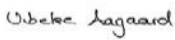



## Appendix B Data Validation

# DATA VALIDATION REPORT December 2010

<b>URS Project number:</b>	42213719	<b>Data verified by:</b>	Bek Aagaard	<b>Date:</b> 23/4/2013
<b>Client:</b>	Darwin Waterfront Corporation	<b>Signed:</b>		
<b>Site:</b>	Waterfront Precinct	<b>Validation by:</b>	Tim Smith	<b>Date:</b> 2/5/2013
<b>URS Project Manager:</b>	Jacques van Rensburg	<b>Signed:</b>		
<b>Matrix type:</b>	Groundwater	<b>Project Manager:</b>		<b>Date:</b>
<b>No Primary samples:</b>	10	<b>Signed:</b>		
<b>Laboratory:</b>	ALS			
	Labmark			
<b>Lab reference:</b>	ES1026649			
	E051579			

## 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.
Frequency of laboratory QC	<p>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:</p> <ul style="list-style-type: none"> <li>Laboratory duplicates were not reported for Total Petroleum Hydrocarbons (TPH) C10-C36. The precision of TPH is considered acceptable based on the data quality assessment of intra- and inter-laboratory duplicates.</li> <li>The inter-laboratory duplicate was not analysed for total anions and cations. The precision of this data is considered acceptable based on the data quality assessment of intra-laboratory field duplicates for total anions and cations.</li> </ul>
Tests requested/reported	Samples were analysed and reported as requested on the COC.
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"> <li>The protocol LOR exceeded ANZECC 2000 Aquatic Ecosystems (Marine 95%) for Chromium, Copper and Lead for samples MW10, MW106, MW111, MW113, MW115 and MW122.</li> <li>The protocol LOR exceeded ANZECC 2000 Aquatic Ecosystems (Irrigation - Long Term Use) for Iron for samples MW115 and MW122.</li> </ul> <p>The limits of reporting exceeded the guideline trigger levels for chromium copper and lead. 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	A random 10% check of the laboratory results identified no anomalies within the electronic data, the laboratory reports, and tables generated by URS.
<b>Sample management</b>	
Chain of Custody	Chain of custody documents completed.
Handling and preservation	<p>Samples were transported with ice bricks and were received at 16.8°C. Samples were received above the recommended temperature range; therefore, some losses through volatilisation may have occurred and sample concentrations for BTEXN and TPH C<sub>6</sub>–C<sub>9</sub> may be biased low. It should be noted that the groundwater temperature across the site throughout the year and at the time of sampling is generally &gt;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 24hr 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 are considered to be limited.</p>
Holding time compliance	<p>Analysis holding times were exceeded by 7-10 days for alkalinity. Given samples were sealed and stored in temperature controlled conditions at the laboratory, reported total alkalinity is not expected to have been affected and will not affect overall interpretation of results.</p>

## Data precision

### Field duplicate RPDs

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

- **MW128\_20/12/10** and QC01\_20/12/10 for Sodium (57.1%).

### Interlaboratory Replicate (Field triplicate) RPDs

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

- MW128\_20/12/10 and **QC02\_20/12/2010** for Potassium (133.3%).
- MW128\_20/12/10 and **QC02\_20/12/2010** for Sodium (41.8%).

### Laboratory duplicate RPDs

RPDs were within control limits.

## Data accuracy

### Laboratory control spike recovery

The following recoveries were outside control limits and should be considered during data interpretation:

Batch	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
ES1026649	Potassium	84.9	89	109	Recovery less than the lower control limit

For LCS recoveries reported less than the lower control limit, the potential exists for reported concentrations to be biased low by up to 4% for potassium. Care should be taken with interpreting results for potassium close to the adopted guidelines. As there is no applicable guideline this is not considered to have an impact on validity of conclusions.

LCS recoveries were not reported for 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, matrix LCS recoveries and surrogate recoveries (where applicable).

### Matrix spike recovery

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

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
MW136_17/12/10	Sulphate as SO4	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW108_17/12/10	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

As the matrix spike issue was 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).

Matrix spike recoveries were not reported for calcium, magnesium, potassium, sodium, alkalinity and total dissolved solids. The accuracy of the data for these compound groups is considered acceptable based on the presence of other quality control data, such as surrogate recoveries (where applicable).

### Surrogate spike recovery

The surrogate spike recoveries were within control limits.

## 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.

## Chromatograms

N/A

## Other observations

- EA015 TDS result has been confirmed by re-analysis for sample ID "MW115\_17/12/10".
- ED093F: LCS recovery for Potassium falls outside ALS Dynamic Control Limit. However, it is within the acceptance criteria based on ALS DQO. No further action is required.
- EG020: LOR's for some samples have been raised due to matrix interference (High sample salinity)
- EP080: Level of Reporting raised for toluene due to ambient background levels in the laboratory.
- The entire bottle is required for PAH, phenol and semi-volatile TPH determinations. Additional sample bottles are required for laboratory reporting of duplicates and matrix spikes.

