

DATA VALIDATION REPORT_GME March 2013

URS Project number: 42213719 Data verified by: Bek Aagaard Date: 13/09/2013
 Client: Darwin Waterfront Corporation
 Site: Waterfront Precinct Signed: 
 URS Project Manager: Jacques van Rensburg
 Matrix type: Water Validation by: Mitchell Bacon Date: 11/11/2013
 Signed: 
 No Primary samples: 5
 Laboratory: ALS / Envirolab
 Lab reference: ES1304969 / 86813

Data quality objectives

Field data comparison

Hydrocarbon odours were noted for MW125_04032013 and slight hydrocarbon odour for MW128_04032013; however, no Total Petroleum Hydrocarbons (TPHs), or benzene, toluene, ethylbenzene and xylenes (BTEX) were reported for MW125_04032013 and MW128_04032013. As TPH and BTEX has not been reported historically in MW125_04032013 and TPH has not been reported since December 2011 (low concentrations), this is not considered to affect the interpretation of the results.

Frequency of field QC

Field QC samples were collected to project specifications with the following exceptions:

		Discrepancy	Frequency	Required 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			1	1

- Intra- and inter-laboratory duplicates were not reported for PAH and VOCs. The precision of VOC results has been considered acceptable based on the presence of laboratory duplicates were available and acceptable. The precision of PAH cannot be assessed as no laboratory duplicates were available and care should be taken when using PAH results for quantitative purposes.

Frequency of laboratory QC

The laboratory reported a sufficient frequency of QC to assess whether the results have been reported to an acceptable accuracy and precision, with the following exceptions:

- Laboratory duplicates were not reported for PAH. The precision of PAH cannot be assessed as no intra- and inter-laboratory duplicates were available and care should be taken when using PAH results for quantitative purposes.

Tests requested/reported

Samples were analysed and reported as requested on the COC for all Batches.

Limits of reporting

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

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

The limits of reporting was above the adopted investigation levels for benzo(a)pyrene. As a result, potential exists for samples to contain concentrations of benzo(a)pyrene 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 this analyte.

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

Samples were analysed and reported as requested on the COC with the following exception:

According to the SRN field filtered metals bottle for MW103 was not supplied to the laboratory; consequently, the 'Clear Plastic Bottle – Natural' was subsampled, laboratory filtered and analysed for dissolved metals and mercury. Hence, due to not being acidified or field filtered, reported metal



concentrations may not accurately represent field concentrations and this should be taken into consideration when interpreting metals results for these samples. The reported metals concentrations are consistent with historical results for the affected monitoring wells and the results for these analytes are assessed as acceptable for interpretation.

Handling and preservation

Samples were transported with ice bricks and were received at the laboratory at 1.6°C.

Holding time compliance

Samples were extracted and analysed within recommended holding times.

Data precision

Field duplicate RPDs

Field duplicate RPDs were within control limits.

Field triplicate RPDs

Field triplicate RPDs were within control limits for batch ES1304969.

Laboratory duplicate RPDs

Laboratory duplicate RPDs were within control limits for batch ES1304969.

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:

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
Anonymous	Manganese	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Anonymous	Zinc	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW128_040313	Ferrous Iron	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW128_040313	Methane	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

The matrix spikes for manganese, zinc, ferrous iron and methane were not determined due to background levels being greater than the spike level. This does not reflect method bias and the accuracy of the results for these analytes are assessed as acceptable due to the presence of other laboratory quality control data including method blanks, LCS recoveries and surrogate recoveries for analytes analysed under the same analytical method.

Matrix spike recoveries were not reported for sulphur as S, pyrene and acenaphthene. 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 ES1304969.

Blank monitoring

Rinsate blank

Concentrations were not detected above the LOR for all analytes tested, with the following exception:

- Concentrations of nickel (0.001 mg/L) were detected above the LOR in the rinsate sample collected on 4th of March 2013 (QCCF01_040313) in batch ES1304969. The nickel concentration (0.001 mg/L) in the rinsate blank was reported at equal or below concentrations as in the primary samples. As the nickel concentrations in samples collected on this sampling day were either below the GAC by more than the concentration reported in the rinsate, or well above, and as no other metals analytes were reported above LOR cross-contamination from inadequate decontamination is not considered to affect the interpretation of the results.

Field blank

Concentrations of all analytes were reported below the LOR.

Trip blank

Concentrations of volatile TPH and TRH were reported below the LOR.

Method blank

Concentrations of all analytes were reported below the LOR.

Other observations

Batch ES1304969

- EA015 TDS may bias high for sample ID MW122 and QAQC01 due to the presence of fine particulate matter, which have passed through the prescribed GF/C paper.

