCs (C16-C34)	342	8	ug/g	105	111	60-140			
F4 PHCs (C34-C50)	395	6	ug/g	263	96.9	60-140			
F4G PHCs (gravimetric)	920	50	ug/g	ND	92.0	80-120			
Metals									
Arsenic	46.9	1.0	ug/g	1.4	91.1	70-130			
Barium	166	1.0	ug/g	133	65.4	70-130			QM-07
Beryllium	47.4	0.5	ug/g	ND	94.0	70-130			
Boron	47.2	5.0	ug/g	5.3	83.8	70-130			
Cadmium	42.6	0.5	ug/g	ND	85.0	70-130			
Chromium (VI)	4.2	0.2	ug/g	0.2	72.5	70-130			
Chromium	68.0	5.0	ug/g	20.0	95.9	70-130			
Cobalt	53.0	1.0	ug/g	6.4	93.2	70-130			
Copper	57.9	5.0	ug/g	12.9	89.9	70-130			
Lead	52.8	1.0	ug/g	9.7	86.2	70-130			
Mercury	1.43	0.1	ug/g	ND	95.0	70-130			
Molybdenum	45.5	1.0	ug/g	ND	90.4	70-130			
Nickel	58.6	5.0	ug/g	13.4	90.5	70-130			
Selenium	45.2	1.0	ug/g	ND	90.1	70-130			
Silver	37.2	0.3	ug/g	ND	74.3	70-130			
Thallium	45.2	1.0	ug/g	ND	90.1	70-130			
Uranium	47.6	1.0	ug/g	ND	94.7	70-130			
Vanadium	74.6	10.0	ug/g	27.5	94.1	70-130			
Zinc	77.2	20.0	ug/g	36.1	82.1	70-130			
Semi-Volatiles									
Acenaphthene	0.221	0.02	ug/g	ND	81.6	50-140			
Acenaphthylene	0.231	0.02	ug/g	ND	85.3	50-140			
Anthracene	0.247	0.02	ug/g	ND	91.3	50-140			
Benzo [a] anthracene	0.195	0.02	ug/g	ND	72.0	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196

Report Date: 31-May-2024

Order Date: 14-May-2024

Project Description: PE6422

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Benzo [a] pyrene	0.159	0.02	ug/g	ND	58.9	50-140			
Benzo [b] fluoranthene	0.178	0.02	ug/g	ND	65.7	50-140			
Benzo [g,h,i] perylene	0.166	0.02	ug/g	ND	61.3	50-140			
Benzo [k] fluoranthene	0.194	0.02	ug/g	ND	71.7	50-140			
Chrysene	0.205	0.02	ug/g	ND	75.9	50-140			
Dibenzo [a,h] anthracene	0.174	0.02	ug/g	ND	64.4	50-140			
Fluoranthene	0.251	0.02	ug/g	ND	92.8	50-140			
Fluorene	0.209	0.02	ug/g	ND	77.1	50-140			
Indeno [1,2,3-cd] pyrene	0.146	0.02	ug/g	ND	54.1	50-140			
1-Methylnaphthalene	0.205	0.02	ug/g	ND	75.7	50-140			
2-Methylnaphthalene	0.173	0.02	ug/g	ND	64.0	50-140			
Naphthalene	0.208	0.01	ug/g	ND	76.9	50-140			
Phenanthrene	0.208	0.02	ug/g	ND	77.0	50-140			
Pyrene	0.249	0.02	ug/g	ND	92.0	50-140			
Surrogate: 2-Fluorobiphenyl	1.34		%		62.0	50-140			
Surrogate: Terphenyl-d14	1.82		%		84.0	50-140			
Volatiles									
Benzene	2.97	0.02	ug/g	ND	74.3	60-130			
Ethylbenzene	3.01	0.05	ug/g	ND	75.4	60-130			
Toluene	2.88	0.05	ug/g	ND	72.0	60-130			
m,p-Xylenes	6.29	0.05	ug/g	ND	78.6	60-130			
o-Xylene	2.73	0.05	ug/g	ND	68.3	60-130			



Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60196 Project Description: PE6422

Qualifier Notes:

Sample Qualifiers:

1: Elevated reporting limit due to dilution required because of high target analyte concentration.

2: Elevated reporting limits due to the nature of the sample matrix.

Applies to Samples: BH4-24-AU1, BH5-24-AU1

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

Applies to Samples: BH1-24-AU1, BH2-24-AU1, BH3-24-AU1, BH4-24-AU1, BH4-24-SS2, BH5-24-AU1

QC Qualifiers:

QM-07 The spike recovery was outside acceptance limits for the MS and/or MSD. The batch was accepted based on other acceptable QC.

QR-05 Duplicate RPDs higher than normally accepted. Remaining batch QA\QC was acceptable. May be sample effect.

Sample Data Revisions:

None

Report Date: 31-May-2024



Report Date: 31-May-2024

Order Date: 14-May-2024

Project Description: PE6422

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Work Order Revisions / Comments:

Revision 1 - Revised report includes additional metals and PAH analyses.

Other Report Notes:

Client PO: 60196

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 unlesss otherwise noted.

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.

Any use of these results implies your agreement that our total liabilty 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.





nt 6hd. 1G 4.JB labs.com

Paracel Order Number (Lab Use Only)

2420225

Chain Of Custody (Lab Use Only)

Clie	ent Name: Paterson O) Jan	29											
_	Paterson Group Inc.			Proje	Project Ref: PE6422								Page <u>1</u> of <u>2</u>						
	Address				Quote #:								Turnaround Time						
, au	9 AURIGA DRIVE			PO#	:	60196							☐ 1 day ☐ 3					□ 3 c	ay
_	OTTAWA ON K2E 71	9		E-ma	il: j	camposarcone@p	patersongroup	.ca										× Re	gular
Tele	phone: 613-226-7381			1		kmunch@patersor								Requ					,
	REG 153/04 REG 406/19	Other Regulation	T	_					-1 25	(1)		77.78	8201100	1101529		111111	1 1 1		
	Table 1 Res/Park Med/Fine					S (Soil/Sed.) GW (G Nater) SS (Storm/Sa						Re	equired Analysis						
	Table 2 ☐ Ind/Comm ☐ Coarse	☐ CCME ☐ MISA		J., (J.	V (Surface Water) SS (Storm/Sanitary Sewer) P (Paint) A (Air) O (Other)														
	Table 3 Agri/Other	SU - Sani SU - Storm		Т	1 9			BTEX											
Пт	able	Mun:		9	Containers	Sample Taken		4 +	s		G G			WS)					
	For RSC: ☐ Yes ☐ No	Other:	ě	Air Volume				F1-F4			ls by								
	Sample ID/Locatio	n Name	Matrix	Air	Jo #	Date	Time	유	VOCs	PAHs	Metals	ρĤ	S Z	B (HWS)	된	2	SAR		
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2	BH1-24-SS3		s		2	5/8/2024		V		V	V	V	~		М	7	V	П	M
3	BH1-24-SS5		s		2	5/8/2024		V	H	V	V	V		H	7	7	V	H	Ħ
4	BH2-24-AU1		s		2	5/8/2024		1	H	V	7			H			V	H	Ħ
5	BH2-24-SS3		s		2	5/8/2024		7	H	\ \ \				H	H	7		믐	Ħ
6	BH3-24-AU1		s		2	5/9/2024		V	H	V	7	7	V	H	H	<i>V</i>	V	H	H
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Chain Of Custody Paracel Order Number (Lab Use Only) (Lab Use Only)

_	LABORATORIES LTD. I RI	10112					2	420	22	25								
	nt Name: Paterson Group Inc.	Project Ref. PE6422								Page 2 of 2								
Conta	act Name: Jeremy Camposarcone		Quote	#:								Turnaround Time						
Addr	ess: 9 AURIGA DRIVE		PO #:	6	60196							1 0	1 day				3 day	
	OTTAWA ON K2E 7T9		E-mail	l: jo	camposarcone@p	patersongroup.ca	1					2 day					■ Regular	
Telep	phone: 613-226-7381		1		kmunch@patersongroup.ca							Date Required:					ga.a.	
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	able 1 Res/Park Med/Fine REG 558 PWQO			rface V	Vater) SS (Storm/Sa	nitary Sewer)					Re	quire	Anal	ysis				
	able 2 ☐ Ind/Comm ☐ Coarse ☐ CCME ☐ MISA			P (P	aint) A (Air) O (Ot	her)	втех										\Box	
	able 3 Agri/Other SU - Sani SU - Storm			ers			+ BT			ICP								
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	For RSC: Yes No Other:	Matrix	Air Volume	of Cor			C F1	S	£	Metals k		_	B (HWS)			æ		
	Sample ID/Location Name	Σ	Ą	#	Date	Time	오	VOCs	PAHs	Mei	롸	CrVI	B (F	펀	EC	SAR		
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elinqu	sished By (Prince Jeremy Camposarcone Date/Time:					Date/Time: W	14.	2000	11	_	Date/Time: May 24 0928							
ate/Ti	ime: 3/3/2024 Temperature:	Section 1			°C		2.7		4,	X XP	pH Ver	ified:	KIA	By:		100		



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

Order Date:

Report Date:

14-May-24

31-May-24

Subcontracted Analysis

Paterson Group Consulting Engineers (Ottawa)

9 Auriga Drive

Ottawa, ON K2E 7T9

Attn: Jeremy Camposarcone

Paracel Report No. 2420225

Client Project(s): **PE6422**Client PO: **60196**

Reference: #24-017 Standing Offer 2024

CoC Number:

Sample(s) from this project were subcontracted for the listed parameters. A copy of the subcontractor's report is attached

Paracel ID Client ID Analysis

2420225-10 BH4-24-SS3 Methyl Mercury - soil

ALS Canada Ltd.



CERTIFICATE OF ANALYSIS

Work Order : WT2413665

Client : Paracel Laboratories Ltd

Contact : Mark Foto

Address : 2319 St. Laurent Blvd. Unit 300

Ottawa ON Canada K1G 4J8

Telephone : 613 731 9577

Project : 2420225

PO : ----C-O-C number · ----

Sampler : CLIENT

Site : ----

Quote number : Standing Offer 2024

No. of samples received : 1
No. of samples analysed : 1

Page : 1 of 2

Laboratory : ALS Environmental - Waterloo

Account Manager : Costas Farassoglou

Address : 60 Northland Road, Unit 1

Waterloo ON Canada N2V 2B8

Telephone : 613 225 8279

Date Samples Received : 28-May-2024 12:10

Date Analysis Commenced : 02-Jun-2024

Issue Date : 13-Jun-2024 09:08

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. This document shall not be reproduced, except in full.

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results

Additional information pertinent to this report will be found in the following separate attachments: Quality Control Report, QC Interpretive report to assist with Quality Review and Sample Receipt Notification (SRN).

Signatories

This document has been electronically signed by the authorized signatories below. Electronic signing is conducted in accordance with US FDA 21 CFR Part 11.

Signatories Position Laboratory Department

Kinny Wu Lab Analyst Metals, Burnaby, British Columbia

Page : 2 of 2

Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



General Comments

The analytical methods used by ALS are developed using internationally recognized reference methods (where available), such as those published by US EPA, APHA Standard Methods, ASTM, ISO, Environment Canada, BC MOE, and Ontario MOE. Refer to the ALS Quality Control Interpretive report (QCI) for applicable references and methodology summaries. Reference methods may incorporate modifications to improve performance.

Where a reported less than (<) result is higher than the LOR, this may be due to primary sample extract/digestate dilution and/or insufficient sample for analysis.

Where the LOR of a reported result differs from standard LOR, this may be due to high moisture content, insufficient sample (reduced weight employed) or matrix interference.

Please refer to Quality Control Interpretive report (QCI) for information regarding Holding Time compliance.

Key: CAS Number: Chemical Abstracts Services number is a unique identifier assigned to discrete substances

LOR: Limit of Reporting (detection limit).

Unit	Description
μg/kg	micrograms per kilogram

<: less than.

>: greater than.

Surrogate: An analyte that is similar in behavior to target analyte(s), but that does not occur naturally in environmental samples. For applicable tests, surrogates are added to samples prior to analysis as a check on recovery.

Test results reported relate only to the samples as received by the laboratory.

UNLESS OTHERWISE STATED on SRN or QCI Report, ALL SAMPLES WERE RECEIVED IN ACCEPTABLE CONDITION.

Analytical Results

Sub-Matrix: Soil/Solid			CI	ient sample ID	BH4-24-SS3		 	
(Matrix: Soil/Solid)								
	Client sampling date / time						 	
Analyte	CAS Number	Method/Lab	LOR	Unit	WT2413665-001		 	
					Result		 	
Speciated Metals								
Methylmercury (as MeHg)	22967-92-6	E538/VA	0.050	μg/kg	<0.050		 	

Please refer to the General Comments section for an explanation of any result qualifiers detected.

Please refer to the Accreditation section for an explanation of analyte accreditations.



QUALITY CONTROL INTERPRETIVE REPORT

Work Order : **WT2413665** Page : 1 of 5

Client : Paracel Laboratories Ltd Laboratory : ALS Environmental - Waterloo

Contact : Mark Foto Account Manager : Costas Farassoglou

Address : 2319 St. Laurent Blvd. Unit 300 Address : 60 Northland Road, Unit 1

Ottawa ON Canada K1G 4J8 Waterloo, Ontario Canada N2V 2B8

 Telephone
 : 613 731 9577
 Telephone
 : 613 225 8279

 Project
 : 2420225
 Date Samples Received
 : 28-May-2024 12:10

PO :--- Issue Date : 13-Jun-2024 09:08 C-O-C number :---

Quote number : Standing Offer 2024

: CLIENT

No. of samples received :1

No. of samples analysed :1

This report is automatically generated by the ALS LIMS (Laboratory Information Management System) through evaluation of Quality Control (QC) results and other QA parameters associated with this submission, and is intended to facilitate rapid data validation by auditors or reviewers. The report highlights any exceptions and outliers to ALS Data Quality Objectives, provides holding time details and exceptions, summarizes QC sample frequencies, and lists applicable methodology references and summaries.

Key

Sampler

Site

Anonymous: Refers to samples which are not part of this work order, but which formed part of the QC process lot.

CAS Number: Chemical Abstracts Service number is a unique identifier assigned to discrete substances.

DQO: Data Quality Objective.

LOR: Limit of Reporting (detection limit).

RPD: Relative Percent Difference.

Workorder Comments

Holding times are displayed as "---" if no guidance exists from CCME, Canadian provinces, or broadly recognized international references.

Summary of Outliers

Outliers: Quality Control Samples

- No Method Blank value outliers occur.
- No Duplicate outliers occur.
- No Laboratory Control Sample (LCS) outliers occur
- No Test sample Surrogate recovery outliers exist.

Outliers: Reference Material (RM) Samples

No Reference Material (RM) Sample outliers occur.

Outliers : Analysis Holding Time Compliance (Breaches)

• Analysis Holding Time Outliers exist - please see following pages for full details.

Outliers : Frequency of Quality Control Samples ■ No Quality Control Sample Frequency Outliers occur.		
		alsglobal.com

Page : 3 of 5 Work Order · WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



Analysis Holding Time Compliance

This report summarizes extraction / preparation and analysis times and compares each with ALS recommended holding times, which are selected to meet known provincial and/or federal requirements. In the absence of regulatory hold times, ALS establishes recommendations based on guidelines published by organizations such as CCME, US EPA, APHA Standard Methods, ASTM, or Environment Canada (where available). Dates and holding times reported below represent the first dates of extraction or analysis. If subsequent tests or dilutions exceeded holding times, qualifiers are added (refer to COA).

If samples are identified below as having been analyzed or extracted outside of recommended holding times, measurement uncertainties may be increased, and this should be taken into consideration when interpreting results.

Where actual sampling date is not provided on the chain of custody, the date of receipt with time at 00:00 is used for calculation purposes.

Where only the sample date without time is provided on the chain of custody, the sampling date at 00:00 is used for calculation purposes.

Matrix: Soil/Solid

Evaluation: **x** = Holding time exceedance; ✓ = Within Holding Time

Matrix Com Cond						diddion.	i loiding time exces	Judinoo ,	* * 1 (111	riolaning riiii
Analyte Group : Analytical Method	Method	Sampling Date	Ext	raction / Pro	eparation			Analysis		
Container / Client Sample ID(s)			Preparation	Holding Times		Eval	Analysis Date	Holding	g Times	Eval
			Date	Rec	Actual			Rec	Actual	
Speciated Metals : Methylmercury in Soil by GCAFS										
Glass soil jar/Teflon lined cap [ON MECP]										
BH4-24-SS3	E538	09-May-2024	07-Jun-2024	28	29	*	10-Jun-2024	28 days	3 days	✓
				days	days	EHT				

Legend & Qualifier Definitions

EHT: Exceeded ALS recommended hold time prior to analysis.

Rec. HT: ALS recommended hold time (see units).

Page : 4 of 5 Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



Quality Control Parameter Frequency Compliance

The following report summarizes the frequency of laboratory QC samples analyzed within the analytical batches (QC lots) in which the submitted samples were processed. The actual frequency should be greater than or equal to the expected frequency.

Matrix: Soil/Solid Evaluation: × = QC frequency outside specification; ✓ = QC frequency within specification										
Quality Control Sample Type			Co	unt)				
Analytical Methods	Method	QC Lot #	QC	Regular	Actual	Expected	Evaluation			
Laboratory Duplicates (DUP)										
Methylmercury in Soil by GCAFS	E538	1479163	1	11	9.0	5.0	✓			
Laboratory Control Samples (LCS)										
Methylmercury in Soil by GCAFS	E538	1479163	2	11	18.1	10.0	✓			
Method Blanks (MB)										
Methylmercury in Soil by GCAFS	E538	1479163	1	11	9.0	5.0	✓			

Page : 5 of 5 Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



Methodology References and Summaries

The analytical methods used by ALS are developed using internationally recognized reference methods (where available), such as those published by US EPA, APHA Standard Methods, ASTM, ISO, Environment Canada, BC MOE, and Ontario MOE. Reference methods may incorporate modifications to improve performance (indicated by "mod").

Analytical Methods	Method / Lab	Matrix	Method Reference	Method Descriptions					
Methylmercury in Soil by GCAFS	E538 ALS Environmental - Vancouver	Soil/Solid	DeWild et al. (2004)/EPA 1630 (mod)	This method follows procedures published by DeWild, Olund, Olsen and Tate (2004) for the US Geological Survey (Techniques and Methods 5A-7). Samples are leached with an acidic copper sulphate solution to solubilize methylmercury for inorganic complexes. The methylmercury is then extracted into dichloromethane and then an aliquot is back extracted into ultra-pure water. The extract is analyzed by aqueous phase ethylation purge and trap, desorption and GC separation. The separated species are the pyrolized to elemental Hg and quantified by cold vapour atomic flourescence spectroscopy. Results are reported "as MeHg".					
Preparation Methods	Method / Lab	Matrix	Method Reference	Method Descriptions					
Methylmercury Soil Digestion	EP538 ALS Environmental - Vancouver	Soil/Solid	DeWild et al. (2004)	This method follows procedures published by DeWild, Olund, Olsen and Tate (2004) for the US Geological Survey (Techniques and Methods 5A-7). Samples are leached with an acidic copper sulphate solution to solubilize methylmercury for inorganic complexes. The methylmercury is then extracted into dichloromethane and then an aliquot is back extracted into ultra-pure water. The extract is analyzed by aqueous phase ethylation, purge and trap, desorption and GC separation. The separated species are then pyrolized to elemental Hg and quantified by cold vapour atomic flourescence spectroscopy. Results are reported "as MeHg".					

ALS Canada Ltd.



QUALITY CONTROL REPORT

Work Order : WT2413665

Client : Paracel Laboratories Ltd

Contact : Mark Foto

Address : 2319 St. Laurent Blvd. Unit 300

Ottawa ON Canada K1G 4J8

Telephone : 613 731 9577

Project : 2420225

PO :----C-O-C number :----

Sampler : CLIENT

Site :----

Quote number : Standing Offer 2024

No. of samples received : 1
No. of samples analysed : 1

Page : 1 of 3

Laboratory : ALS Environmental - Waterloo

Account Manager : Costas Farassoglou

Address : 60 Northland Road, Unit 1

Waterloo, Ontario Canada N2V 2B8

Telephone :613 225 8279

Date Samples Received : 28-May-2024 12:10

Date Analysis Commenced : 02-Jun-2024

Issue Date : 13-Jun-2024 09:08

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. This document shall not be reproduced, except in full.

This Quality Control Report contains the following information:

Laboratory Duplicate (DUP) Report; Relative Percent Difference (RPD) and Data Quality Objectives

- Reference Material (RM) Report; Recovery and Data Quality Objectives
- Method Blank (MB) Report; Recovery and Data Quality Objectives
- Laboratory Control Sample (LCS) Report; Recovery and Data Quality Objectives

Signatories

This document has been electronically signed by the authorized signatories below. Electronic signing is conducted in accordance with US FDA 21 CFR Part 11.

Signatories Position Laboratory Department

Kinny Wu Lab Analyst Vancouver Metals, Burnaby, British Columbia

Page : 2 of 3 Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225

ALS

General Comments

The ALS Quality Control (QC) report is optionally provided to ALS clients upon request. ALS test methods include comprehensive QC checks with every analysis to ensure our high standards of quality are met. Each QC result has a known or expected target value, which is compared against predetermined Data Quality Objectives (DQOs) to provide confidence in the accuracy of associated test results. This report contains detailed results for all QC results applicable to this sample submission. Please refer to the ALS Quality Control Interpretation report (QCI) for applicable method references and methodology summaries.

Key:

Anonymous = Refers to samples which are not part of this work order, but which formed part of the QC process lot.

CAS Number = Chemical Abstracts Service number is a unique identifier assigned to discrete substances.

DQO = Data Quality Objective.

LOR = Limit of Reporting (detection limit).

RPD = Relative Percent Difference

= Indicates a QC result that did not meet the ALS DQO.

Workorder Comments

Holding times are displayed as "---" if no guidance exists from CCME, Canadian provinces, or broadly recognized international references.

Laboratory Duplicate (DUP) Report

A Laboratory Duplicate (DUP) is a randomly selected intralaboratory replicate sample. Laboratory Duplicates provide information regarding method precision and sample heterogeneity. ALS DQOs for Laboratory Duplicates are expressed as test-specific limits for Relative Percent Difference (RPD), or as an absolute difference limit of 2 times the LOR for low concentration duplicates within ~ 4-10 times the LOR (cut-off is test-specific).

Sub-Matrix: Soil/Solid	ub-Matrix: Soil/Solid						Laboratory Duplicate (DUP) Report									
Laboratory sample ID	Client sample ID	Analyte	CAS Number	Method	LOR	Unit	Original Result	Duplicate Result	RPD(%) or Difference	Duplicate Limits	Qualifier					
Speciated Metals (C	C Lot: 1479163)															
VA24B0502-001	Anonymous	Methylmercury (as MeHg)	22967-92-6	E538	0.050	μg/kg	0.230	0.133	0.097	Diff <2x LOR						

Method Blank (MB) Report

A Method Blank is an analyte-free matrix that undergoes sample processing identical to that carried out for test samples. Method Blank results are used to monitor and control for potential contamination from the laboratory environment and reagents. For most tests, the DQO for Method Blanks is for the result to be < LOR.

Sub-Matrix: Soil/Solid

Analyte	CAS Number Method	LOR	Unit	Result	Qualifier
Speciated Metals (QCLot: 1479163)					
Methylmercury (as MeHg)	22967-92-6 E538	0.05	μg/kg	<0.050	

Page 3 of 3 Work Order: WT2413665

Client Paracel Laboratories Ltd

Project 2420225



Laboratory Control Sample (LCS) Report

A Laboratory Control Sample (LCS) is an analyte-free matrix that has been fortified (spiked) with test analytes at known concentration and processed in an identical manner to test samples. LCS results are expressed as percent recovery, and are used to monitor and control test method accuracy and precision, independent of test sample matrix.