Site Name Waterfront Precinct  
Project No. 42213719  
Project Manager Tim Smith  
Matrix Water  
Laboratory ALS - Labmark  
Batch File Number ESI0266449 - ESI1579

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Primary Samples	Holding Time (a)	Limit 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)		Duplicate Sample (1 in 20)		Lab Duplicate Sample (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			
METALS ANALYSIS/ALLEGORGE	Arsenic	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Chromium	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Chromium	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Copper	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Lead	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
METALS ANALYSIS/ALLEGORGE	Manganese	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Nickel	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Vanadium	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Zinc	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Boron	17	17	✓	✓	2	2	2	2	1	1	1	1	1	1	2	2	1	1	1	✓	✓
VOLATILES ANALYSIS/ALLEGORGE	Benzene	17	17	✓	✓	2	2	2	2	1	1	1	1	1	1	2	2	1	1	1	✓	✓
	Ethylbenzene	17	17	✓	✓	2	2	2	2	1	1	1	1	1	1	2	2	1	1	1	✓	✓
	meta & para Xylene	17	17	✓	✓	2	2	2	2	1	1	1	1	1	1	2	2	1	1	1	✓	✓
	ortho Xylene	17	17	✓	✓	2	2	2	2	1	1	1	1	1	1	2	2	1	1	1	✓	✓
	Toluene	17	17	✓	✓	2	2	2	2	1	1	1	1	1	1	2	2	1	1	1	✓	✓
VOLATILES ANALYSIS/ALLEGORGE	2,4-CD Fraction	17	17	✓	✓	2	2	2	2	1	1	1	1	1	1	2	2	1	1	1	✓	✓
	2,4-D Fraction	17	17	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	✓	✓
	2,4,6 Fraction	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	✓	✓
	2,4,6-CD Fraction	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	✓	✓
	2,4,6-CD Fraction	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	✓	✓
METALS ANALYSIS/ALLEGORGE	Sodium	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Magnesium	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Phosphorus	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Sodium	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Sodium	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
INSTRUMENT LABORATORY NON-METALS/ALLEGORGE	Sulfur as SO <sub>4</sub> Tolerant	16	16	✓	✓	2	2	2	2	0	0	1	1	1	1	2	2	1	1	1	-	-
	Mercury	16	16	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Barium Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Carbon Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Pyroxy Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
INSTRUMENT LABORATORY NON-METALS/ALLEGORGE	Total Activity as CaCO <sub>3</sub>	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
INSTRUMENT LABORATORY NON-METALS/ALLEGORGE	Total Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	12	12	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-
	Total Activity as CaCO <sub>3</sub>	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	2	2	1	1	1	-	-

Groundwater Analytical Results - Field Blanks  
December 2010

Location
Sample ID
Date Sampled
Sample Type
Area

QAQC	QAQC	QAQC	QAQC	QAQC	QAQC
QCB01_17/12/10	QCC01_17/12/10	QCA01_20/12/10	QCB02_20/12/10	QCC02_20/12/10	
17/12/2010	17/12/2010	20/12/2010	20/12/2010	20/12/2010	
Field Blank	Rinsate Blank	Trip Blank	Field Blank	Rinsate Blank	
QAQC	QAQC	QAQC	QAQC	QAQC	

Analyte	LOR	Units	ANZECC 2000 - Marine Water - 95%	ANZECC 2000 - Irrigation - LTU
<b>Total Petroleum Hydrocarbons</b>				
C6-C9 fraction	20	µg/L	<20	<20
C10-C14 fraction	50	µg/L	<50	<50
C15-C28 fraction	100	µg/L	<100	<100
C29-C36 fraction	50	µg/L	<50	<50
Sum of TPH C10 - C36	50	µg/L	<50	<50
C10-C36 fraction		µg/L	<200	<200
C6-C36 fraction		µg/L	<220	<220
<b>BTEX Compounds</b>				
Benzene	1	µg/L	<1	<1
Toluene	2	µg/L	<5	<5
Ethylbenzene	2	µg/L	<2	<2
m&p-Xylene	2	µg/L	<2	<2
o-Xylene	2	µg/L	<2	<2
Total Xylenes		µg/L	<4	<4
Total BTEX		µg/L	<12	<12
<b>Metals (Dissolved)</b>				
Arsenic	0.001	mg/L	<0.001	<0.001
Cadmium	0.0001	mg/L	<0.0001	<0.0001
Chromium	0.001	mg/L	<0.001	<0.001
Copper	0.001	mg/L	<0.001	<0.001
Lead	0.001	mg/L	<0.001	<0.001
Manganese	0.001	mg/L	<0.001	<0.001
Mercury	0.0001	mg/L	<0.0001	<0.0001
Nickel	0.001	mg/L	<0.001	<0.001
Vanadium	0.01	mg/L	<0.01	<0.01
Zinc	0.005	mg/L	<0.005	<0.005