DATA VALIDATION REPORT GME March 2013

URS Project number: 42213719 Data verified by: Bek Aagaard Date: 13/09/2013
 Client: Darwin Waterfront Corporation
 Site: Waterfront Precinct Signed: *Bek Aagaard*
 URS Project Manager: Jacques van Rensburg Validation by: Mitchell Bacon Date: 11/11/2013
 Matrix type: Water Signed: *[Signature]*
 No Primary samples: 8
 Laboratory: ALS
 Envirolab
 Lab reference: ES1305174 / 86859

Data quality objectives

Field data comparison

No apparent anomalies were observed between laboratory results and field observations.

Frequency of field QC

Field QC samples were collected to project specifications with the following exceptions:

		Discrepancy	Frequency	Required 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

Intra- and inter-laboratory duplicates were not reported for PAH and VOCs. The precision of VOC results has been considered acceptable based on the presence of laboratory duplicates were available and acceptable. The precision of PAH cannot be assessed as no laboratory duplicates were available and care should be taken when using PAH results for quantitative purposes.

Frequency of laboratory QC

The laboratory reported a sufficient frequency of QC to assess whether the results have been reported to an acceptable accuracy and precision, with the following exceptions:

- Laboratory duplicate were not reported for PAH, TPH C₁₀-C₃₆ and TRH >C₁₀-C₄₀. The precision of TPH C₁₀-C₃₆ and TRH >C₁₀-C₄₀ results has been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable. The precision of PAH cannot be assessed as no intra- and inter-laboratory duplicates were available and care should be taken when using PAH results for quantitative purposes.
- Laboratory QC analysis was reported to project specifications, except for the frequency of laboratory duplicates for dissolved mercury (9.7% actual and 10% required). The precision of dissolved mercury results has been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable, which reported RPDs within acceptable levels.

Tests requested/reported

Samples were analysed and reported as requested on the COC.

Limits of reporting

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

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

The limits of reporting was above the adopted investigation levels for benzo(a)pyrene. As a result, potential exists for samples to contain concentrations of benzo(a)pyrene 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 this analyte.

- Elevated LOR (after dilution) exceeds the GAC for cadmium (0.0002 mg/L)
- Elevated LOR (after dilution) exceeds the GAC for chromium (0.0044 mg/L)
- Elevated LOR (after dilution) exceeds the GAC for copper (0.0013 mg/L)
- Elevated LOR (after dilution) exceeds the GAC for lead (0.0034 mg/L)
- Elevated LOR (after dilution) exceeds the GAC for zinc (0.008 mg/L)

The limits of reporting (after dilution) exceeded the GAC for cadmium, chromium, copper, lead and zinc. 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 for 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

Samples were analysed and reported as requested on the COC, with the following exception:

- The field filtered metals bottles for MW108, MW116, MW129 and MW136 were received as "unspecified" - that is, the filtered box on the bottle was not ticked. Consequently, the 'Clear Plastic Bottle - Natural' was subsampled, laboratory filtered and analysed for dissolved metals and mercury.

Handling and preservation

Samples were transported with ice bricks and were received at the laboratory at 17.3°C. Samples were received above the recommended temperature range for water samples on receipt (4°C, ± 2°C); therefore, some losses through volatilisation may have occurred and sample concentrations for BTEXN and TPH C₆-C₉ (and other volatiles) may be biased low.

It should be noted that the ambient temperature across the site throughout the year and at the time of sampling is above 30°C. Samples are chilled in the field on ice and transported to the laboratory with appropriate cooling medium. Samples remain chilled while cooling medium is present; however, over the 24 h transit period to the laboratory cooling medium will melt and samples may warm. Due to the initial sample temperature and subsequent appropriate cooling of samples, losses due to volatilisation and sample degradation are considered to be limited.

Sample containers for dissolved metals and dissolved mercury did not comply with preservation standards for MW108, MW116, MW129, MW136 and QAQC03. SRN states, "A 'Clear Plastic Bottle - Natural' was provided, and the preferred sample container was 'Clear Plastic Bottle - Nitric Acid; Filtered' ". The correct sample bottles were provided to the lab, however, the "field filtered" box was not ticked, and the laboratory did not confirm the status of the sample. Consequently, the 'Clear Plastic Bottle - Natural' was subsampled, laboratory filtered and analysed for dissolved metals and mercury. Hence, due to not being acidified or field filtered, reported metal concentrations may not accurately represent field concentrations and this should be taken into consideration when interpreting metals results for these samples. The reported metals concentrations are consistent with historical results for the affected monitoring wells and the results for these analytes are assessed as acceptable for interpretation.

Holding time compliance

Samples were extracted and analysed within recommended holding times.

Data precision

Field duplicate RPDs

Field duplicate RPDs were within control limits.

Field triplicate RPDs

RPDs exceeded control limits for the following sample analysis in batch ES1305174 (Samples with higher reported concentrations are in bold):

- **MW129_050313 and QAQC04_050313** for zinc (82%).

The potential imprecision in zinc concentrations should be taken into consideration when interpreting results close to the ILs; hence, care should be taken when using the data for zinc for any quantitative purposes.

Laboratory duplicate RPDs

Laboratory duplicate RPDs were within control limits.

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:

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
MW115_060312	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Anonymous	Ferrous Iron	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

The matrix spike for chloride and ferrous iron were not determined due to background levels being greater than the spike level. This does not reflect method bias and the accuracy of the results for these analytes are assessed as acceptable due to the presence of other laboratory quality control data including method blanks, LCS recoveries and surrogate recoveries for analytes analysed under the same analytical method.

Matrix spike (MS) recovery was reported on an anonymous sample for ferrous iron. Where MS were conducted on an anonymous sample, the reported MS recovery may not be considered representative of



the URS field sample.

Matrix spike recoveries were not reported for sulphur as S, pyrene, acenaphthene and TPH C₁₀-C₃₆ and TRH >C₁₀-C₄₀. 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
MW134_050313	4-Bromofluorobenzene	79.9	80.8	123.7	Recovery less than lower data quality objective

Surrogate recoveries were reported lower than the lower data quality objective for 4-Bromofluorobenzene; hence, there is the potential for the volatile TPH C6-C10, benzene, toluene, ethylbenzene and xylenes (BTEX) and volatile organic carbon (VOC) results to be biased low. This is not considered to affect the interpretation of the results, as volatile TPH C6-C10, BTEX and VOCs were not reported above the laboratory LOR in this sample, and the magnitude of methane reported is assessed as acceptable for interpretation.

Blank monitoring

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 volatile TPH and TRH were reported below the LOR.

Method blank

Concentrations of all analytes were reported below the LOR.

Other observations

Batch ES1305179

- EA015 TDS may bias high for sample ID MW122 and QAQC01 due to the presence of fine particulate matter, which may pass through the prescribed GF/C paper.
- EG020: Some samples were diluted and rerun due to matrix interference and LORs have been raised accordingly (High salinity).



DATA VALIDATION REPORT_GME March 2013

URS Project number: 42213719 Data verified by: Bek Aagaard Date: 13/09/2013
 Client: Darwin Waterfront Corporation
 Site: Waterfront Precinct Signed: Uebeke Aagaard
 URS Project Manager: Jacques van Rensburg
 Matrix type: Water Validation by: Mitchell Bacon Date: 11/11/2013
 No Primary samples: 5 Signed: 
 Laboratory: ALS
 Lab reference: ES1305223

Data quality objectives

Field data comparison No apparent anomalies were observed between laboratory results and field observations.

Frequency of field QC Field QC samples were collected to project specifications with the following exceptions:

- It was not specified on the COC to send the inter-laboratory sample QAQC_06032013 to a secondary lab. Consequently, this sample was analysed at the primary laboratory as an intra-laboratory duplicate. Field QC samples were analysed to project specifications with a total frequency of four intra-laboratory duplicates and two inter-laboratory duplicates for the total of eighteen primary samples for batches ES1304969, ES1305174 and ES1305223.
- Intra- and inter-laboratory duplicates were not reported for PAH and VOCs. The precision of VOC results has been considered acceptable based on the presence of laboratory duplicates were available and acceptable. The precision of PAH cannot be assessed as no laboratory duplicates were available and care should be taken when using PAH results for quantitative purposes.

		Discrepancy	Frequency	Required Frequency
Intra-laboratory Duplicates Taken	1	None noted	1	1
Inter-laboratory Duplicates Taken	1	None noted	1	1
Intra-laboratory Duplicates Analysed	2	None noted	2	1
Inter-laboratory Duplicates Analysed	0	No inter-laboratory duplicate analysed	0	1
Total Frequency			2	2

Frequency of laboratory QC The laboratory reported a sufficient frequency of QC to assess whether the results have been reported to an acceptable accuracy and precision, with the following exceptions:

- Laboratory duplicate were not reported for PAH, TPH C₁₀-C₃₆ and TRH >C₁₀-C₄₀. The precision of TPH C₁₀-C₃₆ and TRH >C₁₀-C₄₀ results has been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable. The precision of PAH cannot be assessed as no intra- and inter-laboratory duplicates were available and care should be taken when using PAH results for quantitative purposes.
- Laboratory QC analysis was reported to project specifications, except for the frequency of laboratory duplicates for dissolved major ions (9.1% actual and 10% required). The precision of major ions results has been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable, which reported RPDs within acceptable levels.

Tests requested/reported Samples were analysed and reported as requested on the COC.

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

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

The limits of reporting was above the adopted investigation levels for benzo(a)pyrene. As a result, potential exists for samples to contain concentrations of benzo(a)pyrene 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 this analyte.

- Elevated LOR (after dilution) exceeds the GAC for cadmium (0.0002 mg/L)



	<ul style="list-style-type: none"> Elevated LOR (after dilution) exceeds the GAC for chromium (0.0044 mg/L) Elevated LOR (after dilution) exceeds the GAC for copper (0.0013 mg/L) Elevated LOR (after dilution) exceeds the GAC for lead (0.0034 mg/L) Elevated LOR (after dilution) exceeds the GAC for zinc (0.008 mg/L) <p>The limits of reporting (after dilution) exceeded the GAC for cadmium, chromium, copper, lead and zinc in MW10. 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 for these metals.</p>
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	Samples were analysed and reported as requested on the COC. It was not specified on the COC to send the inter-laboratory sample QAQC_06032013 to a secondary lab; consequently, this sample was analysed at the primary laboratory as an intra-laboratory duplicate. Field QC samples were analysed to project specifications with a total frequency of four intra-laboratory duplicates and two inter-laboratory duplicates for the total of eighteen primary samples for batches ES1304969, ES1305174 and ES1305223.
Handling and preservation	<p>Samples were transported with ice bricks and were received at the laboratory at 6.6°C. Samples were received slightly above the recommended temperature range for water samples on receipt (4°C, ± 2°C); therefore, some losses through volatilisation may have occurred and sample concentrations for BTEXN and TPH C₆–C₉ (and other volatiles) may be biased low.</p> <p>It should be noted that the ambient temperature across the site throughout the year and at the time of sampling is above 30°C. Samples are chilled in the field on ice and transported to the laboratory with appropriate cooling medium. Samples remain chilled while cooling medium is present; however, over the 24 h transit period to the laboratory cooling medium will melt and samples may warm. Due to the initial sample temperature and subsequent appropriate cooling of samples, losses due to volatilisation and sample degradation are considered to be limited.</p>
Holding time compliance	<p>Samples were extracted and analysed within recommended holding times, with the following exceptions:</p> <ul style="list-style-type: none"> Nitrate as N and sulphite as SO₃²⁻ analysis was exceeded by 4 days for all samples. These analytes are used as indicators of natural attenuation and as such will not affect the overall interpretation of results; however this should be taken into consideration if the results are used for quantitative purposes.
Data precision	
Field duplicate RPDs	Field duplicate RPDs were within control limits.
Field triplicate RPDs	Field Triplicate RPDs were within control limits.
Laboratory duplicate RPDs	Laboratory duplicate RPDs were within control limits.
Data accuracy	
Laboratory control spike recovery	Laboratory control spike recoveries were within control limits.
Matrix spike recovery	<p>Matrix spike recoveries were within control limits.</p> <p>Matrix spike recoveries were not reported for sulphur as S, pyrene, acenaphthene and TPH C₁₀–C₃₆ and TRH >C₁₀–C₄₀. 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).</p>
Surrogate spike recovery	Surrogate recoveries were within control limits.
Blank monitoring	
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 volatile TPH and TRH were reported below the LOR.
Method blank	Concentrations of all analytes were reported below the LOR.
Other observations	<p>Batch ES1305223</p> <ul style="list-style-type: none"> EG020: Some samples (MW10_060313) were diluted and re-run due to matrix interference and LORs have been raised accordingly (High Total Dissolved Solids).

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Number of Primary Samples	Holding Times (h)	Limits of Reporting (lb)	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	Reported	OK
METALS ANALYSES/ED00F	Arsenic	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Calcium	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Chromium	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Copper	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Lead	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Manganese	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Nickel	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Vanadium	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	5	1	3	1	3	-	-
	Zinc	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	1	5	1	3	1	3	-	-
	Zirconium	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
VOLATILES ANALYSES/EP00B	Benzene	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	Benzophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	Diethyl Benzene	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	Naphthalene	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	ortho-Xylene	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	para-Xylene	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	Toluene	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	Volatiles ANALYSES/EP00B/071	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	Volatiles ANALYSES/EP00B/071	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
	Volatiles ANALYSES/EP00B/071	9	9	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	3	1	2	1	2	✓	✓
SEMIVOLATILES ANALYSES/EP00B	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
VOLATILES ANALYSES/EP00B	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
VOLATILES ANALYSES/EP00B	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
VOLATILES ANALYSES/EP00B	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
VOLATILES ANALYSES/EP00B	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	0	1	0	1	0	1	1	✓	✓
	Acetophenone	1	1	1	✓	✓	0																			

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Holding Times (h)	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 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	Reported	OK						
METALS ANALYSES/EP002F	Arsenic	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Calcium	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Chromium	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Copper	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Lead	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Manganese	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Nickel	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Vanadium	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Zinc	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	-	-				
	Zinc	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
VOLATILES ANALYSES/EP008	Benzene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Benzene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Benzene & para-Xylene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Naphthalene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	ortho-Xylene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	ortho-Xylene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Toluene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Volatiles ANALYSES/EP008/071	10	10	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Volatiles ANALYSES/EP008/071	10	10	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Volatiles ANALYSES/EP008/071	10	10	8	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
SEMI-VOLATILES ANALYSES/EP009/071	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	2	1	1	1	1	✓	✓				
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
SEMI-VOLATILES ANALYSES/EP009/071	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
VOLATILES ANALYSES/EP008	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓			
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
VOLATILES ANALYSES/EP008	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓	
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓	
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓	
VOLATILES ANALYSES/EP008	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓		
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓	
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓	
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
VOLATILES ANALYSES/EP008	Acenaphthene	5	5	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1</										

[illegible]

NOTES:

(a)	✓	holding times are within project guideline limits.
	✗	holding times exceed project guideline limits.
(b)	✓	Limits of reporting (LORs) comply with project specifications.
	✗	LORs do not comply with project specifications.

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Primary Samples	Holding Time (a)	Limits of Reporting (b)	Field Blank (1 per day)		Rinseate 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/ED20P	Arsenic	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	0	1	2	1	1	1	1	1	-
	Cadmium	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	0	1	2	1	1	1	1	-
	Chromium	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	0	1	2	1	1	1	1	-
	Copper	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	0	1	2	1	1	1	1	-
	Lead	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	0	1	2	1	1	1	1	-
	Manganese	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	0	1	2	1	1	1	1	-
	Nickel	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	0	1	2	1	1	1	1	-
	Vanadium	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	0	1	2	1	1	1	-
	Zinc	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	0	1	2	1	1	1	-
	Zn	7	7	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	0	1	2	1	1	1	-
NON-METALS ANALYSIS/ALS/EP008	Boron	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Chloride	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Fluoride & Sulfate	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Nitrate & Sulfate	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Nitrite	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Phosphate	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Ammonia	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Ammonium	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Calcium	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
	Iron	9	9	5	✓	✓	1	1	1	1	0	0	1	3	1	1	1	1	0	1	2	1	1	2	✓
NON-METALS ANALYSIS/ALS/EP007H	Fluoride	10	10	5	✓	✓	1	1	1	1	1	1	1	2	1	1	1	1	0	1	3	1	1	2	✓
	Chloride	10	10	5	✓	✓	1	1	1	1	1	1	1	2	1	1	1	1	0	1	3	1	1	2	✓
	Chloride Fraction	10	10	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	0	1	3	1	1	2	✓
	Chloride Fraction	10	10	5	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	0	1	3	1	1	2	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	✓
NON-METALS ANALYSIS/ALS/EP007S/MSB	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	✓
	Acetophenone	1	1	1	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	✓
NON-METALS ANALYSIS/ALS/ED20P	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	9	9	5	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
NON-METALS ANALYSIS/ALS/ED20P	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	-
	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	-
	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	1	1	1	1	1	-
	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
	Acetophenone	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	0	1	1	1	1	1	-
INSTRUMENT LABORATORY (NON-METALS)/ALS/EP007G	Nitrate as N	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	2	1	1	1	1	-
	Sulfate as SO4 ²⁻ (Dichromate)	7	7	5	✓	✓	0	0	0	0	0	0	1	1	1	1	1	0	1	2	1	1	1	1	-

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Holding Time (h)	Limits of Reporting (b)	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)		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		
VOLATILES ANALYSISALS-EPIV43	Bromochloromethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Bromodichloromethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Chloroform	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Dibromochloromethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV48	2-Bromoethanol (MBC)	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	2-Propanol (MBC)	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	4-Methyl-2-pentanol (MBC)	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	WVO Acetate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV44	1,2,4-Trinitrophenol	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	1,3,5-Trinitrophenol	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Isopropylbenzene	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	tert-Butylbenzene	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV46	tert-Butylbenzene	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	1,1,1-Trichloroethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	1,1,2-Trichloroethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	1,2-Dibromochloropropane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV45	1,2-Dichloroethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	1,3-Dichloropropane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Bromonitroethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Carbon Tetrachloride	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV47	Chloroethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Chloromethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	cis-1,2-Dichloroethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	cis-1,4-Dichloro-2-butene	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV42	Dibromomethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Dichlorodifluoromethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Hexachlorobutadiene	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Iodomethane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV41	Perfluorooctane	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Tetraethylenesulfone	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV40	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV49	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV47	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV46	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV44	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV43	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV42	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV41	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV40	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV39	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
VOLATILES ANALYSISALS-EPIV38	Triethylamine 2-chloride sulfate	1	1	✓	✓	0	0	0	0	0	0	1	1	1	0	1	1	0	0	1	✓
	Triethylamine 2-chloride sulfate	1																			

DATA VALIDATION REPORT_GME June 2013

URS Project number:	42213719	Data verified by:	Bek Aagaard	Date:	13/09/2013
Client:	Darwin Waterfront Corporation				
Site:	Waterfront Precinct	Signed:			
URS Project Manager:	Jacques van Rensburg	Validation by:	Mitchell Bacon	Date:	11/11/2013
Matrix type:	Water	Signed:			
No Primary samples:	10				
Laboratory:	ALS / Labmark				
Lab reference:	ES1314375 /384097				
	ES1314611	Project Manager:	Jacques van Rensburg		

Data quality objectives

Field data comparison No apparent anomalies were observed between laboratory results and field observations (batches ES1314375 and ES1314611).

Frequency of field QC Field QC samples were collected at a higher frequency than specified in project specifications.

		Discrepancy	Frequency	Required Frequency
Intra-laboratory Duplicates Taken	2	None noted	2	1
Inter-laboratory Duplicates Taken	2	None noted	1	1
Intra-laboratory Duplicates Analysed	2	None noted	2	1
Inter-laboratory Duplicates Analysed	2	None noted	1	1
Total Frequency			3	2

Frequency of laboratory QC The laboratory reported a sufficient frequency of QC to assess whether the results have been reported to an acceptable accuracy and precision, with the exception of the following:

- Inter-laboratory duplicates were not analysed at the required frequency for bicarbonate alkalinity as CaCO_3 and carbonate alkalinity as CaCO_3 . However, the presence of intra-laboratory duplicates and laboratory duplicates for these analytes at the required frequency, with acceptable RPDs, and appropriate laboratory QC. The precision of data is considered sufficient for interpretation of the results except for bicarbonate alkalinity as CaCO_3 and carbonate alkalinity as CaCO_3 .
- Laboratory duplicates were not reported for TPH $\text{C}_{10}\text{-C}_{36}$ and TRH $\text{C}_{10}\text{-C}_{40}$ for batch ES1314375. The precision of TPH and TPH results has been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable.

Tests requested/reported Samples were analysed and reported as requested on the COC for batches ES1314375 and ES1314611.

Limits of reporting LORs were sufficiently low to enable assessment against adopted guideline criteria.

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 Samples were analysed and reported as requested on the COC with the following exception:

According to the SRN field filtered metals bottles for QCB_260613 (field blank) and QCC_260613 (rinsate blank) were not supplied to the laboratory; consequently, the 'Clear Plastic Bottle - Natural' was subsampled, laboratory filtered and analysed for dissolved metals and mercury. Hence, due to not being acidified or field filtered, reported metal concentrations may not accurately represent field concentrations, and this should be taken into consideration when interpreting metals results for these samples. The reported metals concentrations are consistent with historical results for the affected monitoring wells and the results for these analytes are assessed as acceptable for interpretation.

Handling and preservation Samples were transported with ice bricks and were received at 0.3°C and 0.2°C in batches ES1314375 and ES1314611, respectively. Samples from batches ES1314375 and ES1314611 were received below the recommended temperature range for water samples on receipt (4°C, $\pm 2^\circ\text{C}$).



Holding time compliance	Samples were extracted and analysed within recommended holding times, with the following exceptions: <ul style="list-style-type: none">Dissolved metals analysis was exceeded by 5 days for samples QCB01_250613 (field blank) and QCB01_250613 (rinsate blank) in batch ES1314374. As all results for these analytes were reported below LOR in the field blank and rinsate blank, there is no evidence of cross-contamination and the potential underreporting of dissolved metals in these samples is not considered to affect the interpretation of the results.																								
Data precision																									
Field duplicate RPDs	Field duplicate RPDs were within control limits for batches ES1314375 and ES1314611.																								
Field triplicate RPDs	Field triplicate RPDs were within control limits for batches ES1314375 and ES1314611.																								
Laboratory duplicate RPDs	Laboratory duplicate RPDs were within control limits for batches ES1314375 and ES1314611.																								
Data accuracy																									
Laboratory control spike recovery	Laboratory control spike recoveries were within control limits for batches ES1314375 and ES1314611.																								
Matrix spike recovery	The following recoveries were outside control limits and may affect data interpretation: <table><tr><th>Sample</th><th>Analyte</th><th>Recovery (%)</th><th>LCL (%)</th><th>UCL (%)</th><th>Comment</th></tr><tr><td>Anonymous</td><td>Chloride</td><td>Not determined</td><td>-</td><td>-</td><td>MS recovery not determined, background level greater than or equal to 4x spike level.</td></tr><tr><td>Anonymous</td><td>TPH C₁₀-C₁₄ fraction</td><td>Not determined</td><td>-</td><td>-</td><td>MS recovery not determined, background level greater than or equal to 4x spike level.</td></tr><tr><td>Anonymous</td><td>TRH >C₁₀-C₁₆ fraction</td><td>Not determined</td><td>-</td><td>-</td><td>MS recovery not determined, background level greater than or equal to 4x spike level.</td></tr></table> <p>The matrix spike for chloride (batch ES1314375) and TPH C₁₀-C₁₄ fraction and TRH >C₁₀-C₁₆ fraction (batch ES1314611) were not determined due to background levels being greater than the spike level. As this does not reflect method bias, the accuracy of the results for these analytes are assessed as acceptable due to the presence of other laboratory quality control data, including method blanks, LCS recoveries and surrogate recoveries for analytes analysed under the same analytical method.</p> <p>Matrix spike recoveries were not reported for TPH C₁₀-C₁₄ and TRH >C₁₀-C₁₆ fractions in batch ES1314375. 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).</p>	Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment	Anonymous	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.	Anonymous	TPH C ₁₀ -C ₁₄ fraction	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.	Anonymous	TRH >C ₁₀ -C ₁₆ fraction	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment																				
Anonymous	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.																				
Anonymous	TPH C ₁₀ -C ₁₄ fraction	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.																				
Anonymous	TRH >C ₁₀ -C ₁₆ fraction	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.																				
Surrogate spike recovery	Surrogate recoveries were within control limits.																								
Blank monitoring																									
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 volatile TPH and TRH were reported below the LOR.																								
Method blank	Concentrations of all analytes were reported below the LOR.																								
Other observations	Batch ES1314611 <ul style="list-style-type: none">EG020: LOR's have been raised due to matrix interference (High Total Dissolved Solids).																								

Analytical Method	Analytical Parameter	Number of Tests Requested	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 LCS)		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 ANALYSES EDD06F	Arsenic	12	12	9		1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Cadmium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Chromium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Copper	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Lead	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Nickel	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Vanadium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Zinc	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Zinc	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
VOLATILES ANALYSES EP060	Benzene	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Benzene	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Benzene & para-Xylene	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Naphthalene	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	ortho-Xylene	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Toluene	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Volatiles ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Volatiles ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Volatiles ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	Volatiles ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
SEMIVOLATILES ANALYSES EP060/071	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
	SEMIVOLATILES ANALYSES EP060/071	12	12	9	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	1	✓	✓
METALS ANALYSES EDD06F	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Calcium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
METALS ANALYSES EDD06F	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Magnesium	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
METALS ANALYSES EDD06F	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
	Manganese	12	12	9	✓	1	1	1	1	0	0	1	2	1	1	0	0	2	2	1	1	1	1	-	-
INSTRUMENT LABORATORY NON-METALS EDD06P	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0	0	-	-
	Carbonate Alkalinity as CaCO3	10	10	9	✓	0	0	0	0	0	0	0	0	1	1	0	0	1	2	0	0	0			

Groundwater Analytical Results - March and June 2013 GME

[illegible]

[illegible]

LE0510	Exceedance of Adopted T's based on: Exceeds the ANZECC/ARMCANZ 2000 trigger values for the protection of aquatic ecosystems - Marine Water 29's species protection Exceeds the ANZECC/ARMCANZ 2000 trigger values for the protection of aquatic ecosystems - Freshwater 95% species protection "LOR Exceeds Guideline Trigger Value"
--------	--

Chem. Group	Chemicals	Conc. unit	EQI
Acidity	Carbonate as HCO3	mg/L	24.5
	Carbonate as CO3	mg/L	0.6
	Carbonate Alkalinity as CaCO3	mg/L	182
	Carbonate Alkalinity as CaCO3	mg/L	110
	Carbonate Alkalinity as CaCO3	mg/L	110
	Total Alkalinity as CaCO3	mg/L	110
	Toluene	mg/L	2
	Benzene	mg/L	1
	Ethylbenzene	mg/L	2
	Isobutylene	mg/L	2
BTX&N Compounds	Total Xylenes	mg/L	2
	Total BTEX	mg/L	1
	Naphthalene (VOC)	mg/L	5
	Phenol	mg/L	0.039
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
1,2,4-Hydroxyaromatics	Phenol	mg/L	0.039
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
Phenol-Phenol	Phenol	mg/L	0.039
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
	Phenol	mg/L	0.01
Halogenated Aromatic Compounds	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
Halogenated Aromatic Compounds	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
	1,1,1-Trichloroethane	mg/L	5
Major Ions	Calcium (Dissolved)	mg/L	20
	Magnesium (Dissolved)	mg/L	1
	Potassium (Dissolved)	mg/L	33
	Sulfate (as SO4) (Dissolved)	mg/L	0.1
	Sulfate (as SO4) (Dissolved)	mg/L	2
	Sulfate (as S) (Dissolved)	mg/L	1
	Sulfate (as SO4) (Dissolved)	mg/L	1
	Sulfate (as SO4) (Dissolved)	mg/L	1
	Sulfate (as SO4) (Dissolved)	mg/L	1
	Sulfate (as SO4) (Dissolved)	mg/L	1
Metals	Arsenic (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001
	Chromium (Dissolved)	mg/L	0.001

[illegible]

LEGEND

LEGEND

Frequency of Adopted Ilc bands are

Exceedance of Adopted Ls based on:
Exceeds the ANZECC/ARMCANZ 2000 trigger values for the protection of aquatic ecosystems - Marine Water
95 % species protection

Exceeds the ANZECC/ARMCANZ 2000 trigger values for the protection of aquatic ecosystems - Freshwater 95% species protection

* LOR Exceeds Guideline Trigger Value

[illegible]

Exceedance of Adopted ILs based on

Exceeds the ANZCC/ARMCAN/2000

Exceeds the ANZECC/ARMCANZ 2000

LOW EXCESS QUANTILE RISK VALUE

	Sample ID	Sample Date	Sample Type	SDG
Chem. Group	Benzene	-	-	-
	Toluene	-	-	-
	Ethylbenzene	-	-	-
	m,p-Xylene	-	-	-
	p-Xylene	-	-	-
	Total Xylenes	-	-	-
Metals	Naphthalene (VOC)	-	-	-
	Acetic Acid (Dissolved)	-	-	-
	Chromium (Dissolved)	-	-	-
	Copper (Dissolved)	-	-	-
	Lead	-	-	-
	Manganese (Dissolved)	-	-	-
	Mercury (Dissolved)	-	-	-
	Nickel (Dissolved)	-	-	-
Total Petroleum Hydrocarbons	Vanadium (Dissolved)	-	-	-
	Zinc (Dissolved)	-	-	-
	GC-C9 fraction	<20	<20	<20
	C10-C14 fraction	-	-	-
	C15-C28 fraction	-	-	-
	C29-C38 fraction	-	-	-
	C39-C43 fraction (sum)	-	-	-
Total Recoverable Hydrocarbons	GC-C10 fraction (T minus BTEX)	<20	<20	<20
	GC-C16-C18 fraction	-	-	-
	GC-C19-C24 fraction	-	-	-
	GC-C25-C40 fraction (sum)	-	-	-
	GC-C41-C43 fraction (sum)	-	-	-

10% check for March 2013
BA 9/9/2013

Table 2?
Analytical Results -
Darwin Waterfront - GME
(Client Name)

Chem. Group	Chem Name	Output Unit	EOL	ANZECC 2000 Freshwater 95%	ANZECC 2000 Marine water 95%	Lab Report Number	Monitoring Round	SDG	Sample Type	Sample Date, Time	Location	MW109	MW110	MW101	MW102	MW103	MW108	MW115	MW116	MW119
Alkalinity	Bicarbonate as HCO ₃	mg/L	5	151.04	90.12	151.28	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Hydroxide Alkalinity	mg/L	5	132	286	124	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Bicarbonate Alkalinity as CaCO ₃	mg/L	1	<1	<1	<1	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Hydroxide Alkalinity as CaCO ₃	mg/L	1	132	286	124	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Total Alkalinity as CaCO ₃	mg/L	1	132	286	124	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Toluene	µg/L	2	2	2	2	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Ethylbenzene	µg/L	2	2	2	2	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	m,p-Xylene	µg/L	2	2	2	2	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	o-Xylene	µg/L	2	2	2	2	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Total BTEX	µg/L	1	<1	<1	<1	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
G1 - C4 Hydrocarbon Gases	Methane	mg/L	0.01	<0.01	0.057	<0.01	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,2-Dichloroethane	mg/L	0.05	0.91	11.0	0.06	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Ferrous Iron	mg/L	0.05	<0.05	<0.05	<0.05	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Ferric Iron (Filtered)	mg/L	0.05	<0.05	<0.05	<0.05	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Bromonitrile	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,2-Dibromochloroethane (EUB)	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,2-Dibromochloroethane	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,2-Dibromochloroethane	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,1,1-Trichloroethane	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,1-Dichloroethane	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
Halogenated Aromatic Compounds	Carbon Tetrachloride	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Chlorobenzene	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,2-Dichlorobenzene	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,3-Dichlorobenzene	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,4-Dichlorobenzene	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,1,1-Trichloroethane	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,1,2-Trichloroethane	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,1,2,2-Tetrachloroethane	mg/L	50	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,2,3-Trichloropropane	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	1,2-Dibromo-3-chloropropane	mg/L	5	<5	<5	<5	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
Lead	Lead (Filtered)	mg/L	0.001	0.0034	0.0034	0.0034	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Lead (Unfiltered)	mg/L	0.001	0.0034	0.0034	0.0034	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Calcium (Filtered)	mg/L	1	1430	1430	1430	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Magnesium (Filtered)	mg/L	1	1260	1260	1260	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Potassium (Filtered)	mg/L	1	390	390	390	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sodium (as SO ₄)	mg/L	0.1	0.3	0.3	0.3	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sulphate (as SO ₄)	mg/L	2	84	84	84	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sulphur (as S)	mg/L	0.5	84	84	84	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sulphate (as S) (Filtered)	mg/L	1	260	260	260	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sulphate (as SO ₄) (Filtered)	mg/L	0.01	43	43	43	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
Major Ions	Total Anions	mg/L	0.01	43	43	43	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Total Cations	mg/L	0.01	138	138	138	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Ion Balance	%	0.01	138	138	138	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Lead	mg/L	0.001	0.0034	0.0034	0.0034	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Calcium	mg/L	1	1430	1430	1430	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Magnesium	mg/L	1	1260	1260	1260	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Potassium	mg/L	0.1	0.3	0.3	0.3	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sodium	mg/L	2	84	84	84	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sulphate	mg/L	0.5	84	84	84	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013
	Sulphate (Filtered)	mg/L	1	260	260	260	6/03/2013	Normal	Normal	6/03/2013	Normal	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013	6/03/2013

[illegible]