Sub-Matrix: Soil/Solid					Laboratory Control Sample (LCS) Report									
	Spike	Recovery (%)												
Analyte	CAS Number	Method	LOR	Unit	Target Concentration	LCS	Low	High	Qualifier					
Speciated Metals (QCLot: 1479163)														
Methylmercury (as MeHg)	22967-92-6	E538	0.05	μg/kg	10 μg/kg	92.1	70.0	130						

Reference Material (RM) Report

A Reference Material (RM) is a homogenous material with known and well-established analyte concentrations. RMs are processed in an identical manner to test samples, and are used to monitor and control the accuracy and precision of a test method for a typical sample matrix. RM results are expressed as percent recovery of the target analyte concentration. RM targets may be certified target concentrations provided by the RM supplier, or may be ALS long-term mean values (for empirical test methods).

Sub-Matrix:					Reference Material (RM) Report									
			RM Target	Recovery (%)	Recovery L	imits (%)								
Laboratory sample ID	Reference Material ID	Analyte	CAS Number	Method	Concentration	RM	Low	High	Qualifier					
Speciated Metals	(QCLot: 1479163)													
QC-1479163-003	RM	Methylmercury (as MeHg)	22967-92-6	E538	14.8 μg/kg	108	70.0	130						

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

Subcontract Order

SENDING LABORATORY:

Paracel Laboratories Ltd.

300-2319 St. Laurent Blvd. Ottawa, ON K1G 4J8 Phone: 613-731-9577

Fax: 613-731-9064

Date Requested:

Project Number:

Submitted By:

15-May-24

2420225 Sarah Scullion RECEIVING LABORATORY:

ALS Laboratory Group (Vancouver)

8081 Lougheed Highway Burnaby, BC V5A 1W9 Phone: (604) 253-4188

Fax:

INVOICE TO:

Paracel Laboratories Ltd.

300-2319 St. Laurent Blvd. Ottawa, ON K1G 4J8

Phone: 613-731-9577

Fax: 613-731-9064

Required Regulation	Real 53
Turnaround Time	Standard

Sample ID	Matrix	Analyses Requested:	Sampled	Comments
BH4-24-SS3	Soil	Methyl Mercury - soil	09-May-24 09:00	

BB 05/29/24

Environmental Division Waterloo Work Order Reference



Telephone: +1 519 886 6910

46 Sublet

Please email all results to mfoto@paracellabs.com, dbloom@paracellabs.com, drobertson@paracellabs.com

Temperature prior to Shipping: /

13.5°C -> ICE PACK

12:10





nt 6hd. IG 4.JB labs.com Paracel Order Number (Lab Use Only)

2420225

Chain Of Custody (Lab Use Only)

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L	Paterson Group Inc.			Proje	ect Ref:	PE6422									Pa	ge 1	of 2		
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16	elephone: 613-226-7381			1		kmunch@paterso							-	Requ					
	REG 153/04 REG 406/19	Other Regulation	T	_				72.12	- 1 - 20,	(* - :	20	- 177 85	0000100	er ges	AND SEC		y 1 3 Ar		
	Table 1 Res/Park Med/Fine	REG 558 PWQO				S (Soil/Sed.) GW (G Water) SS (Storm/S						Re	equired Analysis						
	Table 2 Ind/Comm Coarse	□ CCME □ MISA		300 (30		Paint) A (Air) O (O		ВТЕХ	1200		(43)				700	1915		-11	
	Table 3 Agri/Other	SU - Sani SU - Storm		\$2															
	Table	Mun:	l l e l					4+			by ICP								
L	For RSC: Yes No	Other:	ě	Air Volume	Cont			F1-F4	s	, o	ls by			(HWS)					
H	Sample ID/Location Name		Matrix	Air	# of	Date	Time	움	VOCs	PAHs	Metals	Ρ̈́Ξ	S Z	B (H)	표	<u>임</u>	SAR		
\vdash	1 BH1-24-AU1				2	5/8/2024		V	$\overline{\Box}$	V	V	V	V		$\bar{\Box}$	V	V		Г
2			S		2	5/8/2024		V	Ħ	V	V	~	~		H	7	~	H	
3	BH1-24-SS5		s		2	5/8/2024		V	$\overline{\Box}$	V	V	~	7	H	H	7	V	H	늗
4	BH2-24-AU1		s		2	5/8/2024		7	片	7	V	7	7	H	V	7	V	H	H
5	BH2-24-SS3		s		2	5/8/2024			님	V	7		·	H	H	7		ዙ	F
6	BH3-24-AU1		s		2	5/9/2024				V	V	7	V	H	H	·	V	H	
7	BH3-24-SS5		S		2	5/9/2024		<u> </u>	\vdash	H		H	H	H	~	V	<u> </u>	믐	
8	BH4-24-AU1		s		2	5/9/2024		V	\vdash				<u> </u>	H		V	\vdash	H	늗
9	BH4-24-SS2		s		2	5/9/2024			님	V	·		H	믐	H	H	<u> </u>	屵	F
10	BH4-24-SS3		s					<u> </u>	닉	<u> </u>	_		~		닏	_	_	닏	
omn	HOLD - BH4-24-SS3				2	5/9/2024				\Box	Q	7	Ш	Ш	Ш	Ш	Ш	Ш	
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elinq	inquiched B. (b.)						9	SP 98			2.000		ified by: S						
ate/	Time: 5/13/2014						Date/Time: Ho	414,	202	14	_		e/Time; 5 May 24 0928						
ain o	Time: 5/13/2024 f Custody (Riank) visx	Temperature:				°C	Temperature:	27 pH Veri				H Verified: AAA by:							
						Davidson 4.0													





Chain Of Custody Paracel Order Number (Lab Use Only) (Lab Use Only) 2420225

et					/											
Client Name: Paterson Group Inc.		Proje	ct Ref: F	PE6422									Pa	ge 2	of 2	_
Contact Name: Jeremy Camposarcone		Quote	#:									-	Turna	roun	d Tim	ie
Address: 9 AURIGA DRIVE		PO #:	6	60196								1 day				3 day
OTTAWA ON K2E 7T9		E-mai	l: je	camposarcone@p	atersongroup.ca	1						2 day				× Regular
Telephone: 613-226-7381			kmunch@patersongroup.ca								Date Required:					
REG 153/04 REG 406/19 Other Regulation	Τ,	fately 7		r (c-11/c-1) curio			15,	, ,		of M		19-5				
☐ Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558 ☐ PWQO				S (Soil/Sed.) GW (G Vater) SS (Storm/Sa					Required Analysis							
□ Table 2 □ Ind/Comm □ Coarse □ CCME □ MISA			P (Paint) A (Air) O (Other)			втех								Г	Г	
☐ Table 3 ☐ Agri/Other ☐ SU - Sani ☐ SU - Storm			5			+ BT			_							
☐ Table Mun:		Je J	Containers	Sample	Taken	F1-F4 +			y ICP							
For RSC: Yes No Other:	Matrix	Air Volume	Con			E C	SS	φ	Metals by		_	B (HWS)				
Sample ID/Location Name	Σ	Air	# of	Date	Time	오	VOCs	PAHs	Met	롸	CrV	B (F	된	E	SAR	
1 BH4-24-SS6	S		2	5/9/2024		v			v					~	~	
2 BH5-24-AU1	S		2	5/9/2024		v		'	V	V	~		V	~	~	
3 BH5-24-SS2(BOTTOM)	S		23	5/9/2024		V			V					V	V	
4 DUP1	s		2	5/8/2024		~	$\overline{\Box}$	~	~	~	V			V	V	
5 DUP2	s		2	5/9/2024		~	$\overline{\Box}$	靣	~	$\overline{\Box}$	而	而	$\overline{\Box}$	V	V	
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9						H	Ħ	Ħ	H	\exists	\exists	H	H	H	H	
10						Ħ	Ħ	Ħ	H	H		H	H	H	H	
Comments:										Metho	d of De	livery:		Ш		لـــاالــــا
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delinquished By (Sign): Received By Dr	iver/De	pot:			Received at Lab:					Verifie	d By:			unu		- (
elinquished By (Princ Jeremy Camposarcone Date/Time:					Date/Time: (2.6	1-					011	0	220)
Temperature:		M			Mai						Verified: 11 A By:					
ain of Custodu (Blank) ylsv		494.40		Revision 4.0		2.7		A COLOR				MA			Market.	



300-2319 St. Laurent Blvd. Ottawa, Ontario K1G 4J8

p: 1-800-749-1947 e: paracel@paracellabs.com

www.paracellabs.com

Certificate of Analysis

Paterson Group Consulting Engineers

28 Concourse Gate, Unit 1 Phone: (613) 226-7381 Nepean, ON K2E 7T7 Fax: (613) 226-6344

Attn: Eric Leveque

Client PO: 10564 Report Date: 17-May-2011 Project: PE2289 Order Date: 16-May-2011

Custody: 84729 Order #: 1121005

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

Paracel ID Client ID 1121005-01 BH1-GW1

Approved By:

Mark Froto

Mark Foto, M.Sc. For Dale Robertson, BSc Laboratory Director



Certificate of AnalysisReport Date: 17-May-2011Client: Paterson Group Consulting EngineersOrder Date: 16-May-2011

Client PO: 10564 Project Description: PE2289

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date Analysis Date
CCME PHC F1	CWS Tier 1 - P&T GC-FID	16-May-11 16-May-11
CCME PHC F2 - F4	CWS Tier 1 - GC-FID, extraction	16-May-11 17-May-11
VOCs	EPA 624 - P&T GC-MS	16-May-11 16-May-11

NIAGARA FALLS



Client PO: 10564

Order #: 1121005

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Project Description: PE2289

Report Date: 17-May-2011 Order Date:16-May-2011

Client FO. 10304		Froject Descrip	1011.1 L2203		
	Client ID:	BH1-GW1	-	-	-
	Sample Date:	16-May-11	-	-	-
P	Sample ID:	1121005-01	-	-	-
William	MDL/Units	Water	-	-	-
Volatiles	5.0		1		1
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	-	-	-
Chloroethane	1.0 ug/L	<1.0	-	-	-
Chloroform	0.5 ug/L	<0.5	-	-	-
Chloromethane	3.0 ug/L	<3.0	-	-	-
Dibromochloromethane	0.5 ug/L	<0.5	-	-	-
Dichlorodifluoromethane	1.0 ug/L	<1.0	-	-	-
1,2-Dibromoethane	0.2 ug/L	<0.2	-	-	-
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-Dichloroethylene, total	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	-	-	-
Hexane	1.0 ug/L	<1.0	-	-	-
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	-	-	-
Methyl Butyl Ketone (2-Hexanone	10.0 ug/L	<10.0	-	-	-
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	-	-	-
· · · · · · · · · · · · · · · · · · ·			l .		I



F3 + F4 PHCs

Order #: 1121005

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Report Date: 17-May-2011 Order Date:16-May-2011

Client PO: 10564 Project Description: PE2289 Client ID: BH1-GW1 Sample Date: 16-May-11 1121005-01 Sample ID: Water MDL/Units 0.5 ug/L Styrene < 0.5 0.5 ug/L 1,1,1,2-Tetrachloroethane < 0.5 0.5 ug/L 1,1,2,2-Tetrachloroethane < 0.5 0.5 ug/L Tetrachloroethylene < 0.5 0.5 ug/L Toluene < 0.5 _ 0.5 ug/L 1,2,4-Trichlorobenzene <0.5 1,1,1-Trichloroethane 0.5 ug/L <0.5 0.5 ug/L 1,1,2-Trichloroethane < 0.5 0.5 ug/L Trichloroethylene < 0.5 1.0 ug/L Trichlorofluoromethane <1.0 _ -0.5 ug/L 1,2,4- Trimethylbenzene < 0.5 0.5 ug/L 1,3,5-Trimethylbenzene < 0.5 _ _ 0.5 ug/L Vinyl chloride < 0.5 _ 0.5 ug/L m,p-Xylenes < 0.5 0.5 ug/L o-Xylene < 0.5 0.5 ug/L Xylenes, total < 0.5 Surrogate 4-Bromofluorobenzene 94.3% -Dibromofluoromethane Surrogate 95.5% Surrogate Toluene-d8 101% _ _ **Hydrocarbons** 25 ug/L F1 PHCs (C6-C10) <25 _ -100 ug/L F2 PHCs (C10-C16) <100 100 ug/L F3 PHCs (C16-C34) <100 100 ug/L F4 PHCs (C34-C50) <100 125 ug/L F1 + F2 PHCs <125 _ _ 200 ug/L

<200

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Certificate of Analysis

Client: Paterson Group Consulting Engineers
Client PO: 10564

Report Date: 17-May-2011 Order Date:16-May-2011

Project Description: PE2289

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
•									
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
F2 PHCs (C10-C16)	ND	100	ug/L						
F3 PHCs (C16-C34)	ND	100	ug/L						
F4 PHCs (C34-C50)	ND	100	ug/L						
Volatiles									
Acetone	ND	5.0	ug/L						
Benzene	ND	0.5	ug/L						
Bromodichloromethane	ND	0.5	ug/L						
Bromoform	ND	0.5	ug/L						
Bromomethane	ND	0.5	ug/L						
Carbon Tetrachloride	ND	0.2	ug/L						
Chlorobenzene	ND	0.5	ug/L						
Chloroethane	ND	1.0	ug/L						
Chloroform	ND	0.5	ug/L						
Chloromethane	ND	3.0	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dibromoethane	ND	0.2 0.5	ug/L						
1,2-Dichlorobenzene 1,3-Dichlorobenzene	ND ND	0.5 0.5	ug/L						
1,4-Dichlorobenzene	ND	0.5	ug/L ug/L						
1,1-Dichloroethane	ND	0.5	ug/L ug/L						
1,2-Dichloroethane	ND	0.5	ug/L 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-Dichloroethylene, total	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						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Butyl Ketone (2-Hexanone)	ND	10.0	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 1,1,1,2-Tetrachloroethane	ND ND	0.5 0.5	ug/L						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L ug/L						
Tetrachloroethylene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
1,2,4-Trichlorobenzene	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						
1,2,4- Trimethylbenzene	ND	0.5	ug/L						
1,3,5-Trimethylbenzene	ND	0.5	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	34.3		ug/L		107	50-140			
	30.2		ug/L		94.4	50-140			
Surrogate: Dibromofluoromethane Surrogate: Toluene-d8	32.5		ug/L ug/L		102	50-140 50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 10564 Project Description: PE2289

Report Date: 17-May-2011 Order Date:16-May-2011

Method Quality Control: Duplicate

Hydrocarbons				Result	%REC	Limit	RPD	Limit	Notes
								_	
F1 PHCs (C6-C10)	ND	25	ug/L	ND				30	
Volatiles			- 3						
	ND	F 0		ND				00	
Acetone Benzene	ND ND	5.0 0.5	ug/L ug/L	ND ND				30 30	
Bromodichloromethane	ND	0.5	ug/L ug/L	ND				30	
Bromoform	ND	0.5	ug/L ug/L	ND				30	
Bromomethane	ND	0.5	ug/L ug/L	ND				30	
Carbon Tetrachloride	ND	0.2	ug/L	ND				30	
Chlorobenzene	ND	0.5	ug/L	ND				30	
Chloroethane	ND	1.0	ug/L	ND				30	
Chloroform	ND	0.5	ug/L	ND				30	
Chloromethane	ND	3.0	ug/L	ND				30	
Dibromochloromethane	ND	0.5	ug/L	ND				30	
Dichlorodifluoromethane	ND	1.0	ug/L	ND				30	
1,2-Dibromoethane	ND	0.2	ug/L	ND				30	
1,2-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,1-Dichloroethane	ND	0.5	ug/L	ND				30	
1,2-Dichloroethane	ND	0.5	ug/L	ND				30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND				30	
cis-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
trans-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
1,2-Dichloropropane	ND	0.5	ug/L	ND				30	
cis-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30 30	
trans-1,3-Dichloropropylene Ethylbenzene	ND ND	0.5 0.5	ug/L ug/L	ND ND				30	
Hexane	ND	1.0	ug/L ug/L	ND				30	
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L ug/L	ND				30	
Methyl Butyl Ketone (2-Batanone)	ND	10.0	ug/L ug/L	ND				30	
Methyl Isobutyl Ketone	ND	5.0	ug/L	ND				30	
Methyl tert-butyl ether	ND	2.0	ug/L	ND				30	
Methylene Chloride	ND	5.0	ug/L	ND				30	
Styrene	ND	0.5	ug/L	ND				30	
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L	ND				30	
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L	ND				30	
Tetrachloroethylene	ND	0.5	ug/L	ND				30	
Toluene	ND	0.5	ug/L	ND				30	
1,2,4-Trichlorobenzene	ND	0.5	ug/L	ND				30	
1,1,1-Trichloroethane	ND	0.5	ug/L	ND				30	
1,1,2-Trichloroethane	ND	0.5	ug/L	ND				30	
Trichloroethylene	ND	0.5	ug/L	ND				30	
Trichlorofluoromethane	ND	1.0	ug/L	ND				30	
1,2,4- Trimethylbenzene	ND	0.5	ug/L	ND				30	
1,3,5-Trimethylbenzene	ND	0.5	ug/L	ND				30	
Vinyl chloride	ND	0.5	ug/L	ND				30	
m,p-Xylenes	ND	0.5	ug/L	ND				30 30	
o-Xylene Surrogate: 4-Bromofluorobenzene	ND <i>33.4</i>	0.5	ug/L	ND <i>ND</i>	104	50-140		30	
Surrogate: Dibromofluoromethane	33.4 31.5		ug/L	ND ND	98.5	50-140 50-140			
Surrogate: Toluene-d8	31.5 31.4		ug/L	ND ND	98.1	50-140 50-140			
Surroyate. Toluelle-uo	31.4		ug/L	ND	90.1	30-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 10564 Project Description: PE2289 Report Date: 17-May-2011 Order Date:16-May-2011

Method Quality Control: Snike

Analyte	Doorst	Reporting	1.19	Source	0/850	%REC	DDC	RPD	NI-+-
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1980	25	ug/L	ND	98.8	68-117			
F2 PHCs (C10-C16)	1260	100	ug/L	ND	79.0	61-129			
F3 PHCs (C16-C34)	4180	100	ug/L	ND	104	61-129			
F4 PHCs (C34-C50)	3000	100	ug/L	ND	125	61-129			
Volatiles									
Acetone	71.9	5.0	ug/L	ND	71.9	50-140			
Benzene	30.5	0.5	ug/L	ND	76.2	60-130			
Bromodichloromethane	37.8	0.5	ug/L	ND	94.4	60-130			
Bromoform	39.2	0.5	ug/L	ND	98.0	60-130			
Bromomethane	22.5	0.5	ug/L	ND	56.2	50-140			
Carbon Tetrachloride	39.6	0.2	ug/L	ND	99.1	60-130			
Chlorobenzene	32.8	0.5	ug/L	ND	82.1	60-130			
Chloroethane	24.4	1.0	ug/L	ND	61.0	50-140			
Chloroform	36.6	0.5	ug/L	ND	91.5	60-130			
Chloromethane	23.0	3.0	ug/L	ND	57.5	50-140			
Dibromochloromethane	39.0	0.5	ug/L	ND	97.4	60-130			
Dichlorodifluoromethane	34.1	1.0	ug/L	ND	85.3	50-140			
1,2-Dibromoethane	30.4	0.2	ug/L	ND	75.9	60-130			
1,2-Dichlorobenzene	37.8	0.5	ug/L	ND	94.4	60-130			
1,3-Dichlorobenzene	37.4	0.5	ug/L	ND	93.6	60-130			
1,4-Dichlorobenzene	37.5	0.5	ug/L	ND	93.8	60-130			
1,1-Dichloroethane	28.0	0.5	ug/L	ND	70.0	60-130			
1,2-Dichloroethane	37.8	0.5	ug/L	ND	94.4	60-130			
1,1-Dichloroethylene	31.2	0.5	ug/L	ND	78.0	60-130			
cis-1,2-Dichloroethylene	31.5	0.5	ug/L	ND	78.7	60-130			
trans-1,2-Dichloroethylene	33.5	0.5	ug/L	ND	83.7	60-130			
1,2-Dichloropropane	27.7	0.5	ug/L	ND	69.2	60-130			
cis-1,3-Dichloropropylene	33.7	0.5	ug/L	ND	84.3	60-130			
trans-1,3-Dichloropropylene	36.5	0.5	ug/L	ND	91.2	60-130			
Ethylbenzene	30.9	0.5	ug/L	ND	77.3	60-130			
Hexane	29.0	1.0	ug/L	ND	72.6	60-130			
Methyl Ethyl Ketone (2-Butanone)	50.9	5.0	ug/L	ND	50.9	50-140			
Methyl Butyl Ketone (2-Hexanone)	56.0	10.0	ug/L	ND	56.0	50-140			
Methyl Isobutyl Ketone	59.3	5.0	ug/L	ND	59.3	50-140			
Methyl tert-butyl ether	129	2.0	ug/L	ND	129	50-140			
Methylene Chloride	32.6	5.0	ug/L	ND	81.4	60-130			
Styrene	29.1	0.5	ug/L ug/L	ND	72.8	60-130			
1,1,1,2-Tetrachloroethane	38.6	0.5	ug/L	ND	96.6	60-130			
1,1,2,2-Tetrachloroethane	30.3	0.5	ug/L ug/L	ND	75.8	60-130			
Tetrachloroethylene	36.6	0.5	ug/L ug/L	ND	91.5	60-130			
Toluene	32.6	0.5	ug/L ug/L	ND	81.6	60-130			
1,2,4-Trichlorobenzene	34.1	0.5	ug/L ug/L	ND	85.4	60-130			
1,1,1-Trichloroethane	40.6	0.5	ug/L ug/L	ND	101	60-130			
1,1,2-Trichloroethane	32.3	0.5	ug/L ug/L	ND	80.8	60-130			
Trichloroethylene	34.5	0.5	ug/L ug/L	ND	86.3	60-130			
Trichlorofluoromethane	41.5	1.0	ug/L ug/L	ND	104	60-130			
1,2,4- Trimethylbenzene	34.9	0.5	ug/L ug/L	ND	87.2	60-130			
1,3,5-Trimethylbenzene	39.4	0.5	ug/L ug/L	ND	98.6	60-130			
	23.7	0.5		ND	96.6 59.3	50-130			
Vinyl chloride			ug/L						
m,p-Xylenes	65.4	0.5 0.5	ug/L	ND	81.7	60-130			
o-Xylene	32.5	0.5	ug/L	ND	81.2	60-130			
Surrogate: 4-Bromofluorobenzene	29.9		ug/L		93.6	50-140 50-140			
Surrogate: Dibromofluoromethane	31.8		ug/L		99.3	50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 10564 Project Description: PE2289 Report Date: 17-May-2011 Order Date:16-May-2011

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
									-

25.2 78.6 Surrogate: Toluene-d8 ug/L 50-140

NIAGARA FALLS



Client: Paterson Group Consulting Engineers

Client PO: 10564 Project Description: PE2289

Report Date: 17-May-2011 Order Date:16-May-2011

Order #: 1121005

Sample and QC Qualifiers Notes

None

Sample Data Revisions

None

Work Order Revisions/Comments:

None

Other Report Notes:

n/a: not applicable

MDL: Method Detection Limit

Source Result: Data used as source for matrix and duplicate samples

%REC: Percent recovery.

RPD: Relative percent difference.

CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.
- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.

OTTAWA



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Chain of Custody
(lab use only)

Nº 84729

OTTAWA @ NIAGARA FALLS @ MISSISSAUGA @ SARNIA

Client Name:	ONOGA	SOAF	TIVIA			Reg. Drinking W	ater	e; p	aracei@para	cellabs.com	1	N ·	0412	Ü
Client Name: Paterson Group	Project	Ref: P	£228	9		Waterworks Name	e:					Page	⊥ of _[3.
Contact Name: Ecic 19 Vo Gue	Quote	#				Waterworks Num	ber:							The .
Address: 28 Concourse Gate, Unit!	PO#	1				Address:	1						Taken by:	
Ottawa ON	E mail	105 Address:									Print N	ame: /	Robinso	01
	ele	Nouress:	2 Qpat	erson	group.ca	After hours Cont	act:				Signatu	ire:	11.	
Telephone: 613 226-7381	Fax:	V			3 1	Public Health Un				1000	ТА	T: NI I day	/ 2-day	I D
Matrix Types: S-Soil/Sed. GW-Ground Water S	W-Surface	Water	SS-Stor	m/Sanit	ary Sewer D	W-Drinking	Water RD	W-Red	pulated D	rinking W	otor D	Doint A	1 2-day	Keg.
☐ O. Reg 153 (511) Table <u>3</u> ☐ O. Reg 170/03 ☐ O. Reg 318 ☐ CCME ☐ O. Reg 243/07 ☐ O. Reg 319/08 ☐ Otho	3/08 ☐ Privat	e well	Type of	DW Samp	ole: R = Raw: T	' = Treated; D = er; G = Ground	Distribution	, Ke	guiated D		ired An		AIF U-O	ner
Paracel Order Number			0	10					(3)		T			
1121005	Matrix	Air Volume	Type of Sample	# of Containers	Samı	ole Taken	Free / Combined Chlorine Residual mg/L	00.3	PHC S(F,-F					
Sample ID / Location Name		330	L.	#	Date	Time	Ch.	7	PH					
1 BHI-GWI	6-W			3	May16/1	1 gam		V	V					
2					1100/10/1	dan		0	-					-
3												_		
4														
5														
6	-													
7														
8														
9														
10														
Comments:								Pres	servation V	erification:	pH_L	Tem	perature /	0,60
Relinquished By (Print & Sign): 7. Robinson							Lab Use Onl			16	DU	2		
1/1/	Receive					Received	11.	/		Verified	1	71	/	
Date/Time: May 16 2011 Gam	Driver/	menderal kind dan basasas				at Lab:	10	1	1	By:	1/1	0		

ChainOfCustody Rev 2.0, January 2010

Walk-in



Head Office

300-2319 St. Laurent Blvd. Ottawa, Ontario K1G 4J8

p: 1-800-749-1947

e: paracel@paracellabs.com

www.paracellabs.com

Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South

Nepean, ON K2E 7J5

Phone: (613) 226-7381

Fax: (613) 226-6344

Attn: Mark D'Arcy

Client PO: 12261 Report Date: 24-Jul-2012 Project: PE2709 Order Date: 20-Jul-2012

Custody: 94393 Order #: 1229249

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

Paracel ID	Client ID
1229249-01	BH1-GW1
1229249-02	BH2-GW1
1229249-03	BH4-GW1

Approved By:

Mark Froto

Mark Foto, M.Sc. For Dale Robertson, BSc Laboratory Director



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 12261 Project Description: PE2709

Report Date: 24-Jul-2012 Order Date: 20-Jul-2012

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date Analysis Date
CCME PHC F1	CWS Tier 1 - P&T GC-FID	20-Jul-12 23-Jul-12
CCME PHC F2 - F4	CWS Tier 1 - GC-FID, extraction	20-Jul-12 21-Jul-12
PAHs by GC-MS, standard scan	EPA 625 - GC-MS, extraction	23-Jul-12 24-Jul-12
VOCs	EPA 624 - P&T GC-MS	20-Jul-12 23-Jul-12

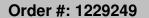


Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 12261 Project Description: PE2709 Report Date: 24-Jul-2012 Order Date:20-Jul-2012

	Client ID: Sample Date: Sample ID: MDL/Units	BH1-GW1 20-Jul-12 1229249-01 Water	BH2-GW1 20-Jul-12 1229249-02 Water	BH4-GW1 20-Jul-12 1229249-03 Water	- - -
Volatiles					
Acetone	5.0 ug/L	82.4	691	40.6	-
Benzene	0.5 ug/L	<0.5	2.4	1.1	-
Bromodichloromethane	0.5 ug/L	<0.5	<0.5	3.8	-
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	-
Chloroethane	1.0 ug/L	<1.0	<1.0	<1.0	-
Chloroform	0.5 ug/L	16.1	14.4	21.8	-
Chloromethane	3.0 ug/L	<3.0	<3.0	<3.0	-
Dibromochloromethane	0.5 ug/L	<0.5	<0.5	<0.5	-
Dichlorodifluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	-
1,2-Dibromoethane	0.2 ug/L	<0.2	<0.2	<0.2	-
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-Dichloroethylene, total	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	1.0	<0.5	-
Hexane	1.0 ug/L	2.1	2.2	<1.0	-
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	10.7	8.0	8.4	-
Methyl Butyl Ketone (2-Hexanone	10.0 ug/L	<10.0	<10.0	<10.0	-
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	-





Client: Paterson Group Consulting Engineers

Client PO: 12261 Project Description: PE2709

Report Date: 24-Jul-2012 Order Date: 20-Jul-2012

Client PO: 12261		Project Description			
	Client ID: Sample Date:	BH1-GW1 20-Jul-12	BH2-GW1 20-Jul-12	BH4-GW1 20-Jul-12	-
	Sample ID:	1229249-01	1229249-02	1229249-03	-
	MDL/Units	Water	Water	Water	-
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	6.2	5.8	5.6	-
1,2,4-Trichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-
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	-
1,3,5-Trimethylbenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
Vinyl chloride	0.5 ug/L	<0.5	<0.5	<0.5	-
m,p-Xylenes	0.5 ug/L	<0.5	1.2	<0.5	-
o-Xylene	0.5 ug/L	<0.5	1.6	<0.5	-
Xylenes, total	0.5 ug/L	<0.5	2.8	<0.5	-
4-Bromofluorobenzene	Surrogate	110%	117%	109%	-
Dibromofluoromethane	Surrogate	106%	110%	106%	-
Toluene-d8	Surrogate	98.4%	94.8%	97.2%	-
Hydrocarbons					_
F1 PHCs (C6-C10)	25 ug/L	45	43	<25	-
F2 PHCs (C10-C16)	100 ug/L	<287	<100	<100	-
F3 PHCs (C16-C34)	100 ug/L	<287	<100	<100	-
F4 PHCs (C34-C50)	100 ug/L	<287	<100	<100	-
Semi-Volatiles					
Acenaphthene	0.05 ug/L	-	<0.05	-	-
Acenaphthylene	0.05 ug/L	-	<0.05	-	-
Anthracene	0.01 ug/L	-	<0.01	-	-
Benzo [a] anthracene	0.01 ug/L	-	<0.01	-	-
Benzo [a] pyrene	0.01 ug/L	-	<0.01	-	-
Benzo [b] fluoranthene	0.05 ug/L	-	<0.05	-	-
Benzo [g,h,i] perylene	0.05 ug/L	-	<0.05	-	-
Benzo [k] fluoranthene	0.05 ug/L	-	<0.05	-	-
Biphenyl	0.05 ug/L	-	0.12	-	-
Chrysene	0.05 ug/L	-	<0.05	-	-
Dibenzo [a,h] anthracene	0.05 ug/L	-	<0.05	-	-
			•		



Client PO: 12261

Order #: 1229249

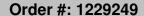
Certificate of Analysis

Client: Paterson Group Consulting Engineers

Project Description: PE2709

Report Date: 24-Jul-2012 Order Date: 20-Jul-2012

	Client ID: Sample Date: Sample ID: MDL/Units	BH1-GW1 20-Jul-12 1229249-01 Water	BH2-GW1 20-Jul-12 1229249-02 Water	BH4-GW1 20-Jul-12 1229249-03 Water	- - -
Fluoranthene	0.01 ug/L	-	<0.01	-	-
Fluorene	0.05 ug/L	-	0.06	-	-
Indeno [1,2,3-cd] pyrene	0.05 ug/L	-	<0.05	-	-
1-Methylnaphthalene	0.05 ug/L	-	0.33	-	-
2-Methylnaphthalene	0.05 ug/L	-	0.40	-	-
Methylnaphthalene (1&2)	0.10 ug/L	-	0.73	-	-
Naphthalene	0.05 ug/L	-	0.82	-	-
Phenanthrene	0.05 ug/L	-	<0.05	-	-
Pyrene	0.01 ug/L	-	<0.01	-	-
2-Fluorobiphenyl	Surrogate	-	73.4%	-	-
Terphenyl-d14	Surrogate	-	80.8%	-	-





Client: Paterson Group Consulting Engineers

Client PO: 12261 Project Description: PE2709

Report Date: 24-Jul-2012 Order Date: 20-Jul-2012

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
F2 PHCs (C10-C16)	ND	100	ug/L						
F3 PHCs (C16-C34)	ND	100	ug/L						
F4 PHCs (C34-C50)	ND	100	ug/L						
Semi-Volatiles									
Acenaphthene	ND	0.05	ug/L						
Acenaphthylene	ND	0.05	ug/L						
Anthracene	ND	0.01	ug/L						
Benzo [a] anthracene	ND	0.01	ug/L						
Benzo [a] pyrene	ND	0.01	ug/L						
Benzo [b] fluoranthene	ND	0.05	ug/L						
Benzo [g,h,i] perylene	ND	0.05	ug/L						
Benzo [k] fluoranthene	ND	0.05	ug/L						
Biphenyl	ND	0.05	ug/L						
Chrysene	ND ND	0.05 0.05	ug/L						
Dibenzo [a,h] anthracene Fluoranthene	ND ND	0.05	ug/L ug/L						
Fluorene	ND ND	0.05	ug/L ug/L						
Indeno [1,2,3-cd] pyrene	ND	0.05	ug/L						
1-Methylnaphthalene	ND	0.05	ug/L						
2-Methylnaphthalene	ND	0.05	ug/L						
Methylnaphthalene (1&2)	ND	0.10	ug/L						
Naphthalene	ND	0.05	ug/L						
Phenanthrene	ND	0.05	ug/L						
Pyrene	ND	0.01	ug/L						
Surrogate: 2-Fluorobiphenyl	19.0		ug/L		95.2	50-140			
Surrogate: Terphenyl-d14	18.0		ug/L		90.0	50-140			
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						
Chloroethane	ND	1.0	ug/L						
Chloroform Chloromethane	ND ND	0.5 3.0	ug/L ug/L						
Dibromochloromethane	ND ND	0.5	ug/L ug/L						
Dichlorodifluoromethane	ND ND	1.0	ug/L ug/L						
1,2-Dibromoethane	ND	0.2	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-Dichloroethylene, total	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						
Hexane	ND	1.0	ug/L						



Certificate of Analysis

Surrogate: 4-Bromofluorobenzene

Surrogate: Dibromofluoromethane

Surrogate: Toluene-d8

Client: Paterson Group Consulting Engineers

Client PO: 12261 Project Description: PE2709

Report Date: 24-Jul-2012 Order Date: 20-Jul-2012

Method Quality Control: Blank Reporting %REC RPD Source Analyte Result RPD Limit Units Result %REC Limit Limit Notes Methyl Ethyl Ketone (2-Butanone) ND 5.0 ug/L Methyl Butyl Ketone (2-Hexanone) ND 10.0 ug/L Methyl Isobutyl Ketone ND 5.0 ug/L Methyl tert-butyl ether ug/L ND 2.0 Methylene Chloride ND 5.0 ug/L Styrene ND 0.5 ug/L 1,1,1,2-Tetrachloroethane ND 0.5 ug/L ug/L 1,1,2,2-Tetrachloroethane ND 0.5 Tetrachloroethylene ND 0.5 ug/L Toluene ND 0.5 ug/L 1,2,4-Trichlorobenzene 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 ug/L Trichlorofluoromethane ND 1.0 1,3,5-Trimethylbenzene 0.5 ug/L ND Vinyl chloride ND 0.5 ug/L m,p-Xylenes o-Xylene ug/L ND 0.5 ND 0.5 ug/L Xylenes, total ND ug/L

ug/L

ug/L

ug/L

37.9

25.4

34.7

50-140

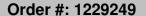
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118

79.4

108





1,3,5-Trimethylbenzene

Surrogate: Toluene-d8

Surrogate: 4-Bromofluorobenzene

Surrogate: Dibromofluoromethane

Vinyl chloride

m,p-Xylenes

o-Xylene

Client: Paterson Group Consulting Engineers

Project Description: PE2709 Client PO: 12261

Report Date: 24-Jul-2012 Order Date:20-Jul-2012

Method Quality Control:									
Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L	ND				30	
Volatiles		-	- 3						
	ND	5 0		ND				00	
Acetone	ND	5.0	ug/L	ND				30	
Benzene	ND	0.5	ug/L	ND				30	
Bromodichloromethane	ND	0.5	ug/L	ND				30	
Bromoform	ND	0.5	ug/L	ND				30	
Bromomethane	ND	0.5	ug/L	ND				30	
Carbon Tetrachloride	ND	0.2	ug/L	ND				30	
Chlorobenzene	ND	0.5	ug/L	ND				30	
Chloroethane	ND	1.0	ug/L	ND				30	
Chloroform	ND	0.5	ug/L	ND				30	
Chloromethane	ND	3.0	ug/L	ND				30	
Dibromochloromethane	ND	0.5	ug/L	ND				30	
Dichlorodifluoromethane	ND	1.0	ug/L	ND				30	
1,2-Dibromoethane	ND	0.2	ug/L	ND				30	
1,2-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,1-Dichloroethane	ND	0.5	ug/L	ND				30	
1,2-Dichloroethane	ND	0.5	ug/L	ND				30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND				30	
cis-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
trans-1,2-Dichloroethylene	ND	0.5	ug/L	ND				30	
1,2-Dichloropropane	ND	0.5	ug/L	ND				30	
cis-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30	
trans-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30	
Ethylbenzene	ND	0.5	ug/L	ND				30	
Hexane	ND	1.0	ug/L	ND				30	
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L	ND				30	
Methyl Butyl Ketone (2-Hexanone)	ND	10.0	ug/L	ND				30	
Methyl Isobutyl Ketone	ND	5.0	ug/L	ND				30	
Methyl tert-butyl ether	ND	2.0	ug/L	ND				30	
Methylene Chloride	ND	5.0	ug/L	ND				30	
Styrene	ND	0.5	ug/L	ND				30	
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L	ND				30	
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L	ND				30	
Tetrachloroethylene	ND	0.5	ug/L	ND				30	
Toluene	ND	0.5	ug/L	ND				30	
1,2,4-Trichlorobenzene	ND	0.5	ug/L	ND				30	
1,1,1-Trichloroethane	ND	0.5	ug/L	ND				30	
1,1,2-Trichloroethane	ND	0.5	ug/L	ND				30	
Trichloroethylene	ND	0.5	ug/L	ND				30	
Trichlorofluoromethane	ND	1.0	ug/L	ND				30	
1 3 5-Trimethylhenzene	ND	0.5	ua/l	ND				30	

ND

ND

ND

ND

38.4

35.1

35.3

0.5

0.5

0.5

0.5

ug/L

ug/L

ug/L

ug/L

ug/L

ug/L

ug/L

ND

ND

ND

ND

ND

ND

ND

120

110

110

50-140

50-140

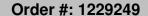
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30

30

30

30





Client: Paterson Group Consulting Engineers

Client PO: 12261 Project Description: PE2709 Report Date: 24-Jul-2012

Order Date:20-Jul-2012

	Spike Result	Reporting	Units	Source	%REC	%REC	RPD	RPD	Notes
Analyte	nesuit	Limit	Offics	Result	/onLU	Limit	NED	Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1910	25	ug/L	ND	95.7	68-117			
F2 PHCs (C10-C16)	1230	100	ug/L	ND	76.7	60-140			
F3 PHCs (C16-C34)	3300	100	ug/L	ND	82.5	60-140			
F4 PHCs (C34-C50)	2390	100	ug/L	ND	99.6	60-140			
Semi-Volatiles									
Acenaphthene	3.44	0.05	ug/L	ND	68.9	50-140			
cenaphthylene	3.49	0.05	ug/L	ND	69.7	50-140			
nthracene	3.55	0.01	ug/L	ND	71.0	50-140			
enzo [a] anthracene	3.88	0.01	ug/L	ND	77.6	50-140			
enzo [a] pyrene	3.44	0.01	ug/L	ND	68.9	50-140			
Senzo [b] fluoranthene	4.66	0.05	ug/L	ND	93.2	50-140			
enzo [g,h,i] perylene	3.22	0.05	ug/L	ND	64.3	50-140			
enzo [k] fluoranthene	4.86	0.05	ug/L	ND	97.3	50-140			
iphenyl	3.54	0.05	ug/L	ND	70.8	50-140			
hrysene	4.10	0.05	ug/L	ND	82.1	50-140			
bibenzo [a,h] anthracene	3.90	0.05	ug/L	ND	78.0	50-140			
luoranthene	4.20	0.01	ug/L	ND	84.1	50-140			
luorene	4.02	0.05	ug/L	ND	80.4	50-140			
ndeno [1,2,3-cd] pyrene	3.04	0.05	ug/L	ND	60.8	50-140			
-Methylnaphthalene	3.13	0.05	ug/L	ND	62.6	50-140			
-Methylnaphthalene	3.44	0.05	ug/L	ND	68.8	50-140			
laphthalene	3.70	0.05	ug/L	ND	74.1	50-140			
Phenanthrene	3.89	0.05	ug/L	ND	77.8	50-140			
lyrene	4.41	0.01	ug/L	ND	88.2	50-140			
Gurrogate: 2-Fluorobiphenyl	17.8	0.0.	ug/L		89.1	50-140			
/olatiles			- 3						
Acetone	75.4	5.0	ug/L	ND	75.4	50-140			
Benzene	43.2	0.5	ug/L	ND	108	50-140			
romodichloromethane	42.0	0.5	ug/L	ND	105	50-140			
romoform	38.1	0.5	ug/L	ND	95.2	50-140			
romomethane	30.1	0.5	ug/L	ND	75.3	50-140			
Carbon Tetrachloride	41.2	0.2	ug/L	ND	103	50-140			
Chlorobenzene	43.1	0.5	ug/L	ND	108	50-140			
Chloroethane	42.2	1.0	ug/L	ND	106	50-140			
Chloroform	43.5	0.5	ug/L ug/L	ND	100	50-140			
Chloromethane	34.1	3.0	ug/L	ND	85.2	50-140			
Dibromochloromethane	40.1	0.5	ug/L ug/L	ND	100	50-140			
Dichlorodifluoromethane	43.5	1.0	ug/L ug/L	ND	100	50-140			
,2-Dibromoethane	36.9	0.2	ug/L ug/L	ND	92.2	50-140			
,2-Dibromoetriane ,2-Dichlorobenzene	32.1	0.5	ug/L ug/L	ND	80.3	50-140			
,3-Dichlorobenzene	32.1	0.5	ug/L ug/L	ND	80.9	50-140			
,4-Dichlorobenzene	33.9	0.5	ug/L ug/L	ND	84.8	50-140			
,1-Dichloroethane	46.0	0.5	ug/L ug/L	ND	04.0 115	50-140			
,2-Dichloroethane	40.8	0.5		ND	102	50-140 50-140			
,1-Dichloroethylene	40.6 27.7	0.5	ug/L ug/L	ND	69.2	50-140 50-140			
•									
is-1,2-Dichloroethylene ans-1,2-Dichloroethylene	35.0 34.6	0.5 0.5	ug/L ug/L	ND ND	87.6 86.4	50-140 50-140			
SUSSI ZELICHIOTOGIOVIANA	34 h	un	11/1/1	INI)	გი.4	50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 12261 Project Description: PE2709 Report Date: 24-Jul-2012 Order Date:20-Jul-2012

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
cis-1,3-Dichloropropylene	33.9	0.5	ug/L	ND	84.8	50-140		•	•
trans-1,3-Dichloropropylene	44.0	0.5	ug/L	ND	110	50-140			
Ethylbenzene	36.4	0.5	ug/L	ND	91.1	50-140			
Hexane	43.3	1.0	ug/L	ND	108	50-140			
Methyl Ethyl Ketone (2-Butanone)	62.5	5.0	ug/L	ND	62.5	50-140			
Methyl Butyl Ketone (2-Hexanone)	67.9	10.0	ug/L	ND	67.9	50-140			
Methyl Isobutyl Ketone	63.7	5.0	ug/L	ND	63.7	50-140			
Methyl tert-butyl ether	72.8	2.0	ug/L	ND	72.8	50-140			
Methylene Chloride	34.3	5.0	ug/L	ND	85.8	50-140			
Styrene	29.9	0.5	ug/L	ND	74.8	50-140			
1,1,1,2-Tetrachloroethane	42.6	0.5	ug/L	ND	107	50-140			
1,1,2,2-Tetrachloroethane	49.5	0.5	ug/L	ND	124	50-140			
Tetrachloroethylene	43.0	0.5	ug/L	ND	108	50-140			
Toluene	40.0	0.5	ug/L	ND	100	50-140			
1,2,4-Trichlorobenzene	43.2	0.5	ug/L	ND	108	50-140			
1,1,1-Trichloroethane	41.4	0.5	ug/L	ND	104	50-140			
1,1,2-Trichloroethane	37.5	0.5	ug/L	ND	93.8	50-140			
Trichloroethylene	36.9	0.5	ug/L	ND	92.4	50-140			
Trichlorofluoromethane	34.8	1.0	ug/L	ND	87.0	50-140			
1,3,5-Trimethylbenzene	28.6	0.5	ug/L	ND	71.5	50-140			
Vinyl chloride	43.7	0.5	ug/L	ND	109	50-140			
m,p-Xylenes	89.0	0.5	ug/L	ND	111	50-140			
o-Xylene	39.6	0.5	ug/L	ND	98.9	50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Report Date: 24-Jul-2012
Order Date: 20-Jul-2012

Client PO: 12261 Project Description: PE2709

Qualifier Notes:

None

Sample Data Revisions

None

Work Order Revisions / Comments:

None

Other Report Notes:

n/a: not applicable

MDL: Method Detection Limit

Source Result: Data used as source for matrix and duplicate samples

%REC: Percent recovery.

RPD: Relative percent difference.

CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.
- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.

6	P	A	R	A	C	ΕL
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Chain of Custody
(Lab Use Only)

Nº 94393

Page _/ of _/

Client Name: VMCLOST GLOUP (NU	Project Reference: PE2744	TAT: Regular 3 Day
Contact Name: HARK - D'ARCV	Quote #	
Address:	PO# 12261	2 Day Day
184 Colonnade food South	Email Address:	Date Required:

Telephone: Criteria: | 10. Reg. 153/04 Table _ | XO. Reg. 153/11 (Current) Table 2 | | RSC Filing | | 0. Reg. 558/00 | | PWQO | | CCME | | SUB (Storm) | | SUB (Sanitary) Municipality:_ []Other: Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) SS (Storm/Sanitary Sewer) P (Paint) A (Air) O (Other) Required Analyses PHCs F1-F4+BTEX Paracel Order Number: # of Containers ICP/MS Air Volume Sample Taken Metals by Matrix PAHS Hg Time Sample ID/Location Name Date 20,2012 1 4 20,2012 X 2 3 4 5 6

Comments: Unelle to recover enough can	ple for 2x vials & 300ml for	BHI-GWI, please first I what you can Method of Delivery:	
Gor RACE 3 VOCS Thank	S. Sial nogents Decount	where nocessary pursean, Walking	
Selinquished By (Print & Sign):	Received by Driver/Depot:	Received at Lat. Venified By	
DEN BOO	Established State St	MC MC	
106GH042 7 90	Date/Time:	Date/Time: / July 20/12 12	1
Date/Time: 2014 20 2012, 1/2	Temperature: °C	Temperature: 99,4°C // O/a pH Verified] By: NA	



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Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South Phone: (613) 226-7381 Nepean, ON K2E 7J5 Fax: (613) 226-6344

Attn: Mark D'Arcy

Client PO: 12263 Report Date: 26-Jul-2012 Project: PE2709 Order Date: 25-Jul-2012

Custody: 94819 Order #: 1230199

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

 Paracel ID
 Client ID

 1230199-01
 BH4-GW2

 1230199-02
 BH2-GW2

Approved By:

Mark Foto

Mark Foto, M.Sc. For Dale Robertson, BSc Laboratory Director



Certificate of Analysis

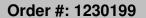
Client: Paterson Group Consulting Engineers

Client PO: 12263 Project Description: PE2709 Report Date: 26-Jul-2012 Order Date:25-Jul-2012

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date Analysis Date
CCME PHC F1	CWS Tier 1 - P&T GC-FID	26-Jul-12 26-Jul-12
CCME PHC F2 - F4	CWS Tier 1 - GC-FID, extraction	26-Jul-12 26-Jul-12
VOCs	EPA 624 - P&T GC-MS	26-Jul-12 26-Jul-12

NIAGARA FALLS 5415 Morning Glory Crt. Niagara Falls, ON L2J 0A3

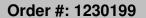




Client: Paterson Group Consulting Engineers

Client PO: 12263 Project Description: PE2709 Report Date: 26-Jul-2012 Order Date:25-Jul-2012

CHETT 6: 12203	о Г	PLIA OMO		I	I
	Client ID:	BH4-GW2 25-Jul-12	BH2-GW2 25-Jul-12	-	-
	Sample Date: Sample ID:	25-Jul-12 1230199-01	1230199-02	-	-
Г	MDL/Units	Water	Water	-	_
Volatiles	WDE/Office				
Acetone	5.0 ug/L	104	1020	-	-
Benzene	0.5 ug/L	<0.5	<0.5	-	-
Bromodichloromethane	0.5 ug/L	<0.5	<0.5	-	-
Bromoform	0.5 ug/L	<0.5	<0.5	-	-
Bromomethane	0.5 ug/L	<0.5	<0.5	-	-
Carbon Tetrachloride	0.2 ug/L	<0.2	<0.2	-	-
Chlorobenzene	0.5 ug/L	<0.5	<0.5	-	-
Chloroethane	1.0 ug/L	<1.0	<1.0	-	-
Chloroform	0.5 ug/L	3.1	9.9	-	-
Chloromethane	3.0 ug/L	<3.0	<3.0	-	-
Dibromochloromethane	0.5 ug/L	<0.5	<0.5	-	-
Dichlorodifluoromethane	1.0 ug/L	<1.0	<1.0	-	-
1,2-Dibromoethane	0.2 ug/L	<0.2	<0.2	-	-
1,2-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	-	-
1,3-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	-	-
1,4-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	-	-
1,1-Dichloroethane	0.5 ug/L	<0.5	<0.5	-	-
1,2-Dichloroethane	0.5 ug/L	<0.5	<0.5	-	-
1,1-Dichloroethylene	0.5 ug/L	<0.5	<0.5	-	-
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	-	-
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	-	-
1,2-Dichloroethylene, total	0.5 ug/L	<0.5	<0.5	-	-
1,2-Dichloropropane	0.5 ug/L	<0.5	<0.5	-	-
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	-	-
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	-	-
1,3-Dichloropropene, total	0.5 ug/L	<0.5	<0.5	-	-
Ethylbenzene	0.5 ug/L	<0.5	<0.5	-	-
Hexane	1.0 ug/L	<1.0	<1.0	-	-
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	26.1	-	-
Methyl Butyl Ketone (2-Hexanone)	10.0 ug/L	<10.0	<10.0	-	-
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	<5.0	-	-
Methyl tert-butyl ether	2.0 ug/L	<2.0	<2.0	-	-
Methylene Chloride	5.0 ug/L	<5.0	<5.0	-	-



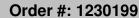


Client: Paterson Group Consulting Engineers

Client PO: 12263 Project Description: PE2709

Report Date: 26-Jul-2012 Order Date: 25-Jul-2012

	Client ID:	BH4-GW2	BH2-GW2	-	-
	Sample Date:	25-Jul-12	25-Jul-12	-	-
	Sample ID:	1230199-01	1230199-02	-	-
T	MDL/Units	Water	Water	-	-
Styrene	0.5 ug/L	<0.5	<0.5	-	-
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	-	-
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	-	-
Tetrachloroethylene	0.5 ug/L	<0.5	<0.5	-	-
Toluene	0.5 ug/L	<0.5	<0.5	-	-
1,2,4-Trichlorobenzene	0.5 ug/L	<0.5	<0.5	-	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	<0.5	-	-
1,1,2-Trichloroethane	0.5 ug/L	<0.5	<0.5	-	-
Trichloroethylene	0.5 ug/L	<0.5	<0.5	-	-
Trichlorofluoromethane	1.0 ug/L	<1.0	<1.0	-	-
1,3,5-Trimethylbenzene	0.5 ug/L	<0.5	<0.5	-	-
Vinyl chloride	0.5 ug/L	<0.5	<0.5	-	-
m,p-Xylenes	0.5 ug/L	<0.5	<0.5	-	-
o-Xylene	0.5 ug/L	<0.5	<0.5	-	-
Xylenes, total	0.5 ug/L	<0.5	<0.5	-	-
4-Bromofluorobenzene	Surrogate	110%	108%	-	-
Dibromofluoromethane	Surrogate	110%	109%	-	-
Toluene-d8	Surrogate	112%	106%	-	-
Hydrocarbons					
F1 PHCs (C6-C10)	25 ug/L	<25	<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	-	-	-





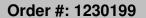
Client: Paterson Group Consulting Engineers

Client PO: 12263 Project Description: PE2709

Report Date: 26-Jul-2012 Order Date: 25-Jul-2012

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
	ND	25	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						
Chloroethane	ND	1.0	ug/L						
Chloroform	ND	0.5	ug/L						
Chloromethane	ND	3.0	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dibromoethane	ND	0.2	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 ND	0.5	ug/L						
1,1-Dichloroethylene	ND ND	0.5	ug/L						
cis-1,2-Dichloroethylene	ND	0.5	ug/L						
trans-1,2-Dichloroethylene 1,2-Dichloroethylene, total	ND	0.5 0.5	ug/L						
1,2-Dichloropropane	ND	0.5	ug/L ug/L						
cis-1,3-Dichloropropylene	ND	0.5	ug/L ug/L						
trans-1,3-Dichloropropylene	ND	0.5							
1,3-Dichloropropene, total	ND	0.5	ug/L ug/L						
Ethylbenzene	ND	0.5	ug/L ug/L						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Butyl Ketone (2-Hexanone)	ND	10.0	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						
1,2,4-Trichlorobenzene	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						
1,3,5-Trimethylbenzene	ND	0.5	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	37.1		ug/L		116	50-140			
Surrogate: Dibromofluoromethane	31.8		ug/L		99.5	50-140			
	38.5		ug/L		120	50-140			





Client: Paterson Group Consulting Engineers

Client PO: 12263 Project Description: PE2709 Report Date: 26-Jul-2012 Order Date:25-Jul-2012

Method Quality Control: Duplicate

Hydrocarbons Fi PHOs (C6-C10)			Reporting		Source		%REC		RPD	
Fi PHCs (Ge-C10)	Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
FI_PHCS (C6-C10)	Hydrocarbons									
Volatiles		ND	25	ua/L	ND				30	
Acetone Benzene ND D D D Seromodichloromethane ND D D S Bromodichloromethane ND D S S S S S S S S S S S S S S S S S				g/ =						
Benzene ND 0.5 ug/L ND 30		ND	- 0	"	ND				00	
Bromotichloromethane										
Bromoferm										
Brommethane										
Carbon Tetrachloride ND 0.2 ug/L ND 30 Chlorobenzene ND 0.5 ug/L ND 30 Chlorotemane ND 1.0 ug/L ND 30 Chlorotom ND 0.5 ug/L ND 30 Chloromethane ND 3.0 ug/L ND 30 Dibromochloromethane ND 0.5 ug/L ND 30 Dibromochloromethane ND 0.5 ug/L ND 30 1,2-Dibromoethane ND 0.5 ug/L ND 30 1,2-Dichlorobenzene ND 0.5 ug/L ND 30 1,3-Dichlorobenzene ND 0.5 ug/L ND 30 1,1-Dichlorobenzene ND 0.5 ug/L ND 30 1,2-Dichlorobethane ND 0.5 ug/L ND 30 1,2-Dichlorobethylene ND 0.5 ug/L ND 30										
Chlorobenzene										
Chloroethane										
Chloroform										
Chloromethane ND 3.0 ug/L ND 30 Dibromochloromethane ND 0.5 ug/L ND 30 Dichlorodifluoromethane ND 1.0 ug/L ND 30 1.2-Dibromoethane ND 0.2 ug/L ND 30 1.3-Dichlorobenzene ND 0.5 ug/L ND 30 1.3-Dichlorobenzene ND 0.5 ug/L ND 30 1.4-Dichlorobenzene ND 0.5 ug/L ND 30 1.1-Dichloroethane ND 0.5 ug/L ND 30 1.2-Dichloroethylene ND 0.5 ug/L ND 30 1.2-Dichloroethylene ND 0.5 ug/L ND 30 1.2-Dichloroethylene ND 0.5 ug/L ND 30 1.2-Dichloropropane ND 0.5 ug/L ND 30 1.2-Dichloropropylene ND 0.5 ug/L ND				ug/L						
Dibromochloromethane										
Dichlorodifluoromethane ND 1.0 ug/L ND 30 1,2-Dichlorobenzene ND 0.2 ug/L ND 30 1,2-Dichlorobenzene ND 0.5 ug/L ND 30 1,3-Dichlorobenzene ND 0.5 ug/L ND 30 1,4-Dichloroethane ND 0.5 ug/L ND 30 1,1-Dichloroethane ND 0.5 ug/L ND 30 1,1-Dichloroethylene ND 0.5 ug/L ND 30 1,1-Dichloroethylene ND 0.5 ug/L ND 30 1,1-Dichloroethylene ND 0.5 ug/L ND 30 1,2-Dichloroptylene ND 0.5 ug/L ND 30 1,2-Dichloropropale ND 0.5 ug/L ND 30 1,2-Dichloropropylene ND 0.5 ug/L ND 30 1,2-Dichloropropylene ND 0.5 ug/L ND <td></td>										
1,2-Dichlorobenzene										
1,2-Dichlorobenzene										
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			0.5			100	50-140		30	
	0									
Surrogate: Dibromofluoromethane 35.1 ug/L ND 110 50-140	· ·									
Surrogate: Toluene-d8 37.2 ug/L ND 116 50-140	Surroyale. Toluelle-uo	3/.2		ug/L	ND	110	30-140			



Client: Paterson Group Consulting Engineers

Client PO: 12263 Project Description: PE2709

Report Date: 26-Jul-2012 Order Date: 25-Jul-2012

Method Quality Control: Spike										
Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes	
Hydrocarbons										
F1 PHCs (C6-C10)	1910	25	ug/L	ND	95.4	68-117				
Volatiles										
Acetone	87.7	5.0	ug/L	ND	87.7	50-140				
Benzene	34.4	0.5	ug/L	ND	86.1	50-140				
Bromodichloromethane	31.0	0.5	ug/L	ND	77.6	50-140				
Bromoform	35.8	0.5	ug/L	ND	89.6	50-140				
Bromomethane	26.3	0.5	ug/L	ND	65.7	50-140				
Carbon Tetrachloride	30.3	0.2	ug/L	ND	75.8	50-140				
Chlorobenzene	42.4	0.5	ug/L	ND	106	50-140				
Chloroethane	37.6	1.0	ug/L	ND	93.9	50-140				
Chloroform	32.3	0.5	ug/L	ND	80.8	50-140				
Chloromethane	30.1	3.0	ug/L	ND	75.2	50-140				
Dibromochloromethane	39.0	0.5	ug/L	ND	97.6	50-140				
Dichlorodifluoromethane	27.9	1.0	ug/L	ND	69.8	50-140				
1,2-Dibromoethane	39.2	0.2	ug/L	ND	98.0	50-140				
1,2-Dichlorobenzene	32.3	0.5	ug/L	ND	80.8	50-140				
1,3-Dichlorobenzene	32.3	0.5	ug/L	ND	80.8	50-140				
1,4-Dichlorobenzene	33.0	0.5	ug/L	ND	82.4	50-140				
1,1-Dichloroethane	43.9	0.5	ug/L	ND	110	50-140				
1,2-Dichloroethane	29.7	0.5	ug/L	ND	74.3	50-140				
1,1-Dichloroethylene	36.2	0.5	ug/L	ND	90.4	50-140				
cis-1,2-Dichloroethylene	29.6	0.5	ug/L	ND	74.1	50-140				
trans-1,2-Dichloroethylene	47.4	0.5	ug/L	ND	119	50-140				
1,2-Dichloropropane	34.1	0.5	ug/L	ND	85.2	50-140				
cis-1,3-Dichloropropylene	49.7	0.5	ug/L	ND	124	50-140				
trans-1,3-Dichloropropylene	46.9	0.5	ug/L	ND	117	50-140				
Ethylbenzene	33.5	0.5	ug/L	ND	83.7	50-140				
Hexane	30.0	1.0	ug/L	ND	75.0	50-140				
Methyl Ethyl Ketone (2-Butanone)	104	5.0	ug/L	ND	104	50-140				
Methyl Butyl Ketone (2-Hexanone)	76.8	10.0	ug/L	ND	76.8	50-140				
Methyl Isobutyl Ketone	74.5	5.0	ug/L	ND	74.5	50-140				
Methyl tert-butyl ether	99.8	2.0	ug/L	ND	99.8	50-140				
Methylene Chloride	37.6	5.0	ug/L	ND	94.1	50-140				
Styrene	29.1	0.5	ug/L	ND	72.7	50-140				
1,1,1,2-Tetrachloroethane	40.3	0.5	ug/L	ND	101	50-140				
1,1,2,2-Tetrachloroethane	51.5	0.5	ug/L	ND	129	50-140				
Tetrachloroethylene	39.5	0.5	ug/L	ND	98.7	50-140				
Toluene	30.0	0.5	ug/L	ND	75.0	50-140				
1,2,4-Trichlorobenzene	27.0	0.5	ug/L	ND	67.6	50-140 50-140				
1,1,1-Trichloroethane 1,1,2-Trichloroethane	30.1	0.5	ug/L	ND	75.2					
* *	30.5	0.5 0.5	ug/L	ND	76.2	50-140				
Trichloroethylene Trichlorofluoromethane	28.5 32.2	0.5 1.0	ug/L ug/L	ND ND	71.2 80.6	50-140 50-140				
1,3,5-Trimethylbenzene	32.2 26.0	0.5	ug/L ug/L	ND	65.0	50-140 50-140				
Vinyl chloride	29.3	0.5	ug/L ug/L	ND	73.2	50-140				
m,p-Xylenes	68.0	0.5	ug/L ug/L	ND	85.0	50-140				
o-Xylene	36.6	0.5	ug/L ug/L	ND	91.5	50-140				
O Aylolio	30.0	0.5	ug/L	שויו	31.3	30-140				

25.8

ug/L

OTTAWA

50-140

80.8

Surrogate: 4-Bromofluorobenzene



PARACEL Order #: 1230199

Certificate of AnalysisReport Date: 26-Jul-2012Client: Paterson Group Consulting EngineersOrder Date: 25-Jul-2012

Client PO: 12263 Project Description: PE2709

Qualifier Notes:

None

Sample Data Revisions

None

Work Order Revisions / Comments:

None

Other Report Notes:

n/a: not applicable ND: Not Detected

MDL: Method Detection Limit

Source Result: Data used as source for matrix and duplicate samples

%REC: Percent recovery. RPD: Relative percent difference.

CCME PHC additional information:

- The method for the analysis of PHCs complies with the Reference Method for the CWS PHC and is validated for use in the laboratory. All prescribed quality criteria identified in the method has been met.
- F1 range corrected for BTEX.
- F2 to F3 ranges corrected for appropriate PAHs where available.
- The gravimetric heavy hydrocarbons (F4G) are not to be added to C6 to C50 hydrocarbons.
- In the case where F4 and F4G are both reported, the greater of the two results is to be used for comparison to CWS PHC criteria.

PARACEL LABORATORIES LTD.		RUST ESPO ELIAE	NSIV	/E .				(Ottaw p: 1-8	319 /a, C	St. Ontai 749-	io K1 1947	nt Blvd. G 4J8 labs.com	Chain of Custody (Lab Use Only) No. 94819				
OTTAWA @ KINGSTON @ NIAGARA @ MISSIS	SAUGA	SAF	RNIA					١	www.	para	acella	abs.co	m		Pa	ge o	f <u>/</u>	
Client Name: PATELSON GLOUP INC. Contact Name: MARK D'ARLY Address: 154 COLONNADE FORD SOUTH Telephone: (613) 226-7381					1263 10 PATER	801				******				Date I	Regu 2 Da Required:	y 1	1 3 Day	
Criteria: [] O. Reg. 153/04 Table [XO. Reg. 153/11 (Curren	nt) Table <u>9</u>	RSC	Filing	O. Reg. 558/00) PWQO	I CCM	E] SU	B (Sto	orm)	[]					I] Other; _	
Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water)	SS (Storm/S	lanitary S	ewer) P	(Paint) A (Air) O	(Other)							I	Required .	Analyse	es			
Paracel Order Number:	rix	Air Volume	Containers	Sampl	e Taken	s FI-F4+BTEX	S	s	ils by ICP/MS			WS)						
Sample ID/Location Name	Matrix	Air	Jo#	Date	Time	PHCs	VOC	PAHs	Metals	Hg	CrVI	B (HWS)						
1 BHH-GWZ	GW			July 25	12:30 pm	X	X											
2 BY2-GW2	GW			July 25	10 am	X	×											
3				V	20.0							1						
4														_			-	
5												4						
6												1						
7												4		_		_	_	
8								_/				4	_	_		-	_	
9						-						_			-	-	_	
10															Mark	d of Deli		
Comments: PHG FI@ BH2, GW	2,														Metho	Ma	JK7	n
Relinquished By (Print & Sign):	Receive	d by Dri	ver/Depo	ot:	Receiv	ed at L	.ab:			L			Veri	ied By: M/C		-;1		
	Date/Ti				Date/T	The state of the s	Ty		25/	دارا		5:4		Time:	Jul	y als	5/12	6:0
Date/Time:	Temper	ature:		C.	Tempe	rature:	17	4	V.C.				pH \	erified [] By:	1	V/A	



Head Office

300-2319 St. Laurent Blvd. Ottawa, Ontario K1G 4J8

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Certificate of Analysis

Paterson Group Consulting Engineers

154 Colonnade Road South Phone: (613) 226-7381 Nepean, ON K2E 7J5 Fax: (613) 226-6344

Attn: Karyn Munch

 Client PO: 15215
 Report Date: 18-Nov-2013

 Project: PE2709
 Order Date: 13-Nov-2013

 Custody: 99019
 Order #: 1346233

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

Paracel ID	Client ID
1346233-01	BH1-GW2
1346233-02	BH2-GW3
1346233-03	BH4-GW3

Approved By:

Mark Foto

Mark Foto, M.Sc. For Dale Robertson, BSc

Laboratory Director



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 15215 Project Description: PE2709 Report Date: 18-Nov-2013

Order Date:13-Nov-2013

Analysis Summary Table

Method Reference/Description Extraction Date Analysis Date Analysis VOCs by P&T GC-MS EPA 624 - P&T GC-MS 15-Nov-13 15-Nov-13

> P: 1-800-749-1947 E: PARACEL@PARACELLABS.COM

NIAGARA FALLS



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 15215 Project Description: PE2709 Report Date: 18-Nov-2013 Order Date:13-Nov-2013

	Client ID: Sample Date: Sample ID:	BH1-GW2 12-Nov-13 1346233-01	BH2-GW3 12-Nov-13 1346233-02	BH4-GW3 12-Nov-13 1346233-03	- - -
	MDL/Units	Water	Water	Water	-
Volatiles					
Acetone	5.0 ug/L	<5.0	15.3	117	-
Benzene	0.5 ug/L	<0.5	11.7	8.0	-
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	-
Chloroethane	1.0 ug/L	<1.0	<1.0	<1.0	-
Chloroform	0.5 ug/L	<0.5	<0.5	<0.5	-
Chloromethane	3.0 ug/L	<3.0	<3.0	<3.0	-
Dibromochloromethane	0.5 ug/L	<0.5	<0.5	<0.5	-
Dichlorodifluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	-
1,2-Dibromoethane	0.2 ug/L	<0.2	<0.2	<0.2	-
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-Dichloroethylene, total	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	5.9	1.2	-
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 Butyl Ketone (2-Hexanone	10.0 ug/L	<10.0	<10.0	<10.0	-
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	-



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 15215 Project Description: PE2709 Report Date: 18-Nov-2013 Order Date:13-Nov-2013

	Client ID: Sample Date: Sample ID:	BH1-GW2 12-Nov-13 1346233-01	BH2-GW3 12-Nov-13 1346233-02	BH4-GW3 12-Nov-13 1346233-03	- - -
	MDL/Units	Water	Water	Water	-
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	1.4	-
1,2,4-Trichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	-
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	-
1,3,5-Trimethylbenzene	0.5 ug/L	<0.5	1.2	<0.5	-
Vinyl chloride	0.5 ug/L	<0.5	<0.5	<0.5	-
m,p-Xylenes	0.5 ug/L	<0.5	<0.5	1.6	-
o-Xylene	0.5 ug/L	<0.5	<0.5	1.1	-
Xylenes, total	0.5 ug/L	<0.5	<0.5	2.8	-
4-Bromofluorobenzene	Surrogate	103%	98.5%	99.7%	-
Dibromofluoromethane	Surrogate	116%	118%	118%	-
Toluene-d8	Surrogate	98.9%	98.5%	97.9%	-



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 15215 Project Description: PE2709 Report Date: 18-Nov-2013 Order Date:13-Nov-2013

Method Quality Control: Blank	7								
Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Volatiles Acetone	ND	5.0	ug/L						

Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Volatiles									
Acetone	ND	5.0	ug/L						
Benzene	ND	0.5	ug/L						
Bromodichloromethane	ND	0.5	ug/L						
Bromoform	ND	0.5	ug/L						
Bromomethane	ND	0.5	ug/L						
Carbon Tetrachloride	ND	0.2	ug/L						
Chlorobenzene	ND	0.5	ug/L						
Chloroethane	ND	1.0	ug/L						
Chloroform	ND	0.5	ug/L						
Chloromethane	ND	3.0	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dibromoethane	ND	0.2	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-Dichloroethylene, total	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						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Butyl Ketone (2-Hexanone)	ND	10.0	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						
1,2,4-Trichlorobenzene	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						
1,3,5-Trimethylbenzene	ND	0.5	ug/L						
Vinyl chloride	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
Xylenes, total	ND	0.5	ug/L						
Surrogate: 4-Bromofluorobenzene	83.4		ug/L		104	50-140			
Surrogate: Dibromofluoromethane	91.8		ug/L		115	50-140			
Surrogate: Toluene-d8	78.0		ug/L		97.5	50-140			
₹			-						



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 15215 Project Description: PE2709

Report Date: 18-Nov-2013 Order Date:13-Nov-2013

Method Quality Control: Duplicate

		Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Volatiles									
Acetone	370	5.0	ug/L	336			9.6	30	
Benzene	1.00	0.5	ug/L	ND			0.0	30	
Bromodichloromethane	ND	0.5	ug/L	ND				30	
Bromoform	ND	0.5	ug/L	ND				30	
Bromomethane	ND	0.5	ug/L	ND				30	
Carbon Tetrachloride	ND	0.2	ug/L	ND			0.0	30	
Chlorobenzene	ND	0.5	ug/L	ND			0.0	30	
Chloroethane	ND	1.0	ug/L	ND				30	
Chloroform	2.25	0.5	ug/L	2.27			0.9	30	
Chloromethane	ND	3.0	ug/L	ND				30	
Dibromochloromethane	ND	0.5	ug/L	ND				30	
Dichlorodifluoromethane	ND	1.0	ug/L	ND				30	
1,2-Dibromoethane	ND	0.2	ug/L	ND				30	
1.2-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND				30	
1,1-Dichloroethane	ND ND	0.5	ug/L	ND				30	
1,2-Dichloroethane	ND ND	0.5	ug/L	ND			0.0	30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND			0.0	30	
cis-1,2-Dichloroethylene	ND ND	0.5	ug/L	ND				30	
trans-1,2-Dichloroethylene	ND ND	0.5	ug/L	ND				30	
1,2-Dichloropropane	ND	0.5	ug/L	ND				30	
cis-1,3-Dichloropropylene	ND ND	0.5	ug/L	ND				30	
rans-1,3-Dichloropropylene	ND	0.5	ug/L	ND				30	
Ethylbenzene	1.06	0.5	ug/L	1.01			4.8	30	
Hexane	25.5	1.0	ug/L	22.0			14.4	30	
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L	ND			17.7	30	
Methyl Butyl Ketone (2-Hexanone)	ND	10.0	ug/L	ND				30	
Methyl Isobutyl Ketone	ND ND	5.0	ug/L ug/L	ND				30	
Methyl tert-butyl ether	ND ND	2.0	ug/L ug/L	ND				30	
Methylene Chloride	ND ND	5.0	ug/L ug/L	ND				30	
Styrene	ND ND	0.5	ug/L ug/L	ND			0.0	30	
1,1,1,2-Tetrachloroethane	ND ND	0.5	ug/L ug/L	ND			0.0	30	
1,1,2,2-Tetrachloroethane	ND ND	0.5	ug/L ug/L	ND				30	
Tetrachloroethylene	ND ND	0.5	ug/L ug/L	ND			0.0	30	
Toluene	254	0.5	ug/L ug/L	253			0.0	30	
1,2,4-Trichlorobenzene	ND	0.5	ug/L ug/L	ND			0.0	30	
1,2,4-Trichlorobenzene 1,1,1-Trichloroethane	ND ND	0.5	ug/L ug/L	ND			0.0	30	
1,1,2-Trichloroethane	ND ND	0.5	ug/L ug/L	ND			0.0	30	
Trichloroethylene	ND ND	0.5	ug/L ug/L	ND				30	
Trichlorofluoromethane	ND ND	1.0	ug/L ug/L	ND			0.0	30	
1,3,5-Trimethylbenzene	ND ND	0.5	ug/L ug/L	ND ND			0.0	30	
Vinyl chloride	ND ND	0.5	ug/L ug/L	ND ND				30	
m,p-Xylenes	7.88	0.5	•	7.21			8.9	30	
o-Xylene	7.00 1.90	0.5	ug/L	1.83			3.8	30	
o-xylene Surrogate: 4-Bromofluorobenzene	1.90 <i>78.9</i>	0.5	ug/L	1.83 ND	98.6	50-140	3.0	30	
S .			ug/L						
Surrogate: Dibromofluoromethane	89.6		ug/L	ND	112	50-140			
Surrogate: Toluene-d8	76.9		ug/L	ND	96.1	50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 15215 Project Description: PE2709

Report Date: 18-Nov-2013 Order Date:13-Nov-2013

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Volatiles									
Acetone	112	5.0	ug/L	ND	112	50-140			
Benzene	43.0	0.5	ug/L	ND	107	60-130			
Bromodichloromethane	29.3	0.5	ug/L	ND	73.2	60-130			
Bromoform	24.4	0.5	ug/L	ND	60.9	60-130			
Bromomethane	37.0	0.5	ug/L	ND	92.6	50-140			
Carbon Tetrachloride	24.7	0.2	ug/L	ND	61.7	60-130			
Chlorobenzene	39.8	0.5	ug/L	ND	99.4	60-130			
Chloroethane	36.9	1.0	ug/L	ND	92.3	50-140			
Chloroform	33.9	0.5	ug/L	ND	84.8	60-130			
Chloromethane	32.9	3.0	ug/L	ND	82.3	50-140			
Dibromochloromethane	25.4	0.5	ug/L	ND	63.4	60-130			
Dichlorodifluoromethane	37.1	1.0	ug/L	ND	92.6	50-140			
1,2-Dibromoethane	38.2	0.2	ug/L	ND	95.4	60-130			
1,2-Dichlorobenzene	34.5	0.5	ug/L	ND	86.3	60-130			
1,3-Dichlorobenzene	34.8	0.5	ug/L	ND	87.0	60-130			
1,4-Dichlorobenzene	31.7	0.5	ug/L	ND	79.2	60-130			
1,1-Dichloroethane	41.2	0.5	ug/L	ND	103	60-130			
1,2-Dichloroethane	31.1	0.5	ug/L	ND	77.7	60-130			
1,1-Dichloroethylene	41.9	0.5	ug/L	ND	105	60-130			
cis-1,2-Dichloroethylene	45.4	0.5	ug/L	ND	114	60-130			
trans-1,2-Dichloroethylene	49.4	0.5	ug/L	ND	124	60-130			
1,2-Dichloropropane	43.4	0.5	ug/L	ND	109	60-130			
cis-1,3-Dichloropropylene	38.3	0.5	ug/L	ND	95.8	60-130			
trans-1,3-Dichloropropylene	34.0	0.5	ug/L	ND	85.0	60-130			
Ethylbenzene	32.4	0.5	ug/L	ND	80.9	60-130			
Hexane	47.7	1.0	ug/L	ND	119	60-130			
Methyl Ethyl Ketone (2-Butanone)	97.2	5.0	ug/L	ND	97.2	50-140			
Methyl Butyl Ketone (2-Hexanone)	81.8	10.0	ug/L	ND	81.8	50-140			
Methyl Isobutyl Ketone	78.1	5.0	ug/L	ND	78.1	50-140			
Methyl tert-butyl ether	83.6	2.0	ug/L	ND	83.6	50-140			
Methylene Chloride	38.7	5.0	ug/L	ND	96.6	60-130			
Styrene	37.7	0.5	ug/L	ND	94.2	60-130			
1,1,1,2-Tetrachloroethane	29.0	0.5	ug/L	ND	72.4	60-130			
1,1,2,2-Tetrachloroethane	40.8	0.5	ug/L	ND	102	60-130			
Tetrachloroethylene	30.4	0.5	ug/L	ND	75.9	60-130			
Toluene	31.6	0.5	ug/L	ND	79.0	60-130			
1,2,4-Trichlorobenzene	26.1	0.5	ug/L	ND	65.2	60-130			
1,1,1-Trichloroethane	24.3	0.5	ug/L	ND	60.8	60-130			
1,1,2-Trichloroethane	41.2	0.5	ug/L	ND	103	60-130			
Trichloroethylene	37.6	0.5	ug/L	ND	94.1	60-130			
Trichlorofluoromethane	27.0	1.0	ug/L	ND	67.5	60-130			
1,3,5-Trimethylbenzene	28.9	0.5	ug/L	ND	72.2	60-130			
Vinyl chloride	36.2	0.5	ug/L	ND	90.5	50-140			
m,p-Xylenes	77.4	0.5	ug/L	ND	96.8	60-130			
o-Xylene	33.0	0.5	ug/L	ND	82.4	60-130			
Surrogate: 4-Bromofluorobenzene	72.5		ug/L		90.6	50-140			



Order #: 1346233

Client: Paterson Group Consulting Engineers

Order Date:13-Nov-2013 Client PO: 15215 Project Description: PE2709

Qualifier Notes:

None

Sample Data Revisions

None

Work Order Revisions / Comments:

None

Other Report Notes:

n/a: not applicable ND: Not Detected

MDL: Method Detection Limit

Source Result: Data used as source for matrix and duplicate samples

%REC: Percent recovery.

RPD: Relative percent difference.

Report Date: 18-Nov-2013

Contact Name: Kary Munch Address: 154 Cotomode Rols. Telephone: 013-226-7381 Criteria: [] O. Reg. 153/04 (As Amended) Table _ [] RSC Filing [] O. Reg. 558/00 [] PWQO [] CCME [] SUB (Storm) [] SUB (Samitary) Municipality: [Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) SS (Storm/Sanitary Sewer) P (Paint) A (Air) O (Other) Required Analyses Paracel Order Number:	Regular [] 2 Day ste Required: [] Other:		3 Day	
Paracel Order Number: 344233 Sample Taken Sa				
1344233 Sample Taken 144-1 do 166			£ 1	5
Sample ID/Location Name \square \square \square \square \square \square Date Time \square		m 7 -		1.5 4
1 BH1-GW2 GW Z NOV.12/13 am 7	. 127 11			
2 BH2-6W3 3 BH4-6W3 GW V V			= A 7	
5				
6				
7				
10				3 7

Received at Lab:

Temperature: 17/8 C

2:36 PH Date/Time: NOV 13 3013

Received by Driver/Depot:

Temperature:

Verified By:

Date/Time:

pH Verified [] By:

Relinquished By (Sign):

Relinquished By (Print):



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

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive Ottawa, ON K2E 7T9 Attn: Karyn Munch

Client PO: 56980 Project: PE2709

Custody:

Report Date: 14-Mar-2023 Order Date: 9-Mar-2023

Order #: 2310387

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

Paracel ID	Client ID
2310387-01	BH2-23-GW1
2310387-02	BH3-23-GW1
2310387-03	BH5-23-GW1
2310387-04	BH6-23-GW1
2310387-05	DUP1-23-GW1

Approved By:

Mark Froto

Mark Foto, M.Sc. Lab Supervisor



Order #: 2310387

Report Date: 14-Mar-2023 Order Date: 9-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 9-Mar-2023

 Client PO:
 56980
 Project Description: PE2709

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
PHC F1	CWS Tier 1 - P&T GC-FID	10-Mar-23	10-Mar-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	10-Mar-23	13-Mar-23
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	10-Mar-23	10-Mar-23



Certificate of Analysis Client: Paterson Group Consulting Engineers

Project Description: PE2709

Report Date: 14-Mar-2023

Order Date: 9-Mar-2023

Client PO: 56980

ſ	Client ID: Sample Date: Sample ID: MDL/Units	BH2-23-GW1 08-Mar-23 00:00 2310387-01 Ground Water	BH3-23-GW1 08-Mar-23 00:00 2310387-02 Ground Water	BH5-23-GW1 08-Mar-23 00:00 2310387-03 Ground Water	BH6-23-GW1 08-Mar-23 00:00 2310387-04 Ground Water
Volatiles			•	•	•
Acetone	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Benzene	0.5 ug/L	<0.5	0.6	<0.5	<0.5
Bromodichloromethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromoform	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Bromomethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Carbon Tetrachloride	0.2 ug/L	<0.2	<0.2	<0.2	<0.2
Chlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Chloroform	0.5 ug/L	<0.5	2.5	<0.5	1.0
Dibromochloromethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Dichlorodifluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
1,2-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	0.7	<0.5
1,3-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,4-Dichlorobenzene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,2-Dichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
cis-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
trans-1,2-Dichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,2-Dichloropropane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
cis-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
trans-1,3-Dichloropropylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,3-Dichloropropene, total	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Ethylbenzene	0.5 ug/L	<0.5	<0.5	17.0	<0.5
Ethylene dibromide (dibromoethane, 1,2-)	0.2 ug/L	<0.2	<0.2	<0.2	<0.2
Hexane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
Methyl Ethyl Ketone (2-Butanone)	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Methyl Isobutyl Ketone	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Methyl tert-butyl ether	2.0 ug/L	<2.0	<2.0	<2.0	<2.0
Methylene Chloride	5.0 ug/L	<5.0	<5.0	<5.0	<5.0
Styrene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,1,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
1,1,2,2-Tetrachloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Tetrachloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Toluene	0.5 ug/L	<0.5	3.4	1.0	<0.5
1,1,1-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5



Certificate of Analysis Client: Paterson Group Consulting Engineers

Order Date: 9-Mar-2023 **Project Description: PE2709**

Report Date: 14-Mar-2023

Client PO: 56980

	Client ID:	BH2-23-GW1	BH3-23-GW1	BH5-23-GW1	BH6-23-GW1
		08-Mar-23 00:00	08-Mar-23 00:00	08-Mar-23 00:00	08-Mar-23 00:00
	Sample Date:	2310387-01	2310387-02	2310387-03	2310387-04
	Sample ID:				
	MDL/Units	Ground Water	Ground Water	Ground Water	Ground Water
1,1,2-Trichloroethane	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Trichloroethylene	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
Trichlorofluoromethane	1.0 ug/L	<1.0	<1.0	<1.0	<1.0
Vinyl chloride	0.5 ug/L	<0.5	<0.5	<0.5	<0.5
m,p-Xylenes	0.5 ug/L	<0.5	3.2	22.0	<0.5
o-Xylene	0.5 ug/L	<0.5	1.0	2.0	<0.5
Xylenes, total	0.5 ug/L	<0.5	4.1	24.0	<0.5
4-Bromofluorobenzene	Surrogate	108%	109%	117%	108%
Dibromofluoromethane	Surrogate	105%	102%	115%	111%
Toluene-d8	Surrogate	110%	110%	108%	110%
Hydrocarbons					
F1 PHCs (C6-C10)	25 ug/L	<25	-	1210	<25
F2 PHCs (C10-C16)	100 ug/L	<100	-	266	<100
F3 PHCs (C16-C34)	100 ug/L	<100	-	<100	<100
F4 PHCs (C34-C50)	100 ug/L	<100	-	<100	<100
			•	•	



Client: Paterson Group Consulting Engineers

Certificate of Analysis

Order #: 2310387

Report Date: 14-Mar-2023

Order Date: 9-Mar-2023

Client PO: 56980 Project Description: PE2709

	Client ID: Sample Date:	DUP1-23-GW1 08-Mar-23 00:00	-	-	
_	Sample ID:	2310387-05	-	-	-
Valentile	MDL/Units	Ground Water	-	-	-
Volatiles	5.0 ug/L	<5.0			
Acetone	0.5 ug/L		-	-	-
Benzene	0.5 ug/L	0.6	-	-	-
Bromodichloromethane	0.5 ug/L	<0.5	-	-	-
Bromoform	0.5 ug/L	<0.5	-	-	-
Bromomethane		<0.5	-	-	-
Carbon Tetrachloride	0.2 ug/L	<0.2	-	-	-
Chlorobenzene	0.5 ug/L	<0.5	-	-	-
Chloroform	0.5 ug/L	2.6	-	-	-
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	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	-	-	-



Order #: 2310387

Report Date: 14-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 9-Mar-2023

 Client PO:
 56980
 Project Description: PE2709

	Client ID:	DUP1-23-GW1	-	_	_
	Sample Date:	08-Mar-23 00:00	_	_	_
	Sample ID:	2310387-05	-	-	-
	MDL/Units	Ground Water	-	-	-
1,1,1-Trichloroethane	0.5 ug/L	<0.5	-	-	-
1,1,2-Trichloroethane	0.5 ug/L	<0.5	-	-	-
Trichloroethylene	0.5 ug/L	<0.5	-	-	-
Trichlorofluoromethane	1.0 ug/L	<1.0	-	-	-
Vinyl chloride	0.5 ug/L	<0.5	-	-	-
m,p-Xylenes	0.5 ug/L	3.2	-	-	-
o-Xylene	0.5 ug/L	1.0	-	-	-
Xylenes, total	0.5 ug/L	4.2	-	-	-
4-Bromofluorobenzene	Surrogate	112%	-	-	-
Dibromofluoromethane	Surrogate	108%	-	-	-
Toluene-d8	Surrogate	109%	-	-	-



Order #: 2310387

Report Date: 14-Mar-2023

Order Date: 9-Mar-2023

Client: Paterson Group Consulting Engineers Client PO: 56980 **Project Description: PE2709**

Method Quality Control: Blank

.		Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
F2 PHCs (C10-C16)	ND	100	ug/L						
F3 PHCs (C16-C34)	ND	100	ug/L						
F4 PHCs (C34-C50)	ND	100	ug/L						
Volatiles									
Acetone	ND	5.0	ug/L						
Benzene	ND	0.5	ug/L						
Bromodichloromethane	ND	0.5	ug/L						
Bromoform	ND	0.5	ug/L						
Bromomethane	ND	0.5	ug/L						
Carbon Tetrachloride	ND	0.2	ug/L						
Chlorobenzene	ND	0.5	ug/L						
Chloroform	ND	0.5	ug/L						
Dibromochloromethane	ND	0.5	ug/L						
Dichlorodifluoromethane	ND	1.0	ug/L						
1,2-Dichlorobenzene	ND	0.5	ug/L						
1,3-Dichlorobenzene	ND	0.5	ug/L						
1,4-Dichlorobenzene	ND	0.5	ug/L						
1,1-Dichloroethane	ND	0.5	ug/L						
1,2-Dichloroethane	ND	0.5	ug/L						
1,1-Dichloroethylene	ND	0.5	ug/L						
cis-1,2-Dichloroethylene	ND	0.5	ug/L						
trans-1,2-Dichloroethylene	ND	0.5	ug/L						
1,2-Dichloropropane	ND	0.5	ug/L						
cis-1,3-Dichloropropylene	ND	0.5	ug/L						
trans-1,3-Dichloropropylene	ND	0.5	ug/L						
1,3-Dichloropropene, total	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Ethylene dibromide (dibromoethane, 1,2-	ND	0.2	ug/L						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Isobutyl Ketone	ND	5.0	ug/L						
Methyl tert-butyl ether	ND	2.0	ug/L						
Methylene Chloride	ND	5.0	ug/L						
Styrene	ND	0.5	ug/L						
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L						
Tetrachloroethylene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
1,1,1-Trichloroethane	ND	0.5	ug/L						
1,1,2-Trichloroethane	ND	0.5	ug/L						
Trichloroethylene	ND	0.5	ug/L						
Trichlorofluoromethane	ND	1.0	ug/L						
Vinyl chloride	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
Xylenes, total	ND	0.5	ug/L						
Surrogate: 4-Bromofluorobenzene	87.8		ug/L		110	50-140			
Surrogate: Dibromofluoromethane	82.0		ug/L		102	50-140			
<u> </u>	88.8		ug/L		111	50-140			



Order #: 2310387

Report Date: 14-Mar-2023 Order Date: 9-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 9-Mar-2023

 Client PO:
 56980
 Project Description: PE2709

Method Quality Control: Duplicate

Availab		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	
/olatiles			-						
Acetone	ND	5.0	ug/L	ND			NC	30	
Benzene	ND ND	0.5	ug/L ug/L	ND ND			NC NC	30	
Bromodichloromethane	ND ND	0.5 0.5	ug/L ug/L	ND ND			NC NC	30 30	
Bromodicnioromethane Bromoform	ND ND	0.5 0.5	ug/L ug/L	ND ND			NC NC	30 30	
Bromomethane	ND ND	0.5 0.5	ug/L ug/L	ND ND			NC NC	30	
Carbon Tetrachloride	ND ND	0.5 0.2	ug/L ug/L	ND ND			NC NC	30 30	
Carbon Tetrachioride Chlorobenzene	ND ND		-	ND ND			NC NC	30 30	
		0.5	ug/L						
Chloroform	ND ND	0.5	ug/L	ND ND			NC NC	30 30	
Dibromochloromethane	ND ND	0.5	ug/L	ND ND			NC NC	30 30	
Dichlorodifluoromethane	ND ND	1.0	ug/L	ND			NC NC	30	
1,2-Dichlorobenzene	ND ND	0.5	ug/L	ND			NC	30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,1-Dichloroethane	ND	0.5	ug/L	ND			NC	30	
1,2-Dichloroethane	ND	0.5	ug/L	ND			NC	30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND			NC	30	
cis-1,2-Dichloroethylene	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 ND	0.5	ug/L ug/L	ND			NC	30	
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L	ND			NC	30	
Tetrachloroethylene	ND ND	0.5	ug/L ug/L	ND			NC	30	
Toluene	ND ND	0.5	ug/L ug/L	ND ND			NC NC	30	
1,1,1-Trichloroethane	ND ND	0.5	ug/L ug/L	ND			NC	30	
1,1,2-Trichloroethane	ND ND	0.5	ug/L ug/L	ND ND			NC NC	30	
1,1,2-11ichloroethane Trichloroethylene	ND ND	0.5 0.5	ug/L ug/L	ND ND			NC NC	30	
Trichloroethylene Trichlorofluoromethane	ND ND	0.5 1.0	ug/L ug/L	ND ND			NC NC	30 30	
	ND ND	1.0 0.5	-	ND ND			NC NC	30 30	
Vinyl chloride	ND ND	0.5 0.5	ug/L	ND ND			NC NC	30 30	
m,p-Xylenes			ug/L						
o-Xylene	ND	0.5	ug/L	ND	400	E0 115	NC	30	
Surrogate: 4-Bromofluorobenzene	87.6		ug/L		109	50-140			
Surrogate: Dibromofluoromethane	85.5		ug/L		107	50-140			
Surrogate: Toluene-d8	87.9		ug/L		110	50-140			



Report Date: 14-Mar-2023 Order Date: 9-Mar-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 56980

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
lydrocarbons									
F1 PHCs (C6-C10)	1950	25	ug/L	ND	97.6	68-117			
F2 PHCs (C10-C16)	1780	100	ug/L	ND	111	60-140			
F3 PHCs (C16-C34)	4660	100	ug/L	ND	119	60-140			
F4 PHCs (C34-C50)	2500	100	ug/L	ND	101	60-140			
olatiles			Ü						
Acetone	103	5.0	ug/L	ND	103	50-140			
Benzene	44.3	0.5	ug/L	ND	111	60-130			
Bromodichloromethane	48.2	0.5	ug/L	ND	120	60-130			
Bromoform	44.0	0.5	ug/L	ND	110	60-130			
Bromomethane	33.6	0.5	ug/L	ND	84.0	50-140			
Carbon Tetrachloride	44.1	0.2	ug/L	ND	110	60-130			
Chlorobenzene	44.0	0.5	ug/L	ND	110	60-130			
Chloroform	33.5	0.5	ug/L	ND	83.7	60-130			
Dibromochloromethane	41.9	0.5	ug/L	ND	105	60-130			
Dichlorodifluoromethane	49.9	1.0	ug/L	ND	125	50-140			
,2-Dichlorobenzene	40.8	0.5	ug/L	ND	102	60-130			
,3-Dichlorobenzene	39.6	0.5	ug/L	ND	99.0	60-130			
,4-Dichlorobenzene	38.1	0.5	ug/L	ND	95.3	60-130			
,1-Dichloroethane	41.9	0.5	ug/L	ND	105	60-130			
,2-Dichloroethane	44.8	0.5	ug/L	ND	112	60-130			
,1-Dichloroethylene	46.7	0.5	ug/L	ND	117	60-130			
sis-1,2-Dichloroethylene	41.4	0.5	ug/L	ND	104	60-130			
rans-1,2-Dichloroethylene	39.0	0.5	ug/L	ND	97.4	60-130			
,2-Dichloropropane	50.2	0.5	ug/L	ND	126	60-130			
sis-1,3-Dichloropropylene	47.6	0.5	ug/L	ND	119	60-130			
rans-1,3-Dichloropropylene	41.8	0.5	ug/L	ND	104	60-130			
Ethylbenzene	47.0	0.5	ug/L	ND	117	60-130			
Ethylene dibromide (dibromoethane, 1,2	47.4	0.2	ug/L	ND	119	60-130			
Hexane	41.0	1.0	ug/L	ND	102	60-130			
Methyl Ethyl Ketone (2-Butanone)	128	5.0	ug/L	ND	128	50-140			
Methyl Isobutyl Ketone	139	5.0	ug/L	ND	139	50-140			
Methyl tert-butyl ether	107	2.0	ug/L	ND	107	50-140			
Methylene Chloride	39.6	5.0	ug/L	ND	99.0	60-130			
Styrene	40.4	0.5	ug/L	ND	101	60-130			
, ,1,1,2-Tetrachloroethane	43.4	0.5	ug/L	ND	108	60-130			
I,1,2,2-Tetrachloroethane	43.3	0.5	ug/L	ND	108	60-130			
Tetrachloroethylene	39.4	0.5	ug/L	ND	98.4	60-130			
Foluene	47.1	0.5	ug/L	ND	118	60-130			
,1,1-Trichloroethane	46.4	0.5	ug/L	ND	116	60-130			
,1,2-Trichloroethane	49.0	0.5	ug/L	ND	122	60-130			
richloroethylene	43.8	0.5	ug/L	ND	110	60-130			
Trichlorofluoromethane	49.8	1.0	ug/L	ND	125	60-130			
/inyl chloride	34.1	0.5	ug/L	ND	85.2	50-140			
n,p-Xylenes	87.7	0.5	ug/L	ND	110	60-130			
-Xylene	44.9	0.5	ug/L	ND	112	60-130			
Surrogate: 4-Bromofluorobenzene	85.5		ug/L		107	50-140			
Surrogate: 4 Diomondologicale Surrogate: Dibromofluoromethane	67.1		ug/L		83.9	50-140			
Surrogate: Toluene-d8	84.3		ug/L		105	50-140			



Client: Paterson Group Consulting Engineers

Order #: 2310387

Report Date: 14-Mar-2023 Order Date: 9-Mar-2023

Client PO: 56980 Project Description: PE2709

Qualifier Notes:

Sample Data Revisions

Certificate of Analysis

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.





Chain Of Custody Paracel Order Number (Lab Use Only) (Lab Use Only)

Client Name: Paterson		Deale		A	, J										
Contact Name: Karyn Munch	-	-	ect Ref:	PE 2709								P	age]	of L	
7991635.		Quot										Turn	naround	Time	
9 Auriga Drive				980							□ 1	day		□ 3	3 day
Telephone: 613 226 7381		E-mai	li: K	Munch @ Pate	rson group.co	1					□ 2	day		∀ F	Regular
Per man			\$	Berube@ pater	son group, co	١					ate Re	equired:			
Table 1 Control Control Control Regulation		Matrix '	Type:	S (Soil/Sed.) GW (Gr	mund Water			12 13	100	10.16		40.7	47.70		
Table 1 Res/Park Med/Fine REG 558 PWQo		SW (Su	ırface \	Water) SS (Storm/Sar	nitary Sewer)					Requ	ired A	nalysis			
Table 2 Ind/Comm Coarse CCME MISA Table 3 Agri/Other SU-Sani SU-Sani			P (F	Paint) A (Air) O (Oth	er)	X				T		100 400	Т	3445.400	To all the
Table 20 - Storm			ers			F1-F4+BTEX			۵						
See Boo Flore Fr		a B	Containers	Sample	Taken	1-F4			Ş						
For RSC: U Yes No Other: Sample ID/Location Name	Matrix	Aīr Volume	of Cor				S	\$	Metals by ICP		_	(HWS)			
1 BH2-23- GW1	+	Ą	#	Date	Time	PHCs	VOCs	PAHs	Met	Ħ.		B (T			
2 BH 3-23 - GW1	GW		3	March 8, 2023		X	X								T
3 BH5-23 - GWI	\sqcup	_	2				X								T
4 BH6 -23 - GUI	Ш		3			X	X				\top				T
	¥		3	+		х	X					\top	\forall	\top	\dagger
5 DUP1-23 - GW1	V		2	•			X			\top	\top	+	+	+	+
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nquished by (Print): Bryce Lee Date/Time:	0	/-		TLOUSE I	Date/Time: 4	/		1.5	10.00	(W) ~		1	_
March 9 2023 Temperature:	4	03	12	3 /3/9	Pate/Time: Mo emperature:				_	ate/Tim	1	100	100	125	819
of Custody (Blank),xlsx	7.66	ARR		Restrict 4.0	emperature;	- 6	2.8		P	H Verific	ed: 🗆	θу	VA		



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

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive Ottawa, ON K2E 7T9 Attn: Sam Berube

Client PO: 57092 Project: PE2709

Custody:

Report Date: 29-Mar-2023 Order Date: 24-Mar-2023

Order #: 2312554

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

 Paracel ID
 Client ID

 2312554-01
 BH3-23-GW2

 2312554-02
 BH5-23-GW2

Approved By:

Mark Froto

Mark Foto, M.Sc. Lab Supervisor



Order #: 2312554

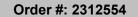
Report Date: 29-Mar-2023 Order Date: 24-Mar-2023

Project Description: PE2709

Client: Paterson Group Consulting Engineers
Client PO: 57092

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 624 - P&T GC-MS	28-Mar-23	28-Mar-23
PHC F1	CWS Tier 1 - P&T GC-FID	27-Mar-23	28-Mar-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	27-Mar-23	28-Mar-23
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	27-Mar-23	28-Mar-23





Client: Paterson Group Consulting Engineers

Client PO: 57092

Report Date: 29-Mar-2023 Order Date: 24-Mar-2023

Project Description: PE2709

Γ	Client ID: Sample Date: Sample ID: MDL/Units	BH3-23-GW2 23-Mar-23 09:00 2312554-01 Ground Water	BH5-23-GW2 23-Mar-23 09:00 2312554-02 Ground 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: 29-Mar-2023

Order Date: 24-Mar-2023

Certificate of Analysis Client: Paterson Group Consulting Engineers

Client PO: 57092 **Project Description: PE2709**

	Client ID:	BH3-23-GW2	BH5-23-GW2	-	-
	Sample Date:	23-Mar-23 09:00	23-Mar-23 09:00	-	-
	Sample ID:	2312554-01	2312554-02	-	-
1	MDL/Units	Ground Water	Ground 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	131%	-	-	-
Dibromofluoromethane	Surrogate	96.5%	-	-	-
Toluene-d8	Surrogate	121%	-	-	-
Benzene	0.5 ug/L	-	<0.5	-	-
Ethylbenzene	0.5 ug/L	-	9.2	-	-
Toluene	0.5 ug/L	-	2.4	-	-
m,p-Xylenes	0.5 ug/L	-	11.5	-	-
o-Xylene	0.5 ug/L	-	1.4	-	-
Xylenes, total	0.5 ug/L	-	12.9	-	-
Toluene-d8	Surrogate	-	115%	-	-
Hydrocarbons	,				
F1 PHCs (C6-C10)	25 ug/L	<25	742	-	-
F2 PHCs (C10-C16)	100 ug/L	<100	<100	-	-
F3 PHCs (C16-C34)	100 ug/L	<100	<100	-	-
F4 PHCs (C34-C50)	100 ug/L	<100	<100	-	-
	•				



Report Date: 29-Mar-2023

Order Date: 24-Mar-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers
Client PO: 57092

Method Quality Control: Blank

Analyte	D	Reporting		Source		%REC		RPD	
Analyte	Result	Limit	Units	Result	%REC	Limit	RPD	Limit	Notes
lydrocarbons									
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						
/olatiles			· ·						
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 ND	1.0	ug/L ug/L						
1,2-Dichlorobenzene	ND ND	0.5	ug/L ug/L						
1,3-Dichlorobenzene	ND ND	0.5	ug/L ug/L						
1,4-Dichlorobenzene	ND ND	0.5	ug/L ug/L						
,			-						
1,1-Dichloroethane	ND	0.5	ug/L						
1,2-Dichloroethane	ND	0.5	ug/L						
1,1-Dichloroethylene	ND	0.5	ug/L						
cis-1,2-Dichloroethylene	ND	0.5	ug/L						
trans-1,2-Dichloroethylene	ND	0.5	ug/L						
1,2-Dichloropropane	ND	0.5	ug/L						
cis-1,3-Dichloropropylene	ND	0.5	ug/L						
trans-1,3-Dichloropropylene	ND	0.5	ug/L						
1,3-Dichloropropene, total	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Ethylene dibromide (dibromoethane, 1,2	ND	0.2	ug/L						
Hexane	ND	1.0	ug/L						
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L						
Methyl Isobutyl Ketone	ND	5.0	ug/L						
Methyl tert-butyl ether	ND	2.0	ug/L						
Methylene Chloride	ND	5.0	ug/L						
Styrene	ND	0.5	ug/L						
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L						
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L						
Tetrachloroethylene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L						
1,1,1-Trichloroethane	ND	0.5	ug/L						
1,1,2-Trichloroethane	ND	0.5	ug/L						
Trichloroethylene	ND	0.5	ug/L						
Trichlorofluoromethane	ND	1.0	ug/L						
Vinyl chloride	ND	0.5	ug/L						
m,p-Xylenes	ND	0.5	ug/L						
o-Xylene	ND	0.5	ug/L						
Xylenes, total	ND	0.5	ug/L						
Surrogate: 4-Bromofluorobenzene	106		ug/L		133	50-140			
Surrogate: Dibromofluoromethane	76.8		ug/L		96.0	50-140			
Surrogate: Toluene-d8	100		ug/L		125	50-140			
Benzene	ND	0.5	ug/L						
Ethylbenzene	ND	0.5	ug/L						
Toluene	ND	0.5	ug/L ug/L						
m,p-Xylenes	ND	0.5	ug/L ug/L						
o-Xylene	ND ND	0.5	ug/L ug/L						
Xylenes, total	ND ND	0.5	-						
Surrogate: Toluene-d8	100	0.5	ug/L		125	50-140			



Order #: 2312554

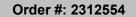
Report Date: 29-Mar-2023

Order Date: 24-Mar-2023

Client: Paterson Group Consulting Engineers Client PO: 57092 **Project Description: PE2709**

Method Quality Control: Duplicate

		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	
/olatiles			o o						
Acetone	ND	5.0	ug/l	ND			NC	30	
Benzene	ND ND	0.5	ug/L ug/L	ND			NC NC	30	
Bromodichloromethane	ND ND	0.5	-	ND ND			NC NC	30	
Bromoform	ND ND	0.5	ug/L	ND			NC	30	
Bromomethane	ND ND	0.5	ug/L ug/L	ND			NC	30	
Carbon Tetrachloride	ND ND	0.2	ug/L ug/L	ND			NC	30	
Chlorobenzene	ND	0.5	ug/L	ND			NC	30	
Chloroform	ND ND	0.5	ug/L ug/L	ND			NC	30	
Dibromochloromethane	ND ND	0.5	ug/L ug/L	ND			NC	30	
Dichlorodifluoromethane	ND ND	1.0	ug/L ug/L	ND			NC	30	
1,2-Dichlorobenzene	ND	0.5	ug/L ug/L	ND			NC	30	
1,3-Dichlorobenzene	ND ND	0.5	ug/L ug/L	ND			NC	30	
1,4-Dichlorobenzene	ND	0.5	ug/L ug/L	ND			NC	30	
1,1-Dichloroethane	ND ND	0.5	ug/L 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	99.7		ug/L		125	50-140			
Surrogate: Dibromofluoromethane	77.2		ug/L		96.6	50-140			
Surrogate: Toluene-d8	100		ug/L		126	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: Toluene-d8	100	-	ug/L		126	50-140	-	-	





Client: Paterson Group Consulting Engineers

Client PO: 57092 Project Description: PE2709

Report Date: 29-Mar-2023 Order Date: 24-Mar-2023

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1950	25	ug/L	ND	97.4	68-117			
F2 PHCs (C10-C16)	1870	100	ug/L	ND	117	60-140			
F3 PHCs (C16-C34)	4600	100	ug/L	ND	117	60-140			
F4 PHCs (C34-C50)	2530	100	ug/L	ND	102	60-140			
Volatiles									
Acetone	120	5.0	ug/L	ND	120	50-140			
Benzene	33.2	0.5	ug/L	ND	83.0	60-130			
Bromodichloromethane	34.5	0.5	ug/L	ND	86.2	60-130			
Bromoform	41.8	0.5	ug/L	ND	104	60-130			
Bromomethane	36.1	0.5	ug/L	ND	90.2	50-140			
Carbon Tetrachloride	35.9	0.2	ug/L	ND	89.7	60-130			
Chlorobenzene	38.6	0.5	ug/L	ND	96.6	60-130			
Chloroform	36.5	0.5	ug/L	ND	91.2	60-130			
Dibromochloromethane	42.2	0.5	ug/L	ND	105	60-130			
Dichlorodifluoromethane	40.9	1.0	ug/L	ND	102	50-140			
1,2-Dichlorobenzene	33.3	0.5	ug/L	ND	83.4	60-130			
1,3-Dichlorobenzene	36.1	0.5	ug/L	ND	90.3	60-130			
1,4-Dichlorobenzene	31.6	0.5	ug/L	ND	79.1	60-130			
1,1-Dichloroethane	38.0	0.5	ug/L	ND	95.1	60-130			
1,2-Dichloroethane	34.7	0.5	ug/L	ND	86.8	60-130			
1,1-Dichloroethylene	36.8	0.5	ug/L	ND	92.0	60-130			
cis-1,2-Dichloroethylene	35.1	0.5	ug/L	ND	87.8	60-130			
trans-1,2-Dichloroethylene	33.9	0.5	ug/L	ND	84.8	60-130			
1,2-Dichloropropane	30.3	0.5	ug/L	ND	75.8	60-130			
cis-1,3-Dichloropropylene	30.2	0.5	ug/L	ND	75.6	60-130			
trans-1,3-Dichloropropylene	32.1	0.5	ug/L	ND	80.3	60-130			
Ethylbenzene	36.7	0.5	ug/L	ND	91.7	60-130			
Ethylene dibromide (dibromoethane, 1,2-	38.9	0.2	ug/L	ND	97.3	60-130			
Hexane	35.4	1.0	ug/L	ND	88.4	60-130			
Methyl Ethyl Ketone (2-Butanone)	107	5.0	ug/L	ND	107	50-140			
Methyl Isobutyl Ketone	83.5	5.0	ug/L	ND	83.5	50-140			
Methyl tert-butyl ether	90.6	2.0	ug/L	ND	90.6	50-140			
Methylene Chloride	34.4	5.0	ug/L	ND	86.0	60-130			
Styrene	37.1	0.5	ug/L	ND	92.7	60-130			
1,1,1,2-Tetrachloroethane	40.6	0.5	ug/L	ND	102	60-130			
1,1,2,2-Tetrachloroethane	45.4	0.5	ug/L	ND	114	60-130			
Tetrachloroethylene	39.4	0.5	ug/L	ND	98.6	60-130			
Toluene	41.4	0.5	ug/L	ND	103	60-130			
1,1,1-Trichloroethane	34.1	0.5	ug/L	ND	85.3	60-130			
1,1,2-Trichloroethane	30.2	0.5	ug/L	ND	75.4	60-130			
Trichloroethylene	33.6	0.5	ug/L	ND	84.1	60-130			
Trichlorofluoromethane	37.9	1.0	ug/L	ND	94.7	60-130			
Vinyl chloride	42.6	0.5	ug/L	ND	107	50-140			
m,p-Xylenes	77.0	0.5	ug/L	ND	96.3	60-130			
o-Xylene	38.7	0.5	ug/L	ND	96.7	60-130			
Surrogate: 4-Bromofluorobenzene	77.8		ug/L		97.3	50-140			
Surrogate: Dibromofluoromethane	75.8		ug/L		94.8	50-140			
Surrogate: Toluene-d8	81.7		ug/L		102	50-140			



Order #: 2312554

Report Date: 29-Mar-2023

Order Date: 24-Mar-2023 **Project Description: PE2709**

Client: Paterson Group Consulting Engineers

Client PO: 57092

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Benzene	33.2	0.5	ug/L	ND	83.0	60-130			
Ethylbenzene	36.7	0.5	ug/L	ND	91.7	60-130			
Toluene	41.4	0.5	ug/L	ND	103	60-130			
m,p-Xylenes	77.0	0.5	ug/L	ND	96.3	60-130			
o-Xylene	38.7	0.5	ug/L	ND	96.7	60-130			
Surrogate: Toluene-d8	81.7		ug/L		102	50-140			



Report Date: 29-Mar-2023 Order Date: 24-Mar-2023

 Client:
 Paterson Group Consulting Engineers
 Order Date: 24-Mar-2023

 Client PO:
 57092
 Project Description: PE2709

Qualifier Notes:

Sample Data Revisions

Certificate of Analysis

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.

@PARACEL |||||||||||||||

Chain of Custody (Blank).xlsx



aurent Blvd. io K1G 4J8 Paracel Order Number (Lab Use Only)

pH Verified:

Chain Of Custody (Lab Use Only)

LABORATORIES LTD

Client Name: Paterson Project Ref: PE 2709 Page of Contact Name: Sam Berube Quote #: Turnaround Time Address: PO#: 57092 9 Auriga Drive ☐ 1 day ☐ 3 day E-mail: 2 day ☑ Regular SBerube @ paterson group.ca Telephone: 613 226 7381 Date Required: REG 153/04 REG 406/19 Other Regulation Matrix Type: S (Soil/Sed.) GW (Ground Water) Required Analysis ☐ Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558 □ PWQo SW (Surface Water) SS (Storm/Sanitary Sewer) ☐ Table 2 ☐ Ind/Comm ☐ Coarse P (Paint) A (Air) O (Other) ☐ CCME ☐ MISA PHCs F1-F4+BTEX ☑ Table 3 ☐ Agri/Other ☐ SU - Sani ☐ SU - Storm # of Containers ПCР ☐ Table Mun: Sample Taken Air Volume Metals by For RSC: Yes No Other: B (HWS) Matrix Vocs PAHs Ç₹ Sample ID/Location Name Date Time 1 BH3-23-GW2 3 94 March 23,2.23 X X BH5-23- 942 l Į L X 3 4 5 6 7 8 9 10 Comments: Method of Deliver Relinquished By (Sign): Blue Received By Driver/Depot: Received at Lab: Verified By: Relinquished By (Print): Bryce Lee Date/Time: Date/Time: Temperature: Temperature:

Revsion 4.0



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

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive Ottawa, ON K2E 7T9 Attn: Sam Berube

Client PO: 57798 Project: PE2709 Custody: 140789

Report Date: 7-Jul-2023 Order Date: 28-Jun-2023

Order #: 2326362

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

Paracel ID Client ID 2326362-01 BH5-23-GW3

Approved By:



Dale Robertson, BSc Laboratory Director



Report Date: 07-Jul-2023 Order Date: 28-Jun-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers
Client PO: 57798

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 624 - P&T GC-MS	2-Jul-23	2-Jul-23
PHC F1	CWS Tier 1 - P&T GC-FID	30-Jun-23	2-Jul-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	5-Jul-23	6-Jul-23



Order #: 2326362

Report Date: 07-Jul-2023 Order Date: 28-Jun-2023 Project Description: PE2709

Client: Paterson Group Consulting Engineers
Client PO: 57798

	Client ID:	BH5-23-GW3	-	-	-
	Sample Date:	27-Jun-23 09:00	-	-	-
	Sample ID:	2326362-01	-	-	-
	MDL/Units	Ground Water	-	-	-
Volatiles			•		
Benzene	0.5 ug/L	1.2	-	-	-
Ethylbenzene	0.5 ug/L	10.7	-	-	-
Toluene	0.5 ug/L	1.0	-	-	-
m,p-Xylenes	0.5 ug/L	12.4	-	-	-
o-Xylene	0.5 ug/L	0.9	-	-	-
Xylenes, total	0.5 ug/L	13.2	-	-	-
Toluene-d8	Surrogate	104%	-	-	-
Hydrocarbons			•		•
F1 PHCs (C6-C10)	25 ug/L	1050	-	-	-
F2 PHCs (C10-C16)	100 ug/L	398	-	-	-
F3 PHCs (C16-C34)	100 ug/L	<100	-	-	-
F4 PHCs (C34-C50)	100 ug/L	<100	-	-	-



Report Date: 07-Jul-2023 Order Date: 28-Jun-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 57798

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L						
F2 PHCs (C10-C16)	ND	100	ug/L						
F3 PHCs (C16-C34)	ND	100	ug/L						
F4 PHCs (C34-C50)	ND	100	ug/L						
Volatiles									
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						
Xylenes, total	ND	0.5	ug/L						
Surrogate: Toluene-d8	84.9		ug/L		106	50-140			



Report Date: 07-Jul-2023 Order Date: 28-Jun-2023

Project Description: PE2709

Certificate of Analysis
Client: Paterson Group Consulting Engineers

Client PO: 57798

Method Quality Control: Duplicate

momou quanty control 2	apnoate								
Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	ND	25	ug/L	ND			NC	30	
Volatiles									
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: Toluene-d8	84.4		ug/L		106	50-140			
Surrogate. Totalette-uo	04.4		ug/L		700	30-140			



Report Date: 07-Jul-2023 Order Date: 28-Jun-2023

Project Description: PE2709

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 57798

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	87.1	68-117			
F2 PHCs (C10-C16)	1930	100	ug/L	ND	121	60-140			
F3 PHCs (C16-C34)	4280	100	ug/L	ND	109	60-140			
F4 PHCs (C34-C50)	2420	100	ug/L	ND	97.6	60-140			
Volatiles									
Benzene	25.7	0.5	ug/L	ND	64.4	60-130			
Ethylbenzene	33.8	0.5	ug/L	ND	84.4	60-130			
Toluene	32.8	0.5	ug/L	ND	82.1	60-130			
m,p-Xylenes	67.5	0.5	ug/L	ND	84.4	60-130			
o-Xylene	32.5	0.5	ug/L	ND	81.2	60-130			
Surrogate: Toluene-d8	80.0		ug/L		100	50-140			



Client: Paterson Group Consulting Engineers

Order #: 2326362

Report Date: 07-Jul-2023 Order Date: 28-Jun-2023 Project Description: PE2709

Client PO: 57798 Proje

Qualifier Notes:

Sample Data Revisions

Certificate of Analysis

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

Paracel ID: 2326362



Paracel Order Number (Lab Use Only)

2.35pm

Chain Of Custody

(Lab Use Only)

	LABORATORI					•••••		n	d30	103	66	7			N	0 1	1078	39
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	9 Aurig	a Drie	10	Sbaubia paterangroup.ca									2 day			Regular		
Telephone:	Patersin Games Samue 9 Aurig 613-226	-7381				S	berubia	Patersar	word	щр		a			Requir	ed:		
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ain of Cus	tody (Env) xlsx						Revision 4.0				0							



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

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

Certificate of Analysis

Paterson Group Consulting Engineers

9 Auriga Drive

Ottawa, ON K2E 7T9

Attn: Sam Berube

Client PO: 58284 Project: PE2709

Report Date: 8-Sep-2023 Order Date: 1-Sep-2023

Order #: 2335491 Custody:

Paracel ID Client ID

2335491-01 BH8-23-GW1

Approved By: Mark Froto Mark Foto, M.Sc.



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58284

Report Date: 08-Sep-2023

Order Date: 1-Sep-2023

Project Description: PE2709

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date Analysis Date
PHC F1	CWS Tier 1 - P&T GC-FID	7-Sep-23 7-Sep-23
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	6-Sep-23 6-Sep-23
REG 153: VOCs by P&T GC/MS	EPA 624 - P&T GC-MS	7-Sep-23 7-Sep-23

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58284 Project Description: PE2709

	-			1		T	
	Client ID:	BH8-23-GW1	-	-	-		
	Sample Date:	31-Aug-23 09:00	-	-	-	-	-
	Sample ID:	2335491-01	-	-	-		
_	Matrix:	Ground Water	-	-	-		
Ĺ	MDL/Units						
Volatiles							
Acetone	5 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	9.6	-	-	-	-	-
Dibromochloromethane	0.5 ug/L	<0.5	-	-	-	-	-
Dichlorodifluoromethane	1 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	-	-	-	-	-
Ethylene dibromide (dibromoethane,	0.2 ug/L	<0.2	-	-	-	-	-
Ethylbenzene	0.5 ug/L	<0.5	-	-	-	-	-
Hexane	1 ug/L	<1.0	-	-	-	-	-

Report Date: 08-Sep-2023

Order Date: 1-Sep-2023

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58284 Project Description: PE2709

	Client ID:	BH8-23-GW1	-	-	-		
	Sample Date:	31-Aug-23 09:00	-	-	-	_	-
	Sample ID:	2335491-01	-	-	-		
	Matrix:	Ground Water	-	-	-		
	MDL/Units						
Volatiles	<u>'</u>			!	!		
Methyl Ethyl Ketone (2-Butanone)	5 ug/L	<5.0	-	-	-	-	-
Methyl Isobutyl Ketone	5 ug/L	<5.0	-	-	-	-	-
Methyl tert-butyl ether	2 ug/L	<2.0	-	-	-	-	-
Methylene Chloride	5 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	-	-	-	-	-
1,1,2-Trichloroethane	0.5 ug/L	<0.5	-	-	-	-	-
Trichloroethylene	0.5 ug/L	<0.5	-	-	-	-	-
Trichlorofluoromethane	1 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	112%	-	-	-	-	-
Dibromofluoromethane	Surrogate	112%	-	-	-	-	-
Toluene-d8	Surrogate	95.0%	-	-	-	-	-
Hydrocarbons				1	1		
F1 PHCs (C6-C10)	25 ug/L	<25	-	-	-	-	-
F2 PHCs (C10-C16)	100 ug/L	<476 [1]	-	-	-	-	-
F3 PHCs (C16-C34)	100 ug/L	<476 [1]	-	-	-	-	-
F4 PHCs (C34-C50)	100 ug/L	<476 [1]	-	-	-	-	-

Report Date: 08-Sep-2023

Order Date: 1-Sep-2023

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58284 Project Description: PE2709

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	%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	ND	1.0	ug/L					
Methyl Ethyl Ketone (2-Butanone)	ND	5.0	ug/L					
Methyl Isobutyl Ketone	ND	5.0	ug/L					

Report Date: 08-Sep-2023

Order Date: 1-Sep-2023



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Report Date: 08-Sep-2023 Order Date: 1-Sep-2023

Client PO: 58284

Project Description: PE2709

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	%REC	%REC Limit	RPD	RPD Limit	Notes
Methyl tert-butyl ether	ND	2.0	ug/L					
Methylene Chloride	ND	5.0	ug/L					
Styrene	ND	0.5	ug/L					
1,1,1,2-Tetrachloroethane	ND	0.5	ug/L					
1,1,2,2-Tetrachloroethane	ND	0.5	ug/L					
Tetrachloroethylene	ND	0.5	ug/L					
Toluene	ND	0.5	ug/L					
1,1,1-Trichloroethane	ND	0.5	ug/L					
1,1,2-Trichloroethane	ND	0.5	ug/L					
Trichloroethylene	ND	0.5	ug/L					
Trichlorofluoromethane	ND	1.0	ug/L					
Vinyl chloride	ND	0.5	ug/L					
m,p-Xylenes	ND	0.5	ug/L					
o-Xylene	ND	0.5	ug/L					
Xylenes, total	ND	0.5	ug/L					
Surrogate: 4-Bromofluorobenzene	90.5		%	113	50-140			
Surrogate: Dibromofluoromethane	87.8		%	110	50-140			
Surrogate: Toluene-d8	77.5		%	96.9	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58284

Report Date: 08-Sep-2023 Order Date: 1-Sep-2023

Project Description: PE2709

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons				ND			NO	00	
F1 PHCs (C6-C10)	ND	25	ug/L	ND			NC	30	
Volatiles			11	ND			NO	00	
Acetone	ND	5.0	ug/L	ND			NC	30	
Benzene	ND	0.5	ug/L	ND			NC	30	
Bromodichloromethane	1.81	0.5	ug/L	1.40			25.5	30	
Bromoform	ND	0.5	ug/L	ND			NC	30	
Bromomethane	ND	0.5	ug/L	ND			NC	30	
Carbon Tetrachloride	ND	0.2	ug/L	ND			NC	30	
Chlorobenzene	ND	0.5	ug/L	ND			NC	30	
Chloroform	3.48	0.5	ug/L	2.81			21.3	30	
Dibromochloromethane	1.95	0.5	ug/L	1.84			5.8	30	
Dichlorodifluoromethane	ND	1.0	ug/L	ND			NC	30	
1,2-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,3-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,4-Dichlorobenzene	ND	0.5	ug/L	ND			NC	30	
1,1-Dichloroethane	ND	0.5	ug/L	ND			NC	30	
1,2-Dichloroethane	ND	0.5	ug/L	ND			NC	30	
1,1-Dichloroethylene	ND	0.5	ug/L	ND			NC	30	
cis-1,2-Dichloroethylene	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	



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Order Date: 1-Sep-2023

Project Description: PE2709

Report Date: 08-Sep-2023

Client PO: 58284

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
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	98.8		%		123	50-140			
Surrogate: Dibromofluoromethane	102		%		128	50-140			
Surrogate: Toluene-d8	79.9		%		99.9	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58284

Report Date: 08-Sep-2023

Order Date: 1-Sep-2023

Project Description: PE2709

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1840	25	ug/L	ND	92.1	85-115			
F2 PHCs (C10-C16)	1540	100	ug/L	ND	96.2	60-140			
F3 PHCs (C16-C34)	4720	100	ug/L	ND	120	60-140			
F4 PHCs (C34-C50)	2710	100	ug/L	ND	109	60-140			
Volatiles									
Acetone	98.3	5.0	ug/L	ND	98.3	50-140			
Benzene	36.7	0.5	ug/L	ND	91.8	60-130			
Bromodichloromethane	38.0	0.5	ug/L	ND	95.0	60-130			
Bromoform	45.1	0.5	ug/L	ND	113	60-130			
Bromomethane	45.5	0.5	ug/L	ND	114	50-140			
Carbon Tetrachloride	38.3	0.2	ug/L	ND	95.6	60-130			
Chlorobenzene	39.5	0.5	ug/L	ND	98.7	60-130			
Chloroform	38.9	0.5	ug/L	ND	97.3	60-130			
Dibromochloromethane	43.3	0.5	ug/L	ND	108	60-130			
Dichlorodifluoromethane	39.8	1.0	ug/L	ND	99.6	50-140			
1,2-Dichlorobenzene	44.2	0.5	ug/L	ND	111	60-130			
1,3-Dichlorobenzene	42.8	0.5	ug/L	ND	107	60-130			
1,4-Dichlorobenzene	42.9	0.5	ug/L	ND	107	60-130			
1,1-Dichloroethane	34.9	0.5	ug/L	ND	87.4	60-130			
1,2-Dichloroethane	35.0	0.5	ug/L	ND	87.6	60-130			
1,1-Dichloroethylene	44.5	0.5	ug/L	ND	111	60-130			
cis-1,2-Dichloroethylene	42.6	0.5	ug/L	ND	106	60-130			
trans-1,2-Dichloroethylene	42.1	0.5	ug/L	ND	105	60-130			
1,2-Dichloropropane	33.7	0.5	ug/L	ND	84.2	60-130			
cis-1,3-Dichloropropylene	41.2	0.5	ug/L	ND	103	60-130			
trans-1,3-Dichloropropylene	42.8	0.5	ug/L	ND	107	60-130			
Ethylbenzene	36.0	0.5	ug/L	ND	90.0	60-130			
Ethylene dibromide (dibromoethane, 1,2-)	41.7	0.2	ug/L	ND	104	60-130			
Hexane	47.5	1.0	ug/L	ND	119	60-130			
Methyl Ethyl Ketone (2-Butanone)	95.4	5.0	ug/L	ND	95.4	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers

Client PO: 58284

Report Date: 08-Sep-2023

Order Date: 1-Sep-2023

Project Description: PE2709

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Methyl Isobutyl Ketone	84.4	5.0	ug/L	ND	84.4	50-140			
Methyl tert-butyl ether	92.8	2.0	ug/L	ND	92.8	50-140			
Methylene Chloride	49.4	5.0	ug/L	ND	123	60-130			
Styrene	37.1	0.5	ug/L	ND	92.8	60-130			
1,1,1,2-Tetrachloroethane	39.5	0.5	ug/L	ND	98.8	60-130			
1,1,2,2-Tetrachloroethane	33.9	0.5	ug/L	ND	84.7	60-130			
Tetrachloroethylene	45.0	0.5	ug/L	ND	113	60-130			
Toluene	37.3	0.5	ug/L	ND	93.2	60-130			
1,1,1-Trichloroethane	39.6	0.5	ug/L	ND	99.0	60-130			
1,1,2-Trichloroethane	38.3	0.5	ug/L	ND	95.8	60-130			
Trichloroethylene	38.1	0.5	ug/L	ND	95.2	60-130			
Trichlorofluoromethane	42.0	1.0	ug/L	ND	105	60-130			
Vinyl chloride	42.8	0.5	ug/L	ND	107	50-140			
m,p-Xylenes	74.5	0.5	ug/L	ND	93.1	60-130			
o-Xylene	34.5	0.5	ug/L	ND	86.2	60-130			
Surrogate: 4-Bromofluorobenzene	78.8		%		98.5	50-140			
Surrogate: Dibromofluoromethane	87.4		%		109	50-140			
Surrogate: Toluene-d8	86.6		%		108	50-140			



Certificate of Analysis

Client: Paterson Group Consulting Engineers

Order Date: 1-Sep-2023

Client PO: 58284 Project Description: PE2709

Qualifier Notes:

Sample Qualifiers:

1: Elevated Reporting Limits due to limited sample volume.

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.

Any use of these results implies your agreement that our total liabilty 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.





Chain Of Custody Paracel Order Number (Lab Use Only) (Lab Use Only) 2220491

ENDORATORIES ETD.										(0)53111										
Client Name: Paterson Group			Proje	ct Ref:	PE2709							1000	N Sec	Par	e 1 c	f 1				
Samuel Berube			Quote									\vdash	Page 1 of 1 Turnaround Time							
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Telephone: 613 226 1381			:	5be	rube@pate	ersongrou	p. ca	L.				☐ 2 day Date Required:			D≰ R	⊠ Regular				
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☐ Table 2 ☐ Ind/Comm ☐ Coarse ☐ CCME	☐ MISA		P (Paint) A (Air) O (Other)				*										\top			
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For RSC: Yes No Other:		.×					F	, s		s by			(S)							
Sample ID/Location Name		Matrix	Air V	# of	Date	Time	PHCs	VOCs	PAHs	Metals by ICP	Нg	CrV	B (HWS)							
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300 - 2319 St. Laurent Blvd Ottawa, ON, K1G 4J8 1-800-749-1947 www.paracellabs.com

Certificate of Analysis

Paterson Group Consulting Engineers (Ottawa)

9 Auriga Drive

Ottawa, ON K2E 7T9

Attn: Karyn Munch

Client PO: 60290

Project: PE6422

Custody:

Approved By:

Report Date: 31-May-2024

Order Date: 27-May-2024

Order #: 2422099

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

Client ID
BH1-24-GW1
BH3-24-GW1
BH4-24-GW1





Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290

Report Date: 31-May-2024

Order Date: 27-May-2024

Project Description: PE6422

Analysis Summary Table

Analysis	Method Reference/Description	Extraction Date	Analysis Date
BTEX by P&T GC-MS	EPA 624 - P&T GC-MS	28-May-24	28-May-24
Chromium, hexavalent - water	MOE E3056 - colourimetric	28-May-24	28-May-24
Mercury by CVAA	EPA 245.2 - Cold Vapour AA	28-May-24	28-May-24
Metals, ICP-MS	EPA 200.8 - ICP-MS	28-May-24	29-May-24
PHC F1	CWS Tier 1 - P&T GC-FID	28-May-24	28-May-24
PHCs F2 to F4	CWS Tier 1 - GC-FID, extraction	30-May-24	31-May-24
REG 153: PAHs by GC-MS	EPA 625 - GC-MS, extraction	30-May-24	31-May-24

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290 Project Description: PE6422

	Client ID:	BH1-24-GW1	BH3-24-GW1	BH4-24-GW1	-		
	Sample Date:	22-May-24 12:20	22-May-24 12:30	22-May-24 12:40	-	-	-
	Sample ID:	2422099-01	2422099-02	2422099-03	-		
	Matrix:	Ground Water	Ground Water	Ground Water	-		
	MDL/Units						
Metals	<u> </u>			•			•
Mercury	0.1 ug/L	<0.1	<0.1	<0.1	-	-	-
Antimony	0.5 ug/L	0.5	0.8	0.7	-	-	-
Arsenic	1 ug/L	<1	<1	1	-	-	-
Barium	1 ug/L	831	882	377	-	-	-
Beryllium	0.5 ug/L	<0.5	<0.5	<0.5	-	-	-
Boron	10 ug/L	403	164	172	-	-	-
Cadmium	0.1 ug/L	<0.1	<0.1	<0.1	-	-	-
Chromium (VI)	10 ug/L	<10	<10	<10	-	-	-
Chromium	1 ug/L	<1	<1	<1	-	-	-
Cobalt	0.5 ug/L	<0.5	<0.5	<0.5	-	-	-
Copper	0.5 ug/L	2.3	2.7	2.6	-	-	-
Lead	0.1 ug/L	1.1	0.4	0.6	-	-	-
Molybdenum	0.5 ug/L	5.7	9.7	12.1	-	-	-
Nickel	1 ug/L	1	2	1	-	-	-
Selenium	1 ug/L	<1	<1	4	-	-	-
Silver	0.1 ug/L	<0.1	<0.1	<0.1	-	-	-
Sodium	200 ug/L	908000	1430000	783000	-	-	-
Thallium	0.1 ug/L	0.1	0.1	0.1	-	-	-
Uranium	0.1 ug/L	1.6	3.9	2.5	-	-	-
Vanadium	0.5 ug/L	1.2	0.6	0.7	-	-	-
Zinc	5 ug/L	<5	<5	<5	-	-	-
Volatiles							
Benzene	0.5 ug/L	<0.5	-	-	-	-	-
Ethylbenzene	0.5 ug/L	<0.5	-	-	-	-	-
Toluene	0.5 ug/L	<0.5	-	-	-	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290 Project Description: PE6422

	Client ID:	BH1-24-GW1	BH3-24-GW1	BH4-24-GW1	-		
	Sample Date:	22-May-24 12:20	22-May-24 12:30	22-May-24 12:40	-	-	-
	Sample ID:	2422099-01	2422099-02	2422099-03	-		
	Matrix:	Ground Water	Ground Water	Ground Water	-		
	MDL/Units						
Volatiles				!			
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	100%	-	-	-	-	-
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	282	-	-	-	-	-
Semi-Volatiles	•						
Acenaphthene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Acenaphthylene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Anthracene	0.01 ug/L	<0.01	<0.01	<0.01	-	-	-
Benzo [a] anthracene	0.01 ug/L	<0.01	<0.01	<0.01	-	•	-
Benzo [a] pyrene	0.01 ug/L	<0.01	<0.01	<0.01	-	-	-
Benzo [b] fluoranthene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Benzo [g,h,i] perylene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Benzo [k] fluoranthene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Chrysene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Dibenzo [a,h] anthracene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Fluoranthene	0.01 ug/L	<0.01	<0.01	<0.01	-	-	-
Fluorene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Indeno [1,2,3-cd] pyrene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
1-Methylnaphthalene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
2-Methylnaphthalene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290 Project Description: PE6422

	Client ID:	BH1-24-GW1	BH3-24-GW1	BH4-24-GW1	-		
	Sample Date:	22-May-24 12:20	22-May-24 12:30	22-May-24 12:40	-	-	-
	Sample ID:	2422099-01	2422099-02	2422099-03	-		
	Matrix:	Ground Water	Ground Water	Ground Water	-		
	MDL/Units						
Semi-Volatiles	•						
Methylnaphthalene (1&2)	0.10 ug/L	<0.10	<0.10	<0.10	-	-	-
Naphthalene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Phenanthrene	0.05 ug/L	<0.05	<0.05	<0.05	-	-	-
Pyrene	0.01 ug/L	<0.01	<0.01	<0.01	-	-	-
2-Fluorobiphenyl	Surrogate	63.0%	60.5%	61.4%	-	-	-
Terphenyl-d14	Surrogate	88.3%	93.3%	96.3%	-	-	-

Report Date: 31-May-2024

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290

Report Date: 31-May-2024

Order Date: 27-May-2024

Project Description: PE6422

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	%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					
Metals								
Mercury	ND	0.1	ug/L					
Antimony	ND	0.5	ug/L					
Arsenic	ND	1	ug/L					
Barium	ND	1	ug/L					
Beryllium	ND	0.5	ug/L					
Boron	ND	10	ug/L					
Cadmium	ND	0.1	ug/L					
Chromium (VI)	ND	10	ug/L					
Chromium	ND	1	ug/L					
Cobalt	ND	0.5	ug/L					
Copper	ND	0.5	ug/L					
Lead	ND	0.1	ug/L					
Molybdenum	ND	0.5	ug/L					
Nickel	ND	1	ug/L					
Selenium	ND	1	ug/L					
Silver	ND	0.1	ug/L					
Sodium	ND	200	ug/L					
Thallium	ND	0.1	ug/L					
Uranium	ND	0.1	ug/L					
Vanadium	ND	0.5	ug/L					
Zinc	ND	5	ug/L					
Semi-Volatiles	112		3/-					
Acenaphthene	ND	0.05	ug/L					
Acenaphthylene	ND	0.05	ug/L					
Anthracene	ND	0.01	ug/L					
Benzo [a] anthracene	ND	0.01	ug/L					
Benzo [a] pyrene	ND	0.01	ug/L					

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290

Report Date: 31-May-2024

Order Date: 27-May-2024

Project Description: PE6422

Method Quality Control: Blank

Analyte	Result	Reporting Limit	Units	%REC	%REC Limit	RPD	RPD Limit	Notes
Benzo [b] fluoranthene	ND	0.05	ug/L					
Benzo [g,h,i] perylene	ND	0.05	ug/L					
Benzo [k] fluoranthene	ND	0.05	ug/L					
Chrysene	ND	0.05	ug/L					
Dibenzo [a,h] anthracene	ND	0.05	ug/L					
Fluoranthene	ND	0.01	ug/L					
Fluorene	ND	0.05	ug/L					
Indeno [1,2,3-cd] pyrene	ND	0.05	ug/L					
1-Methylnaphthalene	ND	0.05	ug/L					
2-Methylnaphthalene	ND	0.05	ug/L					
Methylnaphthalene (1&2)	ND	0.10	ug/L					
Naphthalene	ND	0.05	ug/L					
Phenanthrene	ND	0.05	ug/L					
Pyrene	ND	0.01	ug/L					
Surrogate: 2-Fluorobiphenyl	16.4		%	81.8	50-140			
Surrogate: Terphenyl-d14	25.1		%	125	50-140			
Volatiles								
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					
Xylenes, total	ND	0.5	ug/L					
Surrogate: Toluene-d8	78.1		%	97.6	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290

Report Date: 31-May-2024

Order Date: 27-May-2024

Project Description: PE6422

Method Quality Control: Duplicate

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons	ND	0.5	/!	ND			NC	20	
F1 PHCs (C6-C10)	ND	25	ug/L	ND			NC	30	
Metals	ND	0.4	//	ND			NC	20	
Mercury	ND	0.1	ug/L	ND			NC	20	
Antimony	ND	0.5	ug/L	ND			NC	20	
Arsenic	1.3	1	ug/L	1.3			1.4	20	
Barium	90.7	1	ug/L	86.8			4.4	20	
Beryllium	ND	0.5	ug/L	ND			NC	20	
Boron	36	10	ug/L	38			3.6	20	
Cadmium	ND	0.1	ug/L	ND			NC	20	
Chromium (VI)	ND	10	ug/L	ND			NC	20	
Chromium	ND	1	ug/L	ND			NC	20	
Cobalt	8.75	0.5	ug/L	9.06			3.5	20	
Copper	ND	0.5	ug/L	ND			NC	20	
Lead	0.23	0.1	ug/L	0.32			NC	20	
Molybdenum	0.50	0.5	ug/L	ND			NC	20	
Nickel	1.9	1	ug/L	1.9			3.0	20	
Selenium	ND	1	ug/L	ND			NC	20	
Silver	ND	0.1	ug/L	ND			NC	20	
Sodium	82200	200	ug/L	88500			7.4	20	
Thallium	ND	0.1	ug/L	ND			NC	20	
Uranium	0.3	0.1	ug/L	0.3			4.1	20	
Vanadium	2.17	0.5	ug/L	2.19			1.1	20	
Zinc	ND	5	ug/L	ND			NC	20	
Volatiles									
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: Toluene-d8	79.0		%		98.8	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290

Report Date: 31-May-2024 Order Date: 27-May-2024

Project Description: PE6422

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Hydrocarbons									
F1 PHCs (C6-C10)	1820	25	ug/L	ND	91.0	85-115			
F2 PHCs (C10-C16)	1230	100	ug/L	ND	76.6	60-140			
F3 PHCs (C16-C34)	3360	100	ug/L	ND	85.8	60-140			
F4 PHCs (C34-C50)	2170	100	ug/L	ND	87.5	60-140			
Metals									
Mercury	2.87	0.1	ug/L	ND	95.7	70-130			
Arsenic	53.4	1	ug/L	1.3	104	80-120			
Barium	137	1	ug/L	86.8	99.6	80-120			
Beryllium	50.6	0.5	ug/L	ND	101	80-120			
Boron	80	10	ug/L	38	84.9	80-120			
Cadmium	51.0	0.1	ug/L	ND	102	80-120			
Chromium (VI)	190	10	ug/L	ND	95.0	70-130			
Chromium	53.2	1	ug/L	ND	105	80-120			
Cobalt	59.1	0.5	ug/L	9.06	100	80-120			
Copper	48.5	0.5	ug/L	ND	96.1	80-120			
Lead	43.3	0.1	ug/L	0.32	86.0	80-120			
Molybdenum	46.6	0.5	ug/L	ND	92.3	80-120			
Nickel	51.0	1	ug/L	1.9	98.1	80-120			
Selenium	46.5	1	ug/L	ND	92.1	80-120			
Silver	42.1	0.1	ug/L	ND	84.2	80-120			
Sodium	9910	200	ug/L	ND	99.1	80-120			
Thallium	45.9	0.1	ug/L	ND	91.6	80-120			
Uranium	46.0	0.1	ug/L	0.3	91.4	80-120			
Vanadium	56.8	0.5	ug/L	2.19	109	80-120			
Zinc	46	5	ug/L	ND	90.1	80-120			
Semi-Volatiles									
Acenaphthene	4.43	0.05	ug/L	ND	88.5	50-140			
Acenaphthylene	4.74	0.05	ug/L	ND	94.9	50-140			
Anthracene	5.18	0.01	ug/L	ND	104	50-140			
Benzo [a] anthracene	4.02	0.01	ug/L	ND	80.5	50-140			

Certificate of Analysis

Client: Paterson Group Consulting Engineers (Ottawa)

Client PO: 60290

Report Date: 31-May-2024

Order Date: 27-May-2024

Project Description: PE6422

Method Quality Control: Spike

Analyte	Result	Reporting Limit	Units	Source Result	%REC	%REC Limit	RPD	RPD Limit	Notes
Benzo [a] pyrene	3.84	0.01	ug/L	ND	76.9	50-140			
Benzo [b] fluoranthene	3.98	0.05	ug/L	ND	79.5	50-140			
Benzo [g,h,i] perylene	3.87	0.05	ug/L	ND	77.3	50-140			
Benzo [k] fluoranthene	4.99	0.05	ug/L	ND	99.8	50-140			
Chrysene	4.13	0.05	ug/L	ND	82.5	50-140			
Dibenzo [a,h] anthracene	3.83	0.05	ug/L	ND	76.7	50-140			
Fluoranthene	5.06	0.01	ug/L	ND	101	50-140			
Fluorene	4.22	0.05	ug/L	ND	84.4	50-140			
Indeno [1,2,3-cd] pyrene	3.87	0.05	ug/L	ND	77.3	50-140			
1-Methylnaphthalene	4.01	0.05	ug/L	ND	80.3	50-140			
2-Methylnaphthalene	4.00	0.05	ug/L	ND	80.0	50-140			
Naphthalene	4.19	0.05	ug/L	ND	83.8	50-140			
Phenanthrene	4.55	0.05	ug/L	ND	91.0	50-140			
Pyrene	5.12	0.01	ug/L	ND	102	50-140			
Surrogate: 2-Fluorobiphenyl	15.7		%		78.5	50-140			
Surrogate: Terphenyl-d14	22.4		%		112	50-140			
Volatiles									
Benzene	35.7	0.5	ug/L	ND	89.2	60-130			
Ethylbenzene	37.3	0.5	ug/L	ND	93.3	60-130			
Toluene	37.1	0.5	ug/L	ND	92.8	60-130			
m,p-Xylenes	74.0	0.5	ug/L	ND	92.4	60-130			
o-Xylene	34.4	0.5	ug/L	ND	86.0	60-130			



Client: Paterson Group Consulting Engineers (Ottawa)

Order #: 2422099

Certificate of Analysis

Report Date: 31-May-2024

Client PO: 60290 Project Description: PE6422

Qualifier Notes:

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

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.

Any use of these results implies your agreement that our total liabilty 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.

LABORATORIES L	OPARACEL IIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIII					s.com n				Paracel Order Number (Lab Use Only) 2422099					Chain Of Custody (Lab Use Only)				
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Contact Name: Kary M.w.	uch		Quot	e #:									Т	urnaı	ound	Time			
			PO #:	60	290								1 day			Г	3 day		
9 AUNIGA Dr. NEPE Telephone: G13-226 -				E-mail: 2 day KMUNCH @ PATENSON Group. Ca Date Requ											Ä		Regular		
REG 153/04 REG 406/19	Other Regulation							_	V P	ites de la	Ro Se	Dutte	ricquii	cu.		6,030,00			
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□ Table 2 □ Ind/Comm □ Coarse □	CCME MISA		,		aint) A (Air) O (Ot		m,	1	Т			22.30			\neg				
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300 - 2319 St. Laurent Blvd Ottawa, ON, K1G 4J8 1-800-749-1947 www.paracellabs.com

Order Date:

Report Date:

14-May-24

31-May-24

Subcontracted Analysis

Paterson Group Consulting Engineers (Ottawa)

9 Auriga Drive

Ottawa, ON K2E 7T9

Attn: Jeremy Camposarcone

Paracel Report No. 2420225

Client Project(s): **PE6422**Client PO: **60196**

Reference: #24-017 Standing Offer 2024

CoC Number:

Sample(s) from this project were subcontracted for the listed parameters. A copy of the subcontractor's report is attached

Paracel ID Client ID Analysis

2420225-10 BH4-24-SS3 Methyl Mercury - soil

ALS Canada Ltd.



CERTIFICATE OF ANALYSIS

Work Order : WT2413665

Client : Paracel Laboratories Ltd

Contact : Mark Foto

Address : 2319 St. Laurent Blvd. Unit 300

Ottawa ON Canada K1G 4J8

Telephone : 613 731 9577

Project : 2420225

PO : ----C-O-C number · ----

Sampler : CLIENT

Site : ----

Quote number : Standing Offer 2024

No. of samples received : 1
No. of samples analysed : 1

Page : 1 of 2

Laboratory : ALS Environmental - Waterloo

Account Manager : Costas Farassoglou

Address : 60 Northland Road, Unit 1

Waterloo ON Canada N2V 2B8

Telephone : 613 225 8279

Date Samples Received : 28-May-2024 12:10

Date Analysis Commenced : 02-Jun-2024

Issue Date : 13-Jun-2024 09:08

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. This document shall not be reproduced, except in full.

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results

Additional information pertinent to this report will be found in the following separate attachments: Quality Control Report, QC Interpretive report to assist with Quality Review and Sample Receipt Notification (SRN).

Signatories

This document has been electronically signed by the authorized signatories below. Electronic signing is conducted in accordance with US FDA 21 CFR Part 11.

Signatories Position Laboratory Department

Kinny Wu Lab Analyst Metals, Burnaby, British Columbia

Page : 2 of 2

Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



General Comments

The analytical methods used by ALS are developed using internationally recognized reference methods (where available), such as those published by US EPA, APHA Standard Methods, ASTM, ISO, Environment Canada, BC MOE, and Ontario MOE. Refer to the ALS Quality Control Interpretive report (QCI) for applicable references and methodology summaries. Reference methods may incorporate modifications to improve performance.

Where a reported less than (<) result is higher than the LOR, this may be due to primary sample extract/digestate dilution and/or insufficient sample for analysis.

Where the LOR of a reported result differs from standard LOR, this may be due to high moisture content, insufficient sample (reduced weight employed) or matrix interference.

Please refer to Quality Control Interpretive report (QCI) for information regarding Holding Time compliance.

Key: CAS Number: Chemical Abstracts Services number is a unique identifier assigned to discrete substances

LOR: Limit of Reporting (detection limit).

Unit	Description
μg/kg	micrograms per kilogram

<: less than.

>: greater than.

Surrogate: An analyte that is similar in behavior to target analyte(s), but that does not occur naturally in environmental samples. For applicable tests, surrogates are added to samples prior to analysis as a check on recovery.

Test results reported relate only to the samples as received by the laboratory.

UNLESS OTHERWISE STATED on SRN or QCI Report, ALL SAMPLES WERE RECEIVED IN ACCEPTABLE CONDITION.

Analytical Results

Sub-Matrix: Soil/Solid		ient sample ID	BH4-24-SS3	 	 		
(Matrix: Soil/Solid)							
			Client samp	ling date / time	09-May-2024 09:00	 	
Analyte	CAS Number	Method/Lab	LOR	Unit	WT2413665-001	 	
					Result	 	
Speciated Metals							
Methylmercury (as MeHg)	22967-92-6	E538/VA	0.050	μg/kg	<0.050	 	

Please refer to the General Comments section for an explanation of any result qualifiers detected.

Please refer to the Accreditation section for an explanation of analyte accreditations.



QUALITY CONTROL INTERPRETIVE REPORT

Work Order : **WT2413665** Page : 1 of 5

Client : Paracel Laboratories Ltd Laboratory : ALS Environmental - Waterloo

Contact : Mark Foto Account Manager : Costas Farassoglou

Address : 2319 St. Laurent Blvd. Unit 300 Address : 60 Northland Road, Unit 1

Ottawa ON Canada K1G 4J8 Waterloo, Ontario Canada N2V 2B8

 Telephone
 : 613 731 9577
 Telephone
 : 613 225 8279

 Project
 : 2420225
 Date Samples Received
 : 28-May-2024 12:10

PO :--- Issue Date : 13-Jun-2024 09:08 C-O-C number :---

Quote number : Standing Offer 2024

: CLIENT

No. of samples received :1

No. of samples analysed :1

This report is automatically generated by the ALS LIMS (Laboratory Information Management System) through evaluation of Quality Control (QC) results and other QA parameters associated with this submission, and is intended to facilitate rapid data validation by auditors or reviewers. The report highlights any exceptions and outliers to ALS Data Quality Objectives, provides holding time details and exceptions, summarizes QC sample frequencies, and lists applicable methodology references and summaries.

Key

Sampler

Site

Anonymous: Refers to samples which are not part of this work order, but which formed part of the QC process lot.

CAS Number: Chemical Abstracts Service number is a unique identifier assigned to discrete substances.

DQO: Data Quality Objective.

LOR: Limit of Reporting (detection limit).

RPD: Relative Percent Difference.

Workorder Comments

Holding times are displayed as "---" if no guidance exists from CCME, Canadian provinces, or broadly recognized international references.

Summary of Outliers

Outliers : Quality Control Samples

- No Method Blank value outliers occur.
- No Duplicate outliers occur.
- No Laboratory Control Sample (LCS) outliers occur
- No Test sample Surrogate recovery outliers exist.

Outliers: Reference Material (RM) Samples

No Reference Material (RM) Sample outliers occur.

Outliers : Analysis Holding Time Compliance (Breaches)

• Analysis Holding Time Outliers exist - please see following pages for full details.

Outliers : Frequency of Quality Control Samples ■ No Quality Control Sample Frequency Outliers occur.		
		alsglobal.com

Page : 3 of 5 Work Order · WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



Analysis Holding Time Compliance

This report summarizes extraction / preparation and analysis times and compares each with ALS recommended holding times, which are selected to meet known provincial and/or federal requirements. In the absence of regulatory hold times, ALS establishes recommendations based on guidelines published by organizations such as CCME, US EPA, APHA Standard Methods, ASTM, or Environment Canada (where available). Dates and holding times reported below represent the first dates of extraction or analysis. If subsequent tests or dilutions exceeded holding times, qualifiers are added (refer to COA).

If samples are identified below as having been analyzed or extracted outside of recommended holding times, measurement uncertainties may be increased, and this should be taken into consideration when interpreting results.

Where actual sampling date is not provided on the chain of custody, the date of receipt with time at 00:00 is used for calculation purposes.

Where only the sample date without time is provided on the chain of custody, the sampling date at 00:00 is used for calculation purposes.

Matrix: Soil/Solid

Evaluation: **x** = Holding time exceedance; ✓ = Within Holding Time

Matrix Com Cond						diddion.	i loiding time exces	Judinoo ,	* * 1 (111	riolaning riiii
Analyte Group : Analytical Method	Method	Sampling Date	Extraction / Preparation			Analysis				
Container / Client Sample ID(s)			Preparation	Holding	Times	Eval	Analysis Date	Holding	g Times	Eval
			Date	Rec	Actual			Rec	Actual	
Speciated Metals : Methylmercury in Soil by GCAFS										
Glass soil jar/Teflon lined cap [ON MECP]										
BH4-24-SS3	E538	09-May-2024	07-Jun-2024	28	29	×	10-Jun-2024	28 days	3 days	✓
				days	days	EHT				

Legend & Qualifier Definitions

EHT: Exceeded ALS recommended hold time prior to analysis.

Rec. HT: ALS recommended hold time (see units).

Page : 4 of 5 Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



Quality Control Parameter Frequency Compliance

The following report summarizes the frequency of laboratory QC samples analyzed within the analytical batches (QC lots) in which the submitted samples were processed. The actual frequency should be greater than or equal to the expected frequency.

Matrix: Soil/Solid	Evaluation: × = QC frequency outside specification; ✓ = QC frequency within specification.								
Quality Control Sample Type			Co	unt		Frequency (%)	i .		
Analytical Methods	Method	QC Lot #	QC	Regular	Actual	Expected	Evaluation		
Laboratory Duplicates (DUP)									
Methylmercury in Soil by GCAFS	E538	1479163	1	11	9.0	5.0	✓		
Laboratory Control Samples (LCS)									
Methylmercury in Soil by GCAFS	E538	1479163	2	11	18.1	10.0	✓		
Method Blanks (MB)									
Methylmercury in Soil by GCAFS	E538	1479163	1	11	9.0	5.0	✓		

Page : 5 of 5 Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



Methodology References and Summaries

The analytical methods used by ALS are developed using internationally recognized reference methods (where available), such as those published by US EPA, APHA Standard Methods, ASTM, ISO, Environment Canada, BC MOE, and Ontario MOE. Reference methods may incorporate modifications to improve performance (indicated by "mod").

Analytical Methods	Method / Lab	Matrix	Method Reference	Method Descriptions
Methylmercury in Soil by GCAFS	E538 ALS Environmental - Vancouver	Soil/Solid	DeWild et al. (2004)/EPA 1630 (mod)	This method follows procedures published by DeWild, Olund, Olsen and Tate (2004) for the US Geological Survey (Techniques and Methods 5A-7). Samples are leached with an acidic copper sulphate solution to solubilize methylmercury for inorganic complexes. The methylmercury is then extracted into dichloromethane and then an aliquot is back extracted into ultra-pure water. The extract is analyzed by aqueous phase ethylation, purge and trap, desorption and GC separation. The separated species are then pyrolized to elemental Hg and quantified by cold vapour atomic flourescence spectroscopy. Results are reported "as MeHg".
Preparation Methods	Method / Lab	Matrix	Method Reference	Method Descriptions
Methylmercury Soil Digestion	EP538 ALS Environmental - Vancouver	Soil/Solid	DeWild et al. (2004)	This method follows procedures published by DeWild, Olund, Olsen and Tate (2004) for the US Geological Survey (Techniques and Methods 5A-7). Samples are leached with an acidic copper sulphate solution to solubilize methylmercury for inorganic complexes. The methylmercury is then extracted into dichloromethane and then an aliquot is back extracted into ultra-pure water. The extract is analyzed by aqueous phase ethylation, purge and trap, desorption and GC separation. The separated species are then pyrolized to elemental Hg and quantified by cold vapour atomic flourescence spectroscopy. Results are reported "as MeHg".

ALS Canada Ltd.



QUALITY CONTROL REPORT

Work Order : WT2413665

Client : Paracel Laboratories Ltd

Contact : Mark Foto

Address : 2319 St. Laurent Blvd. Unit 300

Ottawa ON Canada K1G 4J8

Telephone : 613 731 9577

Project : 2420225

PO :----C-O-C number :----

Sampler : CLIENT

Site :----

Quote number : Standing Offer 2024

No. of samples received : 1
No. of samples analysed : 1

Page : 1 of 3

Laboratory : ALS Environmental - Waterloo

Account Manager : Costas Farassoglou

Address : 60 Northland Road, Unit 1

Waterloo, Ontario Canada N2V 2B8

Telephone :613 225 8279

Date Samples Received : 28-May-2024 12:10

Date Analysis Commenced : 02-Jun-2024

Issue Date : 13-Jun-2024 09:08

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. This document shall not be reproduced, except in full.

This Quality Control Report contains the following information:

Laboratory Duplicate (DUP) Report; Relative Percent Difference (RPD) and Data Quality Objectives

- Reference Material (RM) Report; Recovery and Data Quality Objectives
- Method Blank (MB) Report; Recovery and Data Quality Objectives
- Laboratory Control Sample (LCS) Report; Recovery and Data Quality Objectives

Signatories

This document has been electronically signed by the authorized signatories below. Electronic signing is conducted in accordance with US FDA 21 CFR Part 11.

Signatories Position Laboratory Department

Kinny Wu Lab Analyst Vancouver Metals, Burnaby, British Columbia

Page : 2 of 3 Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225

ALS

General Comments

The ALS Quality Control (QC) report is optionally provided to ALS clients upon request. ALS test methods include comprehensive QC checks with every analysis to ensure our high standards of quality are met. Each QC result has a known or expected target value, which is compared against predetermined Data Quality Objectives (DQOs) to provide confidence in the accuracy of associated test results. This report contains detailed results for all QC results applicable to this sample submission. Please refer to the ALS Quality Control Interpretation report (QCI) for applicable method references and methodology summaries.

Key:

Anonymous = Refers to samples which are not part of this work order, but which formed part of the QC process lot.

CAS Number = Chemical Abstracts Service number is a unique identifier assigned to discrete substances.

DQO = Data Quality Objective.

LOR = Limit of Reporting (detection limit).

RPD = Relative Percent Difference

= Indicates a QC result that did not meet the ALS DQO.

Workorder Comments

Holding times are displayed as "---" if no guidance exists from CCME, Canadian provinces, or broadly recognized international references.

Laboratory Duplicate (DUP) Report

A Laboratory Duplicate (DUP) is a randomly selected intralaboratory replicate sample. Laboratory Duplicates provide information regarding method precision and sample heterogeneity. ALS DQOs for Laboratory Duplicates are expressed as test-specific limits for Relative Percent Difference (RPD), or as an absolute difference limit of 2 times the LOR for low concentration duplicates within ~ 4-10 times the LOR (cut-off is test-specific).

Sub-Matrix: Soil/Solid					Laboratory Duplicate (DUP) Report						
Laboratory sample ID	Client sample ID	Analyte	CAS Number	Method	LOR	Unit	Original Result	Duplicate Result	RPD(%) or Difference	Duplicate Limits	Qualifier
Speciated Metals (C	Speciated Metals (QC Lot: 1479163)										
VA24B0502-001	Anonymous	Methylmercury (as MeHg)	22967-92-6	E538	0.050	μg/kg	0.230	0.133	0.097	Diff <2x LOR	

Method Blank (MB) Report

A Method Blank is an analyte-free matrix that undergoes sample processing identical to that carried out for test samples. Method Blank results are used to monitor and control for potential contamination from the laboratory environment and reagents. For most tests, the DQO for Method Blanks is for the result to be < LOR.

Sub-Matrix: Soil/Solid

Analyte	CAS Number Method	LOR	Unit	Result	Qualifier
Speciated Metals (QCLot: 1479163)					
Methylmercury (as MeHg)	22967-92-6 E538	0.05	μg/kg	<0.050	

Page : 3 of 3 Work Order : WT2413665

Client : Paracel Laboratories Ltd

Project : 2420225



Laboratory Control Sample (LCS) Report

A Laboratory Control Sample (LCS) is an analyte-free matrix that has been fortified (spiked) with test analytes at known concentration and processed in an identical manner to test samples. LCS results are expressed as percent recovery, and are used to monitor and control test method accuracy and precision, independent of test sample matrix.

Sub-Matrix: Soil/Solid						Laboratory Co	ontrol Sample (LCS)	Report	port			
					Spike	Recovery (%)	Recovery Limits (%)					
Analyte	CAS Number	Method	LOR	Unit	Target Concentration	LCS	Low	High	Qualifier			
Speciated Metals (QCLot: 1479163)	Speciated Metals (QCLot: 1479163)											
Methylmercury (as MeHg)	22967-92-6	E538	0.05	μg/kg	10 μg/kg	92.1	70.0	130				

Reference Material (RM) Report

A Reference Material (RM) is a homogenous material with known and well-established analyte concentrations. RMs are processed in an identical manner to test samples, and are used to monitor and control the accuracy and precision of a test method for a typical sample matrix. RM results are expressed as percent recovery of the target analyte concentration. RM targets may be certified target concentrations provided by the RM supplier, or may be ALS long-term mean values (for empirical test methods).

Sub-Matrix:					Reference Material (RM) Report					
					RM Target	Recovery (%)	Recovery L	imits (%)		
Laboratory sample ID	Reference Material ID	Analyte	CAS Number	Method	Concentration	RM	Low	High	Qualifier	
Speciated Metals	Speciated Metals (QCLot: 1479163)									
QC-1479163-003	RM	Methylmercury (as MeHg)	22967-92-6	E538	14.8 μg/kg	108	70.0	130		

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

Subcontract Order

SENDING LABORATORY:

Paracel Laboratories Ltd.

300-2319 St. Laurent Blvd. Ottawa, ON K1G 4J8 Phone: 613-731-9577

Fax: 613-731-9064

Date Requested:

Project Number:

Submitted By:

15-May-24

2420225 Sarah Scullion RECEIVING LABORATORY:

ALS Laboratory Group (Vancouver)

8081 Lougheed Highway Burnaby, BC V5A 1W9 Phone: (604) 253-4188

Fax:

INVOICE TO:

Paracel Laboratories Ltd.

300-2319 St. Laurent Blvd. Ottawa, ON K1G 4J8

Phone: 613-731-9577

Fax: 613-731-9064

Required Regulation	Real 53
Turnaround Time	Standland

Sample ID	Matrix	Analyses Requested:	Sampled	Comments
BH4-24-SS3	Soil	Methyl Mercury - soil	09-May-24 09:00	

BB 05/29/24

Environmental Division Waterloo Work Order Reference



Telephone: +1 519 886 6910

46 Sublet

Please email all results to mfoto@paracellabs.com, dbloom@paracellabs.com, drobertson@paracellabs.com

Temperature prior to Shipping: /

13.5°C -> ICE PACK

12:10





nt Bhd. IG 4.JB labs.com Paracel Order Number (Lab Use Only)

2420225

Chain Of Custody (Lab Use Only)

Client Name: Patoreas Crown In																					
Contact Name						Project Ref: PE6422									Page 1 of 2						
Address:					Quote #:									Turnaround Time							
9 AURIGA DRIVE					PO#: 60196									☐ 1 day ☐ 3 day							
OTTAWA ON K2E 7T9					E-mail: jcamposarcone@patersongroup.ca								2 day				▼ Regular				
Telephone: 613-226-7381				kmunch@patersongroup.ca									Date Required:						0		
REG 153/04 REG 406/19 Other Regulation				Y 1 - 3 - 20, 2 - 3 - 20, 2 - 3 - 20, 2 - 3							177.83	- secondarion									
☐ Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558 ☐ PWQO				Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) SS (Storm/Sanitary Sewer)						equired Analysis											
	☐ Table 2 ☐ Ind/Comm ☐ Coarse ☐ CCME ☐ MISA			D (Deint) A (Air) O (Oth -)																	
		SU - Sani SU - Storm		T	2			BTEX													
0	Mun:			e e	Containers	Sample Taken		F1-F4 +			by ICP										
For RSC: ☐ Yes ☐ No ☐ Other:			Matrix	Air Volume					s,	60	ls by			(HWS)							
_	Sample ID/Location Name			Air	jo#	Date	Time	ᅥ옯	VOCs	PAHs	Metals	윤	S Z	B (H	표	S	SAR				
1 BH1-24-AU1			s		2	5/8/2024		V		V	V	V	V			V	V		Г		
2 BH1-24-SS3			S		2	5/8/2024		V	而	V	V	V	V		П	V	V	П			
3 BH1-24-SS5			s		2	5/8/2024		V	一	V	V	V	V		V	7	V	Ħ	F		
4 BH2-24-AU1			s		2	5/8/2024		7	H	7	V	7	V		7	7	V	H	H		
5 BH2-24-SS3			s		2	5/8/2024			퓜	V	7	7	V		H		7	H	F		
6 BH3-24-AU1			s		2	5/9/2024		V	H	V	V	V	V	H	H	7	V	H	-		
7 BH3-24-SS5			s		2	5/9/2024		V	믐	H	V	H	H	H	7	7	V	H			
8 BH4-24-AU1			s		2	5/9/2024		V	H	V	·	7	V	H	H		V	믐	늗		
9 BH4-24-SS2 S			S		2	5/9/2024		V	믬	~	7			H	H		<u>'</u>	ዙ	F		
10 BH4-24-SS3 S					2	5/9/2024		H	H					H	H	Ľ	H	H	L		
mments: HOLD - BH4-24-SS3				-	0/0/2024					\cup		1.65		Ш	Ш	Ш	Ш	L			
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linquished By (Print): Jeremy Camposarcone Date/Time:							SP						0 3								
te/Time: 5/13/2024 Temperature:							Date/Time: May 14, 2024 4,250 Date/						Time; 5 May 24 0928								
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Chain Of Custody Paracel Order Number (Lab Use Only) (Lab Use Only) 2420225

eit					7														
Client Name: Paterson Group Inc.	Project Ref: PE6422									Page 2 of 2									
Contact Name: Jeremy Camposarcone	Quote #:									Turnaround Time									
Address: 9 AURIGA DRIVE	PO #: 60196									☐ 1 day ☐ 3 day									
OTTAWA ON K2E 7T9	E-mail	E-mail: jcamposarcone@patersongroup.ca								☐ 2 day					× Regular				
Telephone: 613-226-7381	kmunch@patersongroup.ca									Date Required:									
☐ REG 153/04 ☐ REG 406/19 Other Regulation	SAGE TO SEE ME ALL COMPANIES							of M											
☐ Table 1 ☐ Res/Park ☐ Med/Fine ☐ REG 558 ☐ PWQO		Matrix Type: S (Soil/Sed.) GW (Ground Water) SW (Surface Water) SS (Storm/Sanitary Sewer)					Required Analysis												
☐ Table 2 ☐ Ind/Comm ☐ Coarse ☐ CCME ☐ MISA		P (Paint) A (Air) O (Other)			er)	X							П	Г					
☐ Table 3 ☐ Agri/Other ☐ SU - Sani ☐ SU - Storm			5			+ BTEX			۵										
□ Table Mun: For RSC: □ Yes □ Other:		Air Volume	f Containers	Sample Taken		F1-F4			y ICP										
							SS	₽	Metals by		_	B (HWS)			<u>«</u>				
Sample ID/Location Name			# of	Date	Time	유	VOCs	PAHs	Met	롸	CrV	B	된	EC	SAR				
1 BH4-24-SS6 S			2	5/9/2024		v			v					v	~				
2 BH5-24-AU1 S			2	5/9/2024		~		~	~	V	V		V	V	V				
3 BH5-24-SS2(BOTTOM) S			2/3	5/9/2024		~	$\overline{\Box}$	$\overline{\Box}$	~		\Box			~	V				
4 DUP1 S			2	5/8/2024		~	同	~	~	~	V	M		~	V				
5 DUP2 S			2	5/9/2024		~	$\overline{\Box}$	Ħ	~	一	Ī	Ħ	Ħ	~	V				
6						Ħ	Ħ	Ħ	Ħ	Ħ	Ħ	Ħ	Ħ	П	H				
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Date/Time: 3/3/2024 Temperature:					Ma							5 May 24 0928							
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