Legend:

Exceeds the ANZECC/ARMCANZ, 2000, Trigger values for marine water ecosystems - Level of protection 95% species

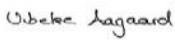

Exceeds the ANZECC/ARMCANZ, 2000, Irrigation

- Not analysed / not calculated

\* LOR Exceeds Guideline Trigger Value



# DATA VALIDATION REPORT February 2011

URS Project number:	42213719	Data verified by:	Bek Aagaard	Date:	24/4/2013
Client:	Darwin Waterfront Corporation				
Site:	Waterfront Precinct	Signed:			
URS Project Manager:	Jacques van Rensburg	Validation by:	Tim Smith	Date:	2/5/2013
Matrix type:	Groundwater				
No Primary samples:	24				
Laboratory:	ALS	Signed:			
Lab reference:	Labmark ES1103572, ES1103571, ES1102970, 291299 and 290561	Project Manager:		Date:	
		Signed:			

## Data quality objectives

Field data comparison	No apparent anomalies were observed between laboratory results and field observations.
Frequency of field QC	Field QA/QC was collected to project specifications (required number of field duplicates and triplicates collected in lab batches ES1103572 and ES1102970).
Frequency of laboratory QC	<p>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:</p> <ul style="list-style-type: none"> <li>Matrix spikes (MS) were reported on anonymous samples for several analytes. Where MS were conducted on anonymous samples, the reported MS recovery may not be considered representative of the URS field sample.</li> <li>Method blanks were not reported for alkalinity and total anions and cations in all batches. The accuracy of these analytes is considered acceptable based on the presence of surrogate recoveries for all batches and LCS recoveries for batch ES1102970.</li> <li>Laboratory duplicates were reported less than the required frequency for: <ul style="list-style-type: none"> <li>Total anions and cations in batch ES1102970;</li> <li>PAHs and Total Petroleum Hydrocarbons (TPH) in batch ES1103571; and</li> <li>PAHs, trihalomethanes, oxygenated compounds, Monocyclic Aromatic Compounds (MAH), carbon disulphide, halogenated aliphatic compounds, total dissolved solids, alkalinity, total anions and cations, sulphur as S, sulphite as <math>\text{SO}_3^{2-}</math> in batch ES1103572. The precision of these analytes is considered acceptable based on the data quality assessment of intra- and inter-laboratory duplicates.</li> </ul> </li> </ul>
Tests requested/reported	Samples were analysed and reported as requested on the COC.
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"> <li>Protocol LOR for anthracene (1 µg/L) above the adopted ILs (0.4 µg/L)</li> <li>Protocol LOR for benzo(a)pyrene (0.5 µg/L) above the adopted ILs (0.2 µg/L)</li> </ul> <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"> <li>The protocol LOR exceeded ANZECC 2000 Aquatic Ecosystems (Marine 95%) for Chromium, Copper and Lead for samples MW09, MW10, MW113 and MW115.</li> <li>The protocol LOR exceeded ANZECC 2000 Aquatic Ecosystems (Irrigation - Long Term Use) for Zinc.</li> </ul> <p>The limits of reporting exceeded the guideline trigger levels for chromium copper and lead. 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	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 for laboratory batch ES1102970 were received at 4.2°C, correctly preserved and chilled with ice. Samples for laboratory batches ES1103572 and ES1103571 were received between 23.9°C and 25.2°C. Samples were received above the recommended temperature range; therefore, some losses through volatilisation may have occurred and sample concentrations for BTEXN and TPH C<sub>6</sub> – C<sub>9</sub> may be biased low.

It should be noted that the groundwater temperature across the site throughout the year and at the time of sampling is generally >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 24hr 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 are considered to be limited.

## Holding time compliance

All samples were extracted and analysed within recommended holding times, with the exception of nitrate and sulphite holding times were exceeded during the 24 hour transit to the laboratory for all samples analysed in laboratory batches ES1103572 and ES1103571. These analytes are used as one of a number of indicators of natural attenuation and as such will not affect the overall interpretation of results.

## Data precision

### Field duplicate RPDs

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

- **MW118\_18/2/11** and QC05\_18/02/11 for Ferrous Iron (181.3%).
- **MW118\_18/2/11** and QC05\_18/02/11 for Ferