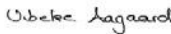



DATA VALIDATION REPORT_June 2012

URS Project number: 42213719 Data verified by: Bek Aagaard Date: 14/03/2013
 Client: Darwin Waterfront Corporation
 Site: Waterfront Precinct Signed: 
 URS Project Manager: Tim Smith Validation by: Mitch Bacon Date: 14/03/2013
 Matrix type: Groundwater
 No Primary samples: 23 Signed: 
 Laboratory: ALS; Envirolab 3/4/2013
 Lab reference: ES1214335, ES1214577, ES1214657, ES1214894, 74692 Project Manager: Tim Smith Date: 14/03/2013
 Signed:

Data quality objectives

Field data comparison No apparent anomalies were observed between laboratory results (Batches ES1214577, ES1214657 and ES12144894) and field observations.

Frequency of field QC Inter-laboratory field QC samples were not analysed to project specifications (1 in 20) for the 23 primary samples collected. Two intra-laboratory duplicates and two inter-laboratory duplicates were taken; however, one of the inter-laboratory duplicates was not forwarded to a secondary laboratory, resulting in three samples were analysed at the primary laboratory.

		Discrepancy	Frequency	Requested Frequency
Intra-laboratory Duplicates Taken	2	None noted	2	2
Inter-laboratory Duplicates Taken	2	None noted	2	2
Intra-laboratory Duplicates Analysed	3	More than target frequency	3	2
Inter-laboratory Duplicates Analysed	1	Less than target frequency	1	2
Total Frequency			4	4

The precision of the data could not adequately be assessed as per AS4482.1 recommendations. However, an assessment of the accuracy and precision of the analytical data can be measured on more than just the repeatability of the results. Although secondary laboratory data is an important part of any data validation, the lack of this data does not signify that the analytical proficiency of the primary laboratory cannot be assessed. As quality control data for the primary laboratory (such as method blanks, matrix spikes, laboratory duplicates, LCS recoveries and surrogate recoveries) is present (in the majority) and reliable, the accuracy of the primary laboratory data can be assessed. As inter-laboratory data (where reported) agrees and validates the primary laboratory data for these samples dates, there is no reason to doubt the accuracy (lack of precision is discussed on a batch by batch basis) of the results provided by the primary laboratory.

Batch ES1214335

Intra- and Inter-laboratory duplicates were not analysed at the required frequency for PAHs, trihalomethanes, oxygenated compounds, monocyclic aromatic hydrocarbons, aliphatic and aromatic halogenated compounds, carbon disulfide and total anions and cations analytes. However, the presence of laboratory duplicates for these analytes is considered sufficient for interpretation of the precision of the results except for PAHs. No laboratory duplicate was performed for PAHs; however, as PAHs were reported below LOR, both for the current GME and historically (except for field sample MW120_29102007), the precision and accuracy of the results for PAH is considered acceptable.

Batch ES1214657

No Inter-Laboratory duplicates were reported for Batch ES1214657; however, as the required frequency of Intra-laboratory field duplicates and laboratory duplicates were performed, the precision of the results is assessed as acceptable.

Frequency of laboratory QC

Batches ES1214335, ES1214577 and ES1214657

No laboratory duplicates were reported for PAHs, TPH C10-C40 and ferric iron. The precision of the data has been considered acceptable based on the presence of intra- and inter-laboratory duplicates where available and acceptable for TPH C10-C40 and ferric iron. However, no Intra- and Inter-

	<p>Laboratory duplicates were reported PAHs. The precision of the data has been considered acceptable based on the presence of other laboratory quality control samples, such as method blanks and LCS for PAHs, and as the reported results were consistent with historical results and field observations.</p> <p>Batch ES1214894</p> <p>No laboratory duplicates were reported for ferric iron. The precision of the data has been considered acceptable based on the presence of intra- and inter-laboratory duplicates were available and acceptable.</p>
Tests requested/reported	<p>Samples were analysed and reported as requested on the COC for ES1214335, ES1214577, ES1214894 and 74692; however, on the COC for Batch ES1214657 field sample MW116 was mislabelled as MW113. This sample was reported as MW116, which corresponds to the field notes. According to the Sample Receipt Notification Report (SRN), this sample was placed on hold. One VOC vial was broken in Batch ES1214657 (MW103); however, sufficient sample was available for analysis of volatile organic compounds.</p>
Limits of reporting	<p>LORs were sufficiently low to enable assessment against adopted guideline criteria with the following limitations:</p> <ul style="list-style-type: none"> • Protocol LOR for anthracene (1 µg/L) above the adopted ILs (0.4 µg/L) • Protocol LOR for benzo(a)pyrene (0.5 µg/L) above the adopted ILs (0.2 µg/L) <p>The limits of reporting were above the adopted investigation levels for anthracene and benzo(a)pyrene. As a result, potential exists for samples to contain concentrations of these analytes above the adopted investigation levels, but below detection limits. This lack of definitive data should be taken into consideration when interpreting analytical results below LOR for these PAHs.</p> <ul style="list-style-type: none"> • Protocol LOR exceeds the guideline trigger value for arsenic (0.1 mg/L) • Protocol LOR exceeds the guideline trigger value for cadmium (0.0055 mg/L) • Protocol LOR exceeds the guideline trigger value for chromium (0.0044 mg/L) • Protocol LOR exceeds the guideline trigger value for copper (0.0013 mg/L) • Protocol LOR exceeds the guideline trigger value for iron (0.2 mg/L) • Protocol LOR exceeds the guideline trigger value for lead (0.0044 mg/L) • Protocol LOR exceeds the guideline trigger value for nickel (0.07 mg/L) • Protocol LOR exceeds the guideline trigger value for zinc (0.015 mg/L) <p>The limits of reporting exceeded the guideline trigger levels for several metals. As a result, potential exists for samples to contain concentrations of these analytes above the adopted investigation levels, but below detection limits. This lack of definitive data should be taken into consideration when interpreting analytical results below these metals.</p>
Data transcription	<p>A random 10% check of the laboratory results identified no anomalies within the electronic data, the laboratory reports, and tables generated by URS.</p>
Sample management	
Chain of Custody	<p>Chain of custody documents completed.</p>
Handling and preservation	<p>Samples were correctly preserved, chilled with ice and received at the laboratory at 1.8-3.6°C.</p>
Holding time compliance	<p>Batch ES1221657</p> <p>The holding time for nitrite as N was exceeded by two days for field samples MW102_13/6/2012, MW103_13/6/2012, MW115_13/6/2012, MW116_13/6/2012, MW125_13/6/2012, MW135_13/6/2012 and MW135_13/6/2012 (ES1214657). Care should be taken with interpreting results for nitrite as N for these field samples, as reported concentrations may be biased low.</p> <p>Batch ES1221894</p> <p>The holding time for ferrous iron was exceeded by one day for field samples MW108_14/6/2012, MW121_14/6/2012 and MW129_14/6/2012 (ES1214894). Care should be taken with interpreting results for ferrous iron for these field samples, as reported concentrations may be biased low.</p>
Data precision	
Field duplicate RPDs	<p>RPDs exceeded control limits for the following sample analysis (Sample with higher reported concentrations are in bold):</p> <ul style="list-style-type: none"> • MW119_7/6/2012 and QAQC01_7/6/2012 for carbonate (74%); and • MW134_12/6/2012 and QAQC04_12/6/2012 for ionic balance (111%). <p>Although carbonate and ionic balance are not considered parameters of potential concern and no ILs has been adopted, care should be taken when using the data for these analytes for any quantitative purposes.</p>
Inter-laboratory Replicate (Field triplicate) RPDs	<p>RPDs exceeded control limits for the following sample analysis (Sample with higher reported concentrations are in bold):</p>

- **MW119_7/6/2012** and **QAQC02_7/6/2012** for manganese (179%), and similar for ionic balance (200%) and sulphate (34%).

The potential imprecision in manganese concentrations should be taken into consideration when interpreting results close to the ILs. Although manganese is not considered a chemical of potential concern and no ILs has been adopted, care should be taken when using the data for this analyte for any quantitative purposes.

Laboratory duplicate RPDs

The following laboratory duplicate RPD exceeded LOR based limits

Batch	Analyte	Data	LCL (%)	UCL (%)	Comment
ES1214894	TPH >C10 - C16	20.7%	0	20	Exceeded by 0.7%. Anonymous sample

Although this indicates a potential lack of precision for TPH >C10- C16, field duplicate RPDs are acceptable and this is not considered to affect the interpretation of the results.

Data accuracy

Laboratory control spike recovery

The following recoveries were outside control limits and may affect data interpretation:

Batch	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
ES1214335	Chromium	85.4	91	111	Recovery 15% less than LCL
ES1214335	Lead	89.3	90	110	Recovery 11% less than LCL
ES1214335	Vanadium	89.6	91	109	Recovery 10% less than LCL
ES1214577	Iodomethane	66.8	70.2	128	Recovery 33% less than LCL
ES1214577	Phenanthrene	119	62.6	116	Recovery 19% greater than UCL
ES1214657	Calcium	115	88	110	Recovery 15% greater than UCL
ES1214657	Potassium	88.1	89	109	Recovery 12% less than LCL
ES1214657	Iodomethane	66.8	70.2	128	Recovery 33% less than LCL

For LCS recoveries reported greater than the upper control limit, the potential exists for reported concentrations to be biased high by up to 19% for phenanthrene and 15% for calcium. Care should be taken with interpreting results for phenanthrene close to the adopted guidelines. As no IL has been adopted for calcium, this potential for over-reporting is not considered to affect the interpretation of the results.

For LCS recoveries reported lower than the lower control limit, the potential exists for reported concentrations to be biased low for chromium (15%), lead (11%), vanadium (10%), iodomethane (33%) and potassium (12%). Care should be taken with interpreting results for chromium and lead close to the adopted guidelines. As no IL has been adopted for vanadium, iodomethane and potassium, this potential for under-reporting is not considered to affect the interpretation of the results.

Matrix spike recovery

The following recoveries were outside control limits and may affect data interpretation:

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
Anonymous (ES1214335)	Sulphate as SO4	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level. Anonymous
MW122_07062012	Sulphate as SO4	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW122_07062012	Chloride	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Anonymous (ES1214335)	Ferrous Iron	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level. Anonymous
Anonymous (ES1214577)	Sulphate as SO4	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level. Anonymous

MW122_12062012	Methane	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW103_13062012	Sulphate as SO4	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW116_13062012	Sulphate as SO4	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Anonymous (ES1214894)	Manganese	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
ND – Not determined					

As all matrix spike issues were related to non-determination due to background levels being greater than the spike level, this is not considered to affect the interpretation of the accuracy of the results, due to the presence of other laboratory quality control data, including method blanks, LCS recoveries and matrix spikes for analytes analysed under the same analytical method (where applicable).

Batches ES1214335 and ES1214577

Matrix spike recoveries were not reported for PAHs, TPH C10-C36, calcium, magnesium, potassium, sodium, ferric iron, oxygenated compounds, monocyclic aromatic hydrocarbons and carbon disulphide. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as method blanks, LCS recoveries and surrogate recoveries (where applicable).

Batch ES1214657

Matrix spike recoveries were not reported for calcium, magnesium, potassium, sodium, oxygenated compound and carbon disulphide. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as method blanks LCS recoveries and surrogate recoveries (where applicable).

Batch ES1214894

Matrix spike recoveries were not reported for calcium, magnesium, potassium, sodium, oxygenated compounds, monocyclic aromatic hydrocarbons and carbon disulphide. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as method blanks LCS recoveries and surrogate recoveries (where applicable).

Surrogate spike recovery

The following recoveries were outside control limits and may affect data interpretation:

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
MW122_07062012	Toluene_D8	132	79	131	Recovery greater than upper data quality objective

Surrogate recoveries were reported greater than the upper data quality objective for Toluene_D8; hence, there is the potential for the volatile TPH C6-C10 and benzene, toluene, ethylbenzene and xylenes (BTEX) results to be biased high. This is not considered to affect the interpretation of the results, as volatile TPH C6-C10 and BTEX were not reported above the laboratory LOR in this sample.

Batches ES1214577, ES1214657 and ES1214894

Nitrobenzene-D5 surrogate recovery was not reported, representing a surrogate for oxygenated compounds, such as ketones. Due to the presence of other laboratory quality controls, including method blanks, matrix spikes and LCS recoveries, this is not considered to affect the interpretation of the results.

Blank monitoring

Equipment rinsate blank

Batches ES1214335 and ESES1214657

Concentrations of all analytes were reported below the LOR.

Batch ES1214577

The rinsate blank reported concentrations for chromium (0.001 mg/L), lead (0.002 mg/L), manganese (0.007 mg/L) and nickel (0.004 mg/L) above LOR in QCC01_12/6/2012. As the manganese concentration in the rinsate blank is reported well below those reported in primary samples, the potential for cross contamination by inadequate decontamination is not considered to affect the results.

The nickel concentration (0.004 mg/L) in the rinsate blank was reported at equal concentrations as in the primary samples. However, as the reported nickel concentrations in the primary samples (0.037

	<p>mg/L) were well below the adopted guidelines (0.07 mg/L) this is not considered to affect the interpretation of the results.</p> <p>As chromium and lead concentrations were reported below LOR in primary samples, it is apparent that these results are not caused by cross contamination during sampling or transportation is not considered to affect the results.</p>
Field blank	Concentrations of all analytes were reported below the LOR.
Trip blank	<p>Concentrations of all analytes were reported below the LOR.</p> <p>Batches ES1214577, ES1214657 and ES1214894</p> <p>No trip blanks were analysed; however, due to the presence of a rinsate blank , which did not report any volatile analytes above laboratory LOR, this is not considered to affect the interpretation of cross-contamination</p>
Method blank	<p>Concentrations of all analytes were reported below the LOR.</p> <p>Batches ES1214335, ES1214577 and ES1214894</p> <p>Method blank recoveries were not reported for ferric iron and alkalinity. The accuracy of alkalinity and ferric iron data could not be assessed. This should be taken into consideration when interpreting the accuracy of the results for these analytes.</p> <p>Batch ES1214657</p> <p>Method blank recoveries were not reported for alkalinity. The accuracy of alkalinity data could not be assessed. This should be taken into consideration when interpreting the accuracy of the results for this analyte.</p>

Site Name	Waerfront Precinct
Project No.	422137719
Project Manager	Tim Smith
Matrix	Water
Laboratory	ALS / EnviroLab
Batch File Number	ES1214335 / 74692

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Number of Primary Samples	Holding Times (h)	Limits of Reporting (h)	Field Blank (1 per day)	Rinsate Blank (1 per day)		Trip Blank (1 per day with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)	Matrix Spike (1 in 20)	LCS (1 per batch)	Surrogates (GC-MS organics)		
								Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required				Number Reported	Number Required	Reported
METALS ANALYSIS (EPA 8210F)	Asenic	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Calcium	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Chromium	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Copper	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Lead	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Manganese	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Nickel	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Silver	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Vanadium	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Zinc	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Zirconium	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
METALS ANALYSIS (EPA 8210F)	Barium	11	11	7	✓	✓	1	1	1	0	0	1	2	1	1	1	1	2	3	1	2	-	-
	Boron	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Chlorine	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Cobalt	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Copper	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Iron	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Manganese	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Nickel	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Silver	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Vanadium	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
	Zinc	12	12	7	✓	✓	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	✓	✓
SEM/EDS ANALYSIS (EPA 8210F)	Asenophenanthrene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Acenaphthylene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Anthracene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Benzo[a]anthracene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Benzo[a]pyrene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Benzo[b]fluoranthene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Benzo[k]fluoranthene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Benzofluoranthene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Chrysene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Dibenz[a,h]anthracene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Fluoranthene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
VOLATILES ANALYSIS (EPA 8210F)	Fluorene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Hexachlorocyclopentadiene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Naphthalene	2	2	2	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Phenanthrene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Pyrene	2	2	2	✓	✓	0	0	0	0	0	1	1	1	0	1	0	1	0	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction minus BTEX (F-1)	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
SEM/EDS ANALYSIS (EPA 8210F)	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	✓	✓
	Car. C10 Fraction	12	12	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	2				

Site Name
Waterfront Precinct
Project No.
42213719
Project Manager
Tim Smith
Matrix
Water
ALS / EnviroLab
Laboratory
Batch File Number
ESI214335 / 74692

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Primary Samples	Holding Times (d)	Limits of Reporting (lb)	Field Blank (1 per day)		Rinseate Blank (1 per day)		Trip Blank (1 per esky with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported
VOLATILES ANALYSIS (SEPI) 74A	1,2,4-Trimethylbenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	1	✓	✓	
	1,3,5-Trimethylbenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Isopropylbenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	m-Biphenylene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	p-Biphenylene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	p-Propylbenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	p-Propyltoluene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	sec-Butylbenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Styrene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	tert-Butylbenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
VOLATILES ANALYSIS (SEPI) 74B	1,1,1,2-Tetrachloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,1,1,2-Tetrachloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,1,2,2-Tetrachloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,1,2-Trichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,1-Dichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,1-Dichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,1-Dichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2-Dichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
VOLATILES ANALYSIS (SEPI) 74C	1,2-Dibromo-3-chloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2-Dibromoethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,3-Dichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Bromomethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Carbon Tetrachloride	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Chloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Chloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	sec-1,2-Dichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Diethylmethanediol	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Heptachlorobutadiene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
VOLATILES ANALYSIS (SEPI) 74D	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	Acetone	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
VOLATILES ANALYSIS (SEPI) 74E	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
VOLATILES ANALYSIS (SEPI) 74F	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
VOLATILES ANALYSIS (SEPI) 74G	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
	1,2,3-Trichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	1	1	1	✓	✓		
VOLATILES ANALYSIS (SEPI) 74H																										

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Number of Primary Samples	Holding Times (h)	Limits of Reporting (g)	Field Blank (1 per day)		Rinse Blank (1 per day)		Trip Blank (1 per day with VOCs)		Method Blank (1 per batch)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported		
METALS ANALYSIS REPORT	Arsenic	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	0	1	2	1	1	1	-	
	Calcium	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
	Chromium	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
	Copper	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
	Lead	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
METALS ANALYSIS REPORT	Manganese	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
	Nickel	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
	Vanadium	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
	Zinc	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
	Mercury	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	-	
VOLATILES ANALYSIS REPORT	Benzene	10	10	6	✓	✓	1	1	1	1	1	1	1	1	1	2	1	0	1	2	1	1	✓	
	Ethylbenzene	10	10	6	✓	✓	1	1	1	1	1	1	1	1	1	2	1	0	1	2	1	1	✓	
	Toluene & para-Xylene	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Napthalene	10	10	6	✓	✓	1	1	1	1	1	1	1	1	1	2	1	0	1	2	1	1	1	✓
	ortho-Xylene	10	10	6	✓	✓	1	1	1	1	1	1	1	1	1	2	1	0	1	2	1	1	1	✓
SEMIVOLATILES ANALYSIS REPORT	Toluene	10	10	6	✓	✓	1	1	1	1	1	1	1	1	1	2	1	0	1	2	1	1	1	✓
	Acenaphthylene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Anthracene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Benzo(a)fluoranthene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Benzo(b)fluoranthene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
VOLATILES ANALYSIS REPORT	Benzodibenzofuran	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Chrysene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Dibenz(a,h)anthracene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Fluoranthene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Fluorene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
VOLATILES ANALYSIS REPORT	Indeno(1,2,3-cd)pyrene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Naphthalene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Phenanthrene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Pyrene	6	6	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Gas - C10 Fraction	11	11	6	✓	✓	1	1	1	1	1	1	1	1	1	1	2	1	0	2	2	1	1	✓
SEMIVOLATILES ANALYSIS REPORT	Gas - C9 Fraction	11	11	6	✓	✓	1	1	1	1	1	1	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C10 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
VOLATILES ANALYSIS REPORT	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
	Gas - C9 Fraction	10	10	6	✓	✓	1	1	1	1	0	0	1	1	1	1	2	1	0	2	2	1	1	✓
METALS ANALYSIS REPORT	Mercury	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Calcium	8	8	6	✓	✓	0	0	0	0	0	0	1	1	2	1	0	1	3	1	0	1	2	-
	Magnesium	8	8	6	✓	✓	0	0	0	0	0	0	1	1	2	1	0	1	3	1	0	1	2	-
	Potassium	8	8	6	✓	✓	0	0	0	0	0	0	1	1	2	1	0	1	3	1	0	1	2	-
	Sodium	8	8	6	✓	✓	0	0	0	0	0	0	1	1	2	1	0	1	3	1	0	1	2	-
NONMETALS ANALYSIS REPORT	Mercury as N	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Nonfluorinated Laboratory Nonmetals as N	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Nonfluorinated Laboratory Nonmetals as S	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Nonfluorinated Laboratory Nonmetals as SO ₂ / Hydrocarbon	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Wet Chemistry and Preparations as SO ₂	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
METALS ANALYSIS REPORT	Sulfur as S	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Sulfur as S	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Nonfluorinated Laboratory Nonmetals as Chloride	8	8	6	✓	✓	0	0	0	0	0	0	1	1	2	1	0	1	3	1	0	1	4	-
	Nonfluorinated Laboratory Nonmetals as Phosphate	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
	Nonfluorinated Laboratory Nonmetals as Phosphate	8	8	6	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	-
VOLATILES ANALYSIS REPORT	Benzodibenzofuran	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Chrysene	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	Dibenzodioxanthene	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	2-Biphenyl NBSO	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	2-Naphthol NBSO	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
VOLATILES ANALYSIS REPORT	2-Naphthol NBSO	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	2-Naphthol NBSO	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
VOLATILES ANALYSIS REPORT	2-Naphthol NBSO	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓
	2-Naphthol NBSO	4	4	2	✓	✓	0	0	0	0	0	0	1	1	1	1	2	1	0	1	2	1	1	✓

Site Name
Project No.
Project Manager
Matrix
Laboratory
Batch File Number

Waterfront Precinct
42213719
Tim Smith
Water
ALS / EnviroLab
ES1214657

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Primary Sampling	Holding Times (a)	Limits of Reporting (b)	Field Blank (1 per day)		Rinse Blank (1 per day)		Trip Blank (1 per esky with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)		
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required
METALS ANALYSIS/ALS/EDP/01	Arsenic	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Chromium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Copper	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Lead	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Manganese	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Nickel	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Vanadium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Zinc	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Mercury	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
METALS ANALYSIS/ALS/EDP/02	Boron	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Benzene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Ethylbenzene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Heptane & para Xylene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Naphthalene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	ortho-Xylene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Toluene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Unidentified Aromatic Compounds	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Acenaphthylene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Anthracene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
VOLATILES ANALYSIS/ALS/EDP/03	Benz[a]anthracene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Benz[a]pyrene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Benz[b]fluoranthene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Benz[b]kijofluorene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Benz[k]fluoranthene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Benz[k]fluoranthene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Chrysene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Dibenz[a,h]anthracene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Fluoranthene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Indeno[1,2,3-cd]pyrene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
VOLATILES ANALYSIS/ALS/EDP/04	Naphthalene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Phenanthrene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Pyrene	4	4	2	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	DBP - C10 Fraction	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	DBP - C10 Fraction minus BTEX (F1)	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	DBP - C9 Fraction	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	DBP - C10 Fraction	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	DBP - C10 Fraction	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	DBP - C10 Fraction	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	DBP - C10 Fraction	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
VOLATILES ANALYSIS/ALS/EDP/05	Methane	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Carbonium	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Magnesium	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Potassium	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Sodium	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Ammonia as N	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Ammonia as N	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Ammonia as N	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Ammonia as N	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Ammonia as N	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
VOLATILES ANALYSIS/ALS/EDP/06	Carbonium	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Magnesium	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	1	1	1	-	-	-
	Potassium	7	7	7	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0									

Site Name Waterfront Precinct
Project No. 42213719
Project Manager Tim Smith
Matrix Water
Laboratory ALS / EnviroLab
Batch File Number ES1214657

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Primary Samples	Holding Times (h)	Limits of Reporting (lb)	Field Blank (1 per day)		Rinsate Blank (1 per day)		Trip Blank (1 per esky with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported		
VOLATILES ANALYSIS (EPA 821-A)	1,2,4-Trinitrobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,3,5-Trinitrobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	m-Bromobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	p-Bromobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	o-Bromobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	p-Bromophenol	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	o-Bromophenol	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Styrene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	tert-Butylbenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,1,1-Trichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (EPA 821-B)	1,1,1-Trichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,1,2-Trichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,1,2-Trichloroethene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,1,2,2-Tetrachloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,1,2,2-Tetrachloroethene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,1,2,2-Tetrachloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,1,2,2-Tetrachloroethene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Bromonitroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Dichlorodifluoromethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Heachlorobutadiene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (EPA 821-C)	Methylene Chloride	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Perchloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Tetrahydrofuran	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Isopropyl Alcohol	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,2-Dichloroethene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,4-Dichloro-2-butene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Trichloroethene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Trichloroethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Trichlorofluoromethane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Vinyl Chloride	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (EPA 821-D)	1,2,3-Trichlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,2,4-Trichlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,2,5-Trichlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,4-Dichlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,4-Dichlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	2-Chlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	4-Chlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Bromobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Chlorobenzene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	Carbon disulfide	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (EPA 821-E)	1,2-Dibromopropane (EPA)	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,2-Dibromopropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	2,2-Dichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	cis-1,2-Dichloropropene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,2-Dichloropropene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,3-Dichloropropene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,3-Dichloropropene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,3-Dichloropropene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,3-Dichloropropene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	trans-1,3-Dichloropropene	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (EPA 821-F)	1,2-Dibromopropane (EPA)	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	1,2-Dibromopropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	2,2-Dichloropropane	2	2	2	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	0	1	1	✓	✓
	cis-1,2-Dichloropropene	2																								

Site Name
Project No.
Project Manager
Matrix
Laboratory
Batch File Number

Waterfront Precinct
42213719
Tim Smith
Water
ALS / EnviroLab
ESI214894

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Primary Tests Reported	Holding Time (a)	Limits of Reporting (b)	Field Blank (1 per day)		Rinse Blank (1 per day)		Trip Blank (1 per day with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
						Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported		
METALS ANALYSIS/ALS/ED00F	Arsenic	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Calcium	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Chromium	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Copper	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Lead	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
METALS ANALYSIS/ALS/ED00F	Manganese	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Nickel	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Vanadium	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Zinc	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
	Mercury	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	
METALS ANALYSIS/ALS/ED00F	Benzene	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
	Ethylbenzene	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
	Toluene & para-Xylene	3	3	✓	✓	0	0	0	0	0	0	1	3	0	0	0	0	1	6	1	2	1	3	✓	✓
	Napthalene	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
	ortho-Xylene	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
VOLATILES ANALYSIS/ALS/ED00F	Toluene	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
	Styrene	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
	Acetophenone	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Benzofuran	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Chrysene	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS/ALS/ED00F	Dibenz(a,h)anthracene	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Fluoranthene	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Fluorene	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Hexachlorobiphenyl	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Naphthalene	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS/ALS/ED00F	Phenanthrene	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Pyrene	1	1	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	✓	✓
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	2	0	0	0	0	1	4	1	2	1	2	✓	✓
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS/ALS/ED00F	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS/ALS/ED00F	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS/ALS/ED00F	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS/ALS/ED00F	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS/ALS/ED00F	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS/ALS/ED00F	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
	Chl. - C10 Fraction	3	3	✓	✓	0	0	0	0</																

Chem Group		ChemName	Output unit	EQI	ANZECC-2000 Marine water 95%										ANZECC-2000 Irrigation LVI																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																									
Total Petroleum Hydrocarbons		Total Dissolved Solids @180 °C	mg/L	10																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																				

Table 1
Analytical Results - June 2012 GME
Darwin Waterfront Corporation

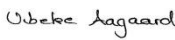

Location Code		MW09	MW10	MW101	MW102	MW103	MW108	MW113	MW114	MW115	MW116	MW117	MW118	MW119
Sample Date Time		03/06/2012	03/06/2012	12/06/2012	13/06/2012	13/06/2012	13/06/2012	12/06/2012	12/06/2012	03/06/2012	03/06/2012	03/06/2012	03/06/2012	03/06/2012
Sample Type		Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335	Prms14335
Batch ID		ES1214335	ES1214335	ES1214577	ES1214657	ES1214657	ES1214834	ES1214877	ES1214577	ES1214657	ES1214657	ES1214335	ES1214335	ES1214335
1,1,1-Trichloroethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,1-Dichloroethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,2-Dichloroethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Carbon Tetrachloride		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Chloroethane		µg/L	50	-	-	-	-	-	-	-	-	-	-	-
Cis-1,2-Dichloroethene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Cis-1,3-Dichloropropene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Hexachlorocyclopentadiene		µg/L	50	-	-	-	-	-	-	-	-	-	-	-
Hexachlorobenzene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Tetrahydrofuran		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
trans-1,2-Dichloroethene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Trichloroethene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Trichlorofluoromethane		µg/L	50	-	-	-	-	-	-	-	-	-	-	-
Vinyl chloride		µg/L	50	-	-	-	-	-	-	-	-	-	-	-
1,1,1,2-Tetrachloroethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,1,2,2-Tetrachloroethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,1-Dichloroethene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,1-Dichloropropane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,2,3-Trichloropropane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,2-Dibromo-3-chloropropane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,3-Dichloropropane		µg/L	50	-	-	-	-	-	-	-	-	-	-	-
Chloromethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Cis-1,4-Dichloro-2-butene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Dibromomethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Iodomethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Perfluoromethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Perfluoroethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Bromomethane		µg/L	50	-	-	-	-	-	-	-	-	-	-	-
1,2-Dibromoethane (EDB)		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
1,2-Dichloropropane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
2,2-Dichloropropane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Cis-1,3-Dichloropropene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
trans-1,3-Dichloropropene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Bromodichloromethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Bromotoluene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Dibromochloromethane		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Chlorobenzene		µg/L	5	-	-	-	-	-	-	-	-	-	-	-
Total Organic Carbon		mg/L	-	-	3	2	-	-	-	3	4	-	-	-
Total Dissolved Solids		mg/L	-	-	782	0.069	-	20,100	6330	-	-	-	-	-
C1 - C4 Hydrocarbon Gases		mg/L	0.005	-	-	0.069	9.88	0.095	0.04	0.631	0.034	0.147	1.22	<0.01
Alkalinity		mg/L	119	254	59.78	173	407	351.36	117.12	801	114	31	250	235
Carbonate as CO3		mg/L	1	<1	0.6	<1	<1	0.6	0.6	<1	<1	<1	<1	13
Total Alkalinity		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
Bicarbonate Alkalinity as CaCO3		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
Carbonate Alkalinity as CaCO3		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
Hydroxide Alkalinity as CaCO3		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
Total Alkalinity as CaCO3		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
Nitrate		mg/L	119	254	49	167	173	288	96	801	114	31	250	248
Nitrite (as N)		mg/L	0.01	<0.01	-	0.92	0.22	0.03	-	<0.01	<0.1	<0.1	<0.01	<0.01
Nitrate & Nitrite (as N)		mg/L	0.005	<0.01	0.19	0.21	0.05	0.01	<0.01	<0.1	<0.02	0.04	<0.01	<0.01
Chloride		mg/L	0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01
Calcium (Dissolved)		mg/L	19,500	18,500	392	1900	357	1680	5000	17,900	622	129	30	231
Calcium (Dissolved) as CaCO3		mg/L	405	400	28	312	54	228	228	185	78	6	52	8
Magnesium (Dissolved)		mg/L	477	465	9	43	21	233	16	122	16	10	31	24
Sodium (Dissolved)		mg/L	11,200	10,900	241	982	203	1110	6400	2490	425	80	34	288
Sulphate (as SO4)		mg/L	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2
Sulphate (as SO3-)		mg/L	885	761	30	220	23	458	187	258	99	10	2	30
Sulphur (as S)		mg/L	2780	2440	-	697	72	-	-	698	334	37	6	114
Sulphate (as SO4-) (Dissolved)		mg/L	1	-	-	-	-	-	-	-	-	-	-	-
Sulphate (as SO4-) (Dissolved) as CaCO3		mg/L	0.01	57	94	71	15	1510	559	555	258	150	57	13.8
Total Calcium		mg/L	839	813	14.5	713	13.6	382	154	565	25.8	4.91	5.7	14.5
Total Calcium as CaCO3		mg/L	247	213	1.64	1.28	4.88	2.25	0.19	2.8	1.83	1.2	0.29	2.24
Total Calcium as CaCO3		mg/L	<0.05	739	<0.05	<0.05	0.34	<0.05	9.37	2.34	22.4	0.24	12.4	<0.05
Ferrous Iron		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
Fentic Iron		mg/L	<0.05	<0.05	<0.05	<0.05	<0.05	11.2	0.93	0.25	2	<0.05	0.3	0.13
Oxygenated Compounds		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
2-Butanone (MEK)		µg/L	50	<50	-	-	-	-	-	-	-	-	<50	-
2-Hexanone (MIBK)		µg/L	50	<50	-	-	-	-	-	-	-	-	<50	-
Sulfolated Compounds		mg/L	-	-	-	-	-	-	-	-	-	-	-	-
Chlorinated Hydrocarbons		µg/L	5	<10	-	-	-	-	-	-	-	-	<10	-

*LOR exceeds Adopted Guidelines

Table 1
Analytical Results - June 2012 GME
Darwin Waterfront Corporation

Location Code		MW135	MW136	QA01	QA02	QA01	QA02	QC01	QC02	QC01	QC02	QC01	QC02
Field ID	Sample Type	Primary sample	Primary sample	Trip Blank	Trip Blank	Trip Blank	Trip Blank	Field Blank	Field Blank	Field Blank	Field Blank	Field Blank	Field Blank
Sample ID	Batch ID	ESI/214657	ESI/214657	ESI/214657	ESI/214657	ESI/214657	ESI/214657	ESI/214385	ESI/214385	ESI/214385	ESI/214385	ESI/214577	ESI/214577
Halogenated Aliphatic Compounds													
1,1,1-Trichloroethane	µg/L	5											
1,1,1-Dichloroethane	µg/L	5											
1,1,2-Dichloroethane	µg/L	5											
1,2-Dichloroethane	µg/L	5											
Chloroethane	µg/L	5											
Chloroethene	µg/L	5											
cis-1,2-Dichloroethane	µg/L	5											
Dichlorodifluoromethane	µg/L	5											
Hexachlorobutadiene	µg/L	5											
trans-1,2-Dichloroethene	µg/L	5											
1,1,2,2-Tetrachloroethane	µg/L	5											
1,1,1,2-Tetrachloroethane	µg/L	5											
1,1,2,2-Tetrachloroethane	µg/L	5											
1,1-Dichloropropene	µg/L	5											
1,2,3-Trichloropropene	µg/L	5											
1,2-Dichloro-3-chloropropene	µg/L	5											
1,2-Dichloropropene	µg/L	5											
Chloropropene	µg/L	5											
cis-1,4-Dichloro-2-butene	µg/L	5											
Dibromomethane	µg/L	5											
Iodomethane	µg/L	5											
Pentachloroethane	µg/L	5											
trans-1,4-Dichloro-2-butene	µg/L	5											
Bromomethane	µg/L	50											
1,2-Dibromodichloroethane [EUB]	µg/L	5											
1,2-Dibromodichloroethane	µg/L	5											
2,2-Dichloropropene	µg/L	5											
cis-1,3-Dichlorocyclopentene	µg/L	5											
trans-1,3-Dichlorocyclopentene	µg/L	5											
Bromodichloromethane	µg/L	5											
Bromochloromethane	µg/L	5											
Chloroform	µg/L	5											
Chloroform	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											
1,1,1,2,2-Pentachloroethane	µg/L	5											

DATA VALIDATION REPORT_September 2012

URS Project number: 42213719 **Data verified by:** Bek Aagaard **Date:** 26/3/2013
Client: Darwin Waterfront Corporation
Site: Waterfront Precinct **Signed:** 
URS Project Manager: Tim Smith **Validation by:** Mitch Bacon **Date:** 3/4/2013
Matrix type: Groundwater
No Primary samples: 23
Laboratory: ALS; Envirolab **Signed:** 
Lab reference: ES1221357, ES1221483, ES1221566, 78400, 78502
Project Manager: Tim Smith **Date:**
Signed:

Data quality objectives

Field data comparison

No apparent anomalies were observed between laboratory results and field observations (Batches ES1221483 and ES1221566). However, slight hydrocarbon odours were noted for MW125_4092012 and MW128_4092012, but no Total Petroleum Hydrocarbons (TPHs) or benzene, toluene, ethylbenzene and xylenes (BTEX) were reported for MW125_4092012 and MW128_4092012 in Batch ES1214357.

Frequency of field QC

Field QC samples were collected to project specifications.

		Discrepancy	Frequency	Requested Frequency
Intra-laboratory Duplicates Taken	2	None noted	2	2
Inter-laboratory Duplicates Taken	2	None noted	2	2
Intra-laboratory Duplicates Analysed	2	None noted	2	2
Inter-laboratory Duplicates Analysed	2	None noted	2	2
Total Frequency			4	4

Batch ES1221357

Inter-laboratory duplicates were not analysed at the required frequency for carbon disulfide and total anions and cations analytes. However, the presence of laboratory duplicates for these analytes is considered sufficient for interpretation of the precision of the results.

Batch ES1221483

Intra- and inter-laboratory duplicates were not analysed at the required frequency for PAHs, trihalomethanes, oxygenated compounds, monocyclic aromatic hydrocarbons, aliphatic and aromatic halogenated compounds, carbon disulfide and total anions and cations analytes. However, the presence of laboratory duplicates for these analytes, which reported RPDs within acceptable levels and acceptable accuracy data, the reported data is considered sufficient for interpretation of the precision of the results except for PAHs. No laboratory duplicate was performed for PAHs; however, as PAHs were reported below LOR, both for the current GME and historically (except for field sample MW120_29102007), the precision and accuracy of the results for PAH is considered acceptable.

Frequency of laboratory QC

Batch ES1221357

No laboratory duplicates were reported for PAHs and total anions and cations. The precision of the data has been considered acceptable based on the presence of intra- and inter-laboratory duplicates were available and acceptable for PAHs. However, no inter- or intra-laboratory duplicates were reported for total anions and cations. The precision of the data has been considered acceptable based on the presence of other laboratory quality control samples, such as LCS and method blank for total anions and cations, and surrogate recoveries (where applicable), LCS recoveries and method blank for PAHs, and as the reported results were consistent with historical results and field observations.

Batches ES1221483 and ES1221566

No laboratory duplicates were reported for PAHs, TPH >C10-C40 and ferric iron. The precision of the data has been considered acceptable based on the presence of intra- and inter-laboratory duplicates

Tests requested/reported	<p>were available and acceptable for TPH >C10-C40 and ferric iron. However, no inter-laboratory duplicates were reported for PAHs. The precision of the data has been considered acceptable based on the presence of other laboratory quality control samples, such as method blanks, LCS and surrogate recoveries for PAHs, and as the reported results were consistent with historical results and field observations.</p> <p>Samples were analysed and reported as requested on the COC for batches ES1221483 and ES1221566.</p> <p><u>Batch ES1221357</u> No amber bottle was received for MW121_04092012, therefore TPH C10-C39 and PAHs analysis could not be conducted.</p>
Limits of reporting	<p>LORs were sufficiently low to enable assessment against adopted guideline criteria with the following limitations:</p> <ul style="list-style-type: none"> • Protocol LOR for anthracene (1 µg/L) above the adopted ILs (0.4 µg/L) • Protocol LOR for benzo(a)pyrene (0.5 µg/L) above the adopted ILs (0.2 µg/L) <p>The limits of reporting were above the adopted investigation levels for anthracene and benzo(a)pyrene. As a result, potential exists for samples to contain concentrations of these analytes above the adopted investigation levels, but below detection limits. This lack of definitive data should be taken into consideration when interpreting analytical results below LOR for these PAHs.</p> <ul style="list-style-type: none"> • Protocol LOR exceeds the guideline trigger value for arsenic (0.1 mg/L) • Protocol LOR exceeds the guideline trigger value for cadmium (0.0055 mg/L) • Protocol LOR exceeds the guideline trigger value for chromium (0.0044 mg/L) • Protocol LOR exceeds the guideline trigger value for copper (0.0013 mg/L) • Protocol LOR exceeds the guideline trigger value for iron (0.2 mg/L) • Protocol LOR exceeds the guideline trigger value for lead (0.0044 mg/L) • Protocol LOR exceeds the guideline trigger value for nickel (0.07 mg/L) • Protocol LOR exceeds the guideline trigger value for zinc (0.015 mg/L) <p>The limits of reporting exceeded the guideline trigger levels for several metals. As a result, potential exists for samples to contain concentrations of these analytes above the adopted investigation levels, but below detection limits.</p> <p><u>Batch ES1221357</u> EG020: Some samples were diluted and rerun due to matrix interference and LOR's have been raised accordingly. (High sample salinity).</p> <p><u>Batch ES1221483</u> EG020A: LOR 's have been raised for some samples due to matrix interference (High sample salinity) It has been noted that Nitrite is greater than NoX for MW114; however, this difference is within the limits of experimental variation.</p> <p><u>Batch ES1221566</u> EG020: Some samples were diluted and rerun due to matrix interference and LOR's have been raised accordingly. (High sample salinity).</p>
Data transcription	<p>A random 10% check of the laboratory results identified no anomalies within the electronic data, the laboratory reports, and tables generated by URS.</p>
Sample management	
Chain of Custody	<p>Chain of custody documents completed. The sampling date on the COC for batch ES1221357 (dated 4/8/12) does not correspond to labelling in field notes (dated 4/9/12). The reported samples were labelled according to field notes (4/9/12).</p>
Handling and preservation	<p>Samples were correctly preserved, chilled with ice and received at the laboratory at 3.9-12.8°C (ice present). While the recorded temperature for batch ES1221566 was slightly above the recommended temperature, potentially resulting in some loss of volatiles and changes in natural attenuation parameters, and therefore possible under-reporting of associated analytes. This is not considered to affect the reported results, as concentrations of volatile analytes for the current and historical investigation have reported below the LOR and analytes were below the LOR in the trip blank (QCA_6092012), and the majority of the samples were received within the suggested range, and samples were cooled on packaging and on arrival at the laboratory.</p>
Holding time compliance	<p><u>Batch ES1221566</u> The holding time for nitrite as N was exceeded by one day and ferrous iron by two days for field samples MW134_5/9/2012. Care should be taken with interpreting results for ferrous iron for these field samples, as reported concentrations may be biased low.</p>

Data precision

Field duplicate RPDs

RPDs exceeded control limits for the following sample analysis. (Samples with higher reported concentrations are in bold).

- MW127_4/9/2012 and **QAQC01_4/9/2012** for ferrous iron in batch ES1221357 (185%); and
- **MW127_4/9/2012** and QAQC01_4/9/2012 for ferric iron batch ES1221357 (196%).

The reported result for ferric iron in the primary sample MW127_4/9/2012 is considered unlikely in a filtered sample. It is likely this individual sample was not field filtered, or the sample has been reduced in transit, therefore it is not considered to affect the interpretation of other ferric iron results.

EG051G: Ferrous Iron results have been confirmed for sample ID's 'MW127' and 'QAQC01' from the HCl preserved bottle.

Although ferrous iron and ferric iron are not considered parameters of potential concern and no ILs have been adopted, care should be taken when using the data for these analytes for any quantitative purposes, particularly in relation to ferric iron concentrations reported in MW127_4/9/2012, as concentrations were orders of magnitude higher than those reported in field duplicates.

Inter-laboratory Replicate (Field triplicate) RPDs

RPDs exceeded control limits for the following sample analysis (Sample with higher reported concentrations are in bold):

- MW127_4/9/2012 and **QAQC02_4/9/2012** for ferrous iron in batch ES1221357 (172%).

Although ferrous iron is not considered a parameter of potential concern and no ILs has been adopted, care should be taken when using the data for this analyte for any quantitative purposes, as the concentrations was orders of magnitude higher than the concentration reported in the primary sample.

Laboratory duplicate RPDs

RPDs were within control limits.

Data accuracy

Laboratory control spike recovery

The following recoveries were outside control limits and may affect data interpretation:

Batch	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
ES1221357	Dibromochloromethane	64.7	65	115	Recovery less than lower control limit
ES1221566	TPH C6-C10 fraction	128	75	127	Recovery greater than upper control limit

For LCS recoveries reported lower than the lower control limit, the potential exists for reported concentrations to be biased low for dibromochloromethane (35%). As no IL has been adopted for dibromochloromethane, this potential for under-reporting is not considered to affect the interpretation of the results, and dibromochloromethane was not reported above LOR.

For LCS recoveries reported greater than the upper control limit, the potential exists for reported concentrations to be biased high by up to 27% for TPH C6-C10 fraction. As no IL has been adopted for TPH C6-C10 fraction, this potential for over-reporting is not considered to affect the interpretation of the results.

Matrix spike recovery

The following recoveries were outside control limits and may affect data interpretation:

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
Anonymous (ES1221357)	Sulfate as SO ₄ - Turbidimetric	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW125_04092012	Cadmium	23.4	70	130	Recovery less than lower data quality objective
MW125_04092012	Copper	35	70	130	Recovery less than lower data quality objective
MW125_04092012	Lead	68.7	70	130	Recovery less than lower data quality objective
MW125_04092012	Zinc	39.6	70	130	Recovery less than lower data quality objective
MW128_04092012	Methane	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Anonymous (ES1221483)	Sulfate as SO ₄ - Turbidimetric	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW129_05092012	Methane	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Anonymous (ES1221566)	Methane	ND	-	-	MS recovery not determined, background level greater than

					or equal to 4x spike level.
Anonymous (ES1221566)	Methane	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
ND – Not determined					

Batches ES1221357, ES1221483 and ES1221566

For the anonymous samples, as the lack of matrix spike for these analytes was due to elevated sample concentrations was in anonymous samples, and due to the presence of other quality control data, such as method blanks and LCS, this is not considered to affect the interpretation of the results.

Matrix spike recoveries were reported less than the lower data quality objective for cadmium, copper, lead and zinc; hence, there is potential for reported concentrations of these analytes to be biased low by up to 77%, 65%, 31% and 60%, respectively. As ILs have been adopted for cadmium, copper, lead and zinc, care should be taken with interpreting results close to the adopted guidelines.

Batch ES1221357

Matrix spike recoveries were not reported for PAHs, calcium, magnesium, potassium, sodium, halogenated methanes, oxygenated compounds, monocyclic aromatic hydrocarbons, carbon disulphide, chlorinated aliphatics, ionic balance, TDS and alkalinity. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as method blanks, LCS recoveries and surrogate recoveries (where applicable).

Batch ES1221483

Matrix spike recoveries were not reported for PAHs, TPH >C10-C40, calcium, magnesium, potassium, sodium, ferric iron, halogenated methanes, oxygenated compounds, monocyclic aromatic hydrocarbons, carbon disulphide, chlorinated aliphatics, ionic balance, TDS and alkalinity. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as method blanks, LCS recoveries and surrogate recoveries (where applicable).

Batch ES1221566

Matrix spike recoveries were not reported for ferric iron and alkalinity. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as method blanks, LCS recoveries and surrogate recoveries (where applicable).

Surrogate spike recovery

The following recoveries were outside control limits and may affect data interpretation:

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
QC01_04092012	4-Bromofluorobenzene	124	80.8	123.7	Recovery greater than upper data quality objective
MW135_04092012	4-Bromofluorobenzene	126	80.8	123.7	Recovery greater than upper data quality objective
MW136_04092012	4-Bromofluorobenzene	126	80.8	123.7	Recovery greater than upper data quality objective
MW121_04092012	Toluene-D8	133	79	131	Recovery greater than upper data quality objective
QC01_04092012	4-Bromofluorobenzene	136	70	128	Recovery greater than upper data quality objective
MW136_04092012	4-Bromofluorobenzene	132	70	128	Recovery greater than upper data quality objective

Surrogate recoveries were reported greater than the upper data quality objective for 4-Bromofluorobenzene in field samples QC01_04092012 (VOC surrogate and TPH/BTEX surrogate);, MW135_04092012, MW136_04092012 (VOC surrogate and TPH/BTEX surrogate); hence, there is the potential for the volatile TPH C6-C10 and benzene, toluene, ethylbenzene and xylenes (BTEX) results to be biased high. This is not considered to affect the interpretation of the results, as volatiles were not reported above the laboratory LOR in this sample.

Surrogate recoveries were reported greater than the upper data quality objective for Toluene_D8; hence, there is the potential for the volatile analytes results to be biased high. This is not considered to affect the interpretation of the results, as volatile analytes were not reported above the laboratory LOR in these samples.

Blank monitoring

Equipment rinsate blank

Batches ES1221357 and ES1221566

Concentrations of all analytes were reported below the LOR.

Batch ES1221483

	The rinsate blank reported concentrations for manganese (0.003 mg/L) above LOR in QCC01_5/09/2012. As the manganese concentration in the rinsate blank is reported well below those reported in primary samples, the potential for cross contamination by inadequate decontamination is not considered to affect the results.
Field blank	Concentrations of all analytes were reported below the LOR.
Trip blank	Concentrations of all analytes were reported below the LOR.
Method blank	Concentrations of all analytes were reported below the LOR.
Other observations	<p><u>Batch ES1221357</u> EG020: Poor matrix spike recovery was obtained for some elements on sample ES1221357 #004 due to matrix interference. Confirmed by reanalysis.</p> <p><u>Batch ES1221483</u> EG020: Positive Manganese results have been confirmed by reanalysis for sample ES1221483 #011.</p>

Site Name
Waterfront Precinct
Project No.
42213719
Project Manager
Tim Smith
Matrix
Water
Laboratory
ALS / EnviroLab
Batch File Number
ESI221357

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Primary Samples	Holding Times (h)	Limits of Reporting (g)	Field Blank (1 per day)		Rinseate Blank (1 per day)		Trip Blank (1 per esky with VOCs)		Method Blank (1 per batch)		Inter-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported
VOLATILES ANALYSIS (ALSEP/DA)	1,2,4-Trimethylbenzene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,3,5-Trimethylbenzene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	Isopropylbenzene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	m-Xylylbenzene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	p-Xylylbenzene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	sec-Butylbenzene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	Styrene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	Isobutylbenzene	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1,1,2-Tetrachloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1,1-Trichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
VOLATILES ANALYSIS (ALSEP/DA)	1,1,2-Trichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1,2,2-Tetrachloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1,2,3-Tetrachloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1,2,3-Tetrachloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1-Dichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1-Dichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1-Dichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1-Dichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1-Dichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,1-Dichloroethane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
VOLATILES ANALYSIS (ALSEP/DA)	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2-Dibromochloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
VOLATILES ANALYSIS (ALSEP/DA)	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
VOLATILES ANALYSIS (ALSEP/DA)	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
VOLATILES ANALYSIS (ALSEP/DA)	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
VOLATILES ANALYSIS (ALSEP/DA)	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1	✓	✓	
	1,2,3-Trichloropropane	9	9	6	✓	✓	1	1	1	1	0	0	1	1	1	0	0	1	2	1	0	1	1			

Site Name
Project No.
Project Manager
Matrix
Laboratory
Batch File Number

Waterfront Precinct
42213719
Tim Smith
Water
ALS / EnviroLab
ESI221483

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Number of Primary Samples	Holding Times (h)	Limits of Reporting (g)	Field Blank (1 per day)		Rinse Blank (1 per day)		Trip Blank (1 per day with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	
METALS ANALYSIS (EOD)F	Asmet	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Calcium	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Chromium	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Copper	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Lead	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
VOLATILES ANALYSIS (EOD)F	Manganese	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Nickel	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Vanadium	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Zinc	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
	Mercury	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-
METALS ANALYSIS (EOD)F	Benzene	12	12	7	✓	✓	1	1	1	1	1	1	2	1	1	1	1	1	1	2	3	1	2	1	✓
	Ethylbenzene	12	12	7	✓	✓	1	1	1	1	1	1	2	1	1	1	1	1	1	2	3	1	2	1	✓
	m,p-Xylene	12	12	7	✓	✓	1	1	1	1	1	1	2	1	1	1	1	1	1	2	3	1	2	1	✓
	Naphthalene	12	12	7	✓	✓	1	1	1	1	1	1	2	1	1	1	1	1	1	2	3	1	2	1	✓
	ortho-Xylene	12	12	7	✓	✓	1	1	1	1	1	1	2	1	1	1	1	1	1	2	3	1	2	1	✓
VOLATILES ANALYSIS (EOD)F	Toluene	12	12	7	✓	✓	1	1	1	1	1	1	2	1	1	1	1	1	1	2	3	1	2	1	✓
	Nonhalogenated	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Acetophenone	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Aniline	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzofuran	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Chrysene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Dibenz[a,h]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[a]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Benzo[b]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[k]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[e]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[a]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Chrysene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Dibenz[a,h]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[a]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[b]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Benzo[k]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[e]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[a]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Chrysene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Dibenz[a,h]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[a]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[b]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[k]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Benzo[e]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[a]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Chrysene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Dibenz[a,h]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[a]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[b]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[k]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[e]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Benzo[a]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Chrysene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Dibenz[a,h]anthracene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
VOLATILES ANALYSIS (EOD)F	Benzo[a]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[b]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[k]fluoranthene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓
	Benzo[e]pyrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	1	1	1			

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Primary Samples	Holding Times (h)	Limits of Reporting (g)	Field Blank (1 per day)		Rinse Blank (1 per day)		Trip Blank (1 per esky with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported
VOLATILES ANALYSIS (ALSEP/924)	1,2,4-Trimethylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	1,3,5-Trimethylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Isopropylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	m-Xylylene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	p-Xylylene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	n-Propylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	o-Propylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	p-Propylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	sec-Butylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Styrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/924)	tert-Butylbenzene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	1,1,2-Trichloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	1,1,1-Trichloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	1,1,2,2-Tetrachloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	1,1,2,2-Tetrachloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	1,1-Dichloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	1,2-Dichloropropane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Bromonitrene	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Carbon Tetrachloride	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/924)	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/924)	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/924)	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/924)	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/924)	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0	0	1	1	✓	✓
	Chloroethane	3	3	1	✓	✓	1	1	1	1	0	0	1	1	1	0	1	0	1	1	0					

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Primary Samples	Holding Times (a)	Limits of Reporting (b)	Field Blank (1 per day)		Rinse Blank (1 per day)		Trip Blank (1 per day with VOCs)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 10)		LCS (1 per batch)		Surrogates (GC-MS organics)	
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Reported	OK
METALS ANALYSIS (EPCO)F	Ammonia	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Calcium	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Chromium	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Copper	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Lead	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
METALS ANALYSIS (EPCO)F	Manganese	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Nickel	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Vanadium	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Zinc	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
	Mercury	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	2	2	1	1	2	-
METALS ANALYSIS (EPCO)F	Benzene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
	Ethylbenzene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
	m,p-xylene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
	o-xylene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
	Toluene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
SEMIVOLATILES ANALYSIS (EPCO)F	Acetophenone	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Acetone	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Acrylonitrile	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Benzofuran	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Benzofuran	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Benzofuran	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Benzofuran	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Fluorene	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Hexachlorocyclopentadiene	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Naphthalene	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Phenanthrene	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Pyrene	6	6	4	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
	Chlorobenzene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
	Chlorobenzene	10	10	7	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	2	✓
SEMIVOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
VOLATILES ANALYSIS (EPCO)F	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0	1	1	0	0	1	1	1	✓
	Chlorobenzene	9	9	7	✓	✓	1	1	1	1	0	0	0	0	0	0								

Table 2
Analytical Results - September 2012 GME
Darwin Waterfront Corporation

Chem. Group		ChemName	EQL	output unit	ANZECC 2000 Marine water 95%	ANZECC 2000 Irrigation 1 & 2	Location Code	MW09	MW10	MW101	MW102	MW102	MW102	MW103	MW108	MW113	MW114	MW115
							Field ID Date Time Sample Type Lab Report Number	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012	6095012 P-6095012
								ESI1221466	ESI1221466	ESI1221466	ESI1221466	ESI1221466	ESI1221466	ESI1221466	ESI1221466	ESI1221466	ESI1221466	ESI1221466
Total Petroleum Hydrocarbons		Total Dissolved Solids @180°C	10	mg/L														
		C6-C9 fraction	10	mg/L														
		C10-C14 fraction	50	mg/L														
		C15-C28 fraction	100	mg/L														
		C29-C34 fraction	50	mg/L														
Total Recoverable Hydrocarbons		C6-C10 fraction (sum)	20	mg/L														
		C6-C10 fraction (F minus BTEX)	20	mg/L														
		C6-C16 fraction	20	mg/L														
		C17-C34 fraction	100	mg/L														
		C35-C40 fraction (sum)	100	mg/L														
Monocyclic Aromatic Hydrocarbons		Benzene	1	mg/L	700													
		Toluene	1	mg/L														
		o-Xylene	1	mg/L														
		m-Xylene	2	mg/L														
		p-Xylene	2	mg/L														
		Total Xylenes	2	mg/L														
		1,2,4-trimethylbenzene	1	mg/L														
		Isopropylbenzene	1	mg/L														
		Styrene	1	mg/L														
		sec-butylbenzene	1	mg/L														
		n-Butylbenzene	1	mg/L														
		Sum of BTEX	1	mg/L														
		Total BTEX	1	mg/L														
		1,2,3-trichlorobenzene	1	mg/L														
Naphthalene		Naphthalene (VOC)	5	mg/L	70													
		Acenaphthylene	1	mg/L														
		Acenaphthene	1	mg/L														
		Fluorene	1	mg/L														
		Phenanthrene	1	mg/L														
Polynuclear Aromatic Hydrocarbons		Benzo[a]anthracene	1	mg/L														
		Benzo[b]fluoranthene	1	mg/L														
		Benzo[k]fluoranthene	1	mg/L														
		Benzo[a]pyrene	2	mg/L														
		Benzo[e]pyrene	0.5	mg/L														
		Chrysene	0.5	mg/L														
		Pyrene	1	mg/L														
		Benzo[a]anthracene	1	mg/L														
		Benzo[b]fluoranthene	1	mg/L														
		Benzo[k]fluoranthene	1	mg/L														
		Benzo[a]pyrene	2	mg/L														
		Benzo[e]pyrene	0.5	mg/L														
		Chrysene	0.5	mg/L														
		Pyrene	1	mg/L														
Metals		Sum of PAHs	0.5	mg/L														
		Arsenic (Dissolved)	0.001	mg/L														
		Cadmium	0.0001	mg/L														
		Cadmium (Dissolved)	0.0001	mg/L														
		Copper	0.001	mg/L														
		Copper (Dissolved)	0.001	mg/L														
		Lead	0.001	mg/L														
		Lead (Dissolved)	0.001	mg/L														
		Manganese	0.001	mg/L														
		Manganese (Dissolved)	0.001	mg/L														
		Mercury	0.0001	mg/L														
		Mercury (Dissolved)	0.0001	mg/L														
		Nickel	0.001	mg/L														
		Nickel (Dissolved)	0.001	mg/L														
Halogenated Aromatic Compounds		Zinc	0.005	mg/L														
		Zinc (Dissolved)	0.005	mg/L														
		Vanadium	0.01	mg/L														
		Vanadium (Dissolved)	0.01	mg/L														
		Chlorobenzene	1	mg/L														
		n-Butylbenzene	1	mg/L														
		Isopropylbenzene	1	mg/L														
		sec-Butylbenzene	1	mg/L														
		n-Butylbenzene	1	mg/L														
		1,2,3-Trichlorobenzene	1	mg/L														
		1,2,4-Trichlorobenzene	1	mg/L														
		1,3-Dichlorobenzene	1	mg/L														
		1,4-Dichlorobenzene	1	mg/L														
		2-Chlorotoluene	1	mg/L														
		Bromobenzene	1	mg/L														

Table 2
Analytical Results - September 2012 GME
Darwin Waterfront Corporation

Chem. Group	ChemName	output unit	EQL	ANZECC 2000 Marine water 95%												ANZECC 2000 Irrigation L v Y
				Total Dissolved Solids @180°C	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	Cr-O3 fraction	
Total Petroleum Hydrocarbons	Cr-O3 fraction	mg/L	10	1770	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	10	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	50	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
Total Recoverable Hydrocarbons	Cr-O3 fraction (sum)	mg/L	20	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction (sum)	mg/L	20	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction (sum)	mg/L	20	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction (sum)	mg/L	20	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction (sum)	mg/L	20	-	-	-	-	-	-	-	-	-	-	-	-	-
Monocyclic Aromatic Hydrocarbons	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
Naphthalene	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
Polynuclear Aromatic Hydrocarbons	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
Metals	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
Halogenated Aromatic Compounds	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-
	Cr-O3 fraction	mg/L	100	-	-	-	-	-	-	-	-	-	-	-	-	-

DATA VALIDATION REPORT _December 2012

URS Project number:	42213719	Data verified by:	Bek Aagaard	Date: 25/03/2013
Client:	Darwin Waterfront Corporation	Signed:		
Site:	Waterfront Precinct	Validation by:	Mitch Bacon	Date:
URS Project Manager:	Tim Smith	Signed:		3/4/2013
Matrix type:	Groundwater	Project Manager:	Tim Smith	Date:
No Primary samples:	10	Signed:		
Laboratory:	ALS; Envirolab			
Lab reference:	ES1229380, 83189			

Data quality objectives

Field data comparison Slight hydrocarbon odours were noted for MW127_11122012, but no Total Petroleum Hydrocarbons (TPHs) or benzene, toluene, ethylbenzene and xylenes (BTEX) were reported for MW127_11122012 in Batch ES1229380.

Frequency of field QC Field QC samples were collected to project specifications.

		Discrepancy	Frequency	Requested Frequency
Intra-laboratory Duplicates Taken	1	None noted	1	1
Inter-laboratory Duplicates Taken	1	None noted	1	1
Intra-laboratory Duplicates Analysed	1	None noted	1	1
Inter-laboratory Duplicates Analysed	1	None noted	1	1
Total Frequency			2	2

Inter-laboratory duplicates were not analysed at the required frequency for oxygenated compounds, carbon disulfide and halogenated aliphatic analytes. However, the presence of intra-laboratory duplicates and laboratory duplicates for these analytes is considered sufficient for interpretation of the precision of the results except for these analytes.

Frequency of laboratory QC Laboratory QC analysis was reported to project specifications, except for the frequency of laboratory duplicates for total metals. The rate was 8.1% and was supposed to be 10%.

Tests requested/reported Samples were analysed and reported as requested on the COC, except for MW135 was added twice on the COC.

Limits of reporting LORs were sufficiently low to enable assessment against adopted guideline criteria with the following limitations:

- Protocol LOR for anthracene (1 µg/L) above the adopted ILs (0.4 µg/L)
- Protocol LOR for benzo(a)pyrene (0.5 µg/L) above the adopted ILs (0.2 µg/L)

The limits of reporting were above the adopted investigation levels for anthracene and benzo(a)pyrene. As a result, potential exists for samples to contain concentrations of these analytes above the adopted investigation levels, but below detection limits. This lack of definitive data should be taken into consideration when interpreting analytical results below LOR for these PAHs.

- Protocol LOR exceeds the guideline trigger value for arsenic (0.1 mg/L)
- Protocol LOR exceeds the guideline trigger value for cadmium (0.0055 mg/L)
- Protocol LOR exceeds the guideline trigger value for chromium (0.0044 mg/L)
- Protocol LOR exceeds the guideline trigger value for copper (0.0013 mg/L)
- Protocol LOR exceeds the guideline trigger value for iron (0.2 mg/L)
- Protocol LOR exceeds the guideline trigger value for lead (0.0044 mg/L)
- Protocol LOR exceeds the guideline trigger value for nickel (0.07 mg/L)
- Protocol LOR exceeds the guideline trigger value for zinc (0.015 mg/L)

The limits of reporting exceeded the guideline trigger levels for several metals. As a result, potential exists for samples to contain concentrations of these analytes above the adopted investigation levels, but below detection limits. This lack of definitive data should be taken into consideration

	when interpreting analytical results below these metals.																		
Data transcription	A random 10% check of the laboratory results identified no anomalies within the electronic data, the laboratory reports, and tables generated by URS.																		
Sample management																			
Chain of Custody	Chain of custody documents completed.																		
Handling and preservation	Samples were correctly preserved, chilled with ice and received at the laboratory at 12.9°C (ice present). The recorded temperature for batch ES1229380 was slightly above the recommended temperature, potentially resulting in some loss of volatiles, and therefore possible under-reporting of associated analytes. This is not considered to affect the reported results, as concentrations of volatile analytes for the current and historical investigation have reported below the LOR and analytes were below the LOR in the trip blank (QCA_12122012).																		
Holding time compliance	Samples were extracted and analysed within recommended holding times.																		
Data precision																			
Field duplicate RPDs	RPDs were within control limits for Batch ES1229380.																		
Interlaboratory Replicate (Field triplicate) RPDs	RPDs were within control limits for Batch ES1229380.																		
Laboratory duplicate RPDs	RPDs were within control limits for Batch ES1229380.																		
Data accuracy																			
Laboratory control spike recovery	Laboratory control spike recoveries were within control limits.																		
Matrix spike recovery	The following recoveries were outside control limits and may affect data interpretation: <table><thead><tr><th>Sample</th><th>Analyte</th><th>Recovery (%)</th><th>LCL (%)</th><th>UCL (%)</th><th>Comment</th></tr></thead><tbody><tr><td>MW136_11122012</td><td>Manganese</td><td>ND</td><td>-</td><td>-</td><td>MS recovery not determined, background level greater than or equal to 4x spike level.</td></tr><tr><td colspan="6">ND – Not determined</td></tr></tbody></table>	Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment	MW136_11122012	Manganese	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.	ND – Not determined					
Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment														
MW136_11122012	Manganese	ND	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.														
ND – Not determined																			
	Matrix spike recovery was not determined for manganese due to elevated sample concentrations. . The accuracy of the data for manganese is considered acceptable based on the presence of other quality control data, such as method blanks, LCS recoveries and matrix spike recoveries for other metals analysed under the same analytical method.																		
	Matrix spike recoveries were not reported for PAHs, calcium, magnesium, potassium, sodium, halogenated methanes, oxygenated compounds, monocyclic aromatic hydrocarbons, carbon disulphide, halogenated aliphatics and alkalinity. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as method blanks, LCS recoveries and surrogate recoveries (where applicable).																		
Surrogate spike recovery	Surrogate recoveries were within control limits for Batch ES1229380.																		
Blank monitoring																			
Equipment rinsate blank	Concentrations of all analytes were reported below the LOR.																		
Field blank	Concentrations of all analytes were reported below the LOR.																		
Trip blank	Concentrations of all analytes were reported below the LOR.																		
Method blank	Concentrations of all analytes were reported below the LOR.																		
	Method blank recoveries were not reported for alkalinity. The accuracy of alkalinity data could not be assessed. This should be taken into consideration when interpreting the accuracy of the results for this analyte.																		
Other observations	None reported.																		

Site Name
Waterfront Precinct
Project No.
42213719
Project Manager
Tim Smith
Matrix
Water
ALS / EnviroLab
Laboratory
Batch File Number
ESI1239380

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Holding Times (a)	Limits of Reporting (b)	Field Blank (1 per day)		Rinsate Blank (1 per day)		Trip Blank (1 per day with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
						Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Reported	OK
METALS ANALYSIS (EPA 8230F)	Arsenic	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	-	-
	Calcium	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
	Chromium	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
	Copper	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
	Lead	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
METALS ANALYSIS (EPA 8210)	Magnesium	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
	Sodium	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
	Vanadium	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
	Zinc	12	12	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	2	5	1	3	1	3	-	-
	Mercury	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-
VOLATILES ANALYSIS (EPA 8210)	Benzene	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	Ethylbenzene	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	m-Xylene & p-Xylene	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	Naphthalene	12	12	✓	✓	2	2	2	2	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	ortho-Xylene	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS (EPA 8210)	Toluene	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	o-C8: C10 Fraction	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	o-C8: C10 Fraction: naph BTEX (F1)	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	o-C8: C10 Fraction	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	o-C8: C10 Fraction	12	12	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
SEMI-VOLATILES ANALYSIS (EPA 8210)	Acenaphthene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Acenaphthylene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Anthracene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[a]anthracene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[b]fluoranthene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
SEMI-VOLATILES ANALYSIS (EPA 8210)	Benzo[k]fluoranthene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[e]pyrene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[a]pyrene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[a]anthracene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[b]fluoranthene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
SEMI-VOLATILES ANALYSIS (EPA 8210)	Benzo[a]pyrene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[a]anthracene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[b]fluoranthene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[e]pyrene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	Benzo[a]pyrene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
SEMI-VOLATILES ANALYSIS (EPA 8210)	Pyrene	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	o-C8: C10 Fraction	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	o-C8: C10 Fraction	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	o-C8: C10 Fraction	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
	o-C8: C10 Fraction	12	12	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	1	1	2	✓	✓
METALS ANALYSIS (EPA 8230F)	Calcium	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	-	-
	Magnesium	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	-	-
	Potassium	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	-	-
	Sodium	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	-	-
	Boron	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	-	-
VOLATILES ANALYSIS (EPA 8210)	Benzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	Chloroform	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	Bromochloroform	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	Bromobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	2-Bromobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS (EPA 8210)	2-Heptanol (MBO)	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	4-Methyl-2-pentanol (MBO)	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	2-Heptanol (MBO)	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	4-Methyl-2-pentanol (MBO)	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	2-Heptanol (MBO)	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS (EPA 8210)	1,2,4-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS (EPA 8210)	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS (EPA 8210)	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓
	1,3,5-Trinitrobenzene	9	9	✓	✓	0	0	0	0	0	0	1	1</												

Site Name
Waterfront Precinct
Project No.
42213719
Project Manager
Tim Smith
Matrix
Water
ALS / EnviroLab
Laboratory
Batch File Number
ESI229380

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Performed	Primary Samples	Holding Times (h)	Limits of Reporting (g)	Field Blank (1 per day)		Rinsate Blank (1 per day)		Trip Blank (1 per easy with VOCs)		Method Blank (1 per batch)		Intra-Laboratory Duplicate Sample (1 in 20)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (GC-MS organics)	
							Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Reported	OK
VOLATILES ANALYSIS (ALSEP/42)	1,1,1,2-Tetra chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,1,1 Trichloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,1,2,2-Tetrachloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,1,2 Trichloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dichloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dichloroethene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2,3-Trichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dibromo-3-chloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dichloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/42)	1,2-Dichloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Bromonaphthalene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Carbon Tetrachloride	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/42)	1,2,3-Trichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2,4-Trichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dichlorobenzene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,3-Dichlorobenzene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,4-Dichlorobenzene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2-Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	4-Chloroethane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Bromobenzene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chlorobenzene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Chlorobenzene	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
VOLATILES ANALYSIS (ALSEP/42)	Carbon disulfide	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dibromoethane (DOB)	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	1,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	2,2-Dichloropropane	9	9	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
INSTRUMENT LABORATORY (NON-METALS)	1,2-Dichloropropane	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbonate Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbonate Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Hydroxide Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Total Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Total Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Total Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Total Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Total Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Total Alkalinity as CaCO3	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
INSTRUMENT LABORATORY (NON-METALS)	Carbon disulfide	12	12	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	1	2	2	1	0	1	0	-	-
	Carbon disulfide	12	12	10	✓	✓</																				

Table 3
Analytical Results - December 2012 GME
Darwin Waterfront Corporation

[illegible]

*LOR exceeds Adopted Guidelines

Chem Group		ChemName	EQL	Output unit	ANZECC 2000 Marine water 95%	ANZECC 2000 Irrigation LV	Location Code	MM18	MM18	QC001	QC001	QC001
Total Petroleum Hydrocarbons		C6-C9 fraction	10	µg/L		<20						
Total Recoverable Hydrocarbons		C10-C14 fraction	50	µg/L		<50						
		C15-C28 fraction	100	µg/L		<100						
		C29-C36 fraction	50	µg/L		<50						
		C37-C40 fraction (sum)	10	µg/L		<10						
		C6-C10 fraction (sum)	10	µg/L		<20						
		C6-C10 fraction (sum)	10	µg/L		<20						
		C6-C10 fraction (sum)	10	µg/L		<20						
		C6-C10 fraction (sum)	10	µg/L		<20						
		C6-C10 fraction (sum)	10	µg/L		<20						
		C6-C10 fraction (sum)	10	µg/L		<20						
Monocyclic Aromatic Hydrocarbons		<C10-C16 fraction	50	µg/L		<100						
		<C10-C16 (less Naphthalene)	50	µg/L		<100						
		<C16-C34 fraction	100	µg/L		<100						
		<C34-C40 fraction	100	µg/L		<100						
		<C10-C40 fraction (sum)	100	µg/L		<100						
		Benzene	1	µg/L		<1						
		Toluene	1	µg/L		<1						
		o-Xylene	1	µg/L		<1						
		m-Xylene	2	µg/L		<2						
		p-Xylene	2	µg/L		<2						
Naphthalene		Total Xylenes	2	µg/L		<2						
		1,2,4-trimethylbenzene	1	µg/L		<1						
		Isopropylbenzene	1	µg/L		<1						
		Styrene	1	µg/L		<1						
		sec-butylbenzene	1	µg/L		<1						
		1,3,5-trimethylbenzene	1	µg/L		<1						
		1,3,5-trimethylbenzene	1	µg/L		<1						
		1,3,5-trimethylbenzene	1	µg/L		<1						
		1,3,5-trimethylbenzene	1	µg/L		<1						
		1,3,5-trimethylbenzene	1	µg/L		<1						
Polynuclear Aromatic Hydrocarbons		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
		Acenaphthene	1	µg/L		<1						
Metals		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
Metals		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						
		Asenic (Filtered)	0.01	mg/L		<0.01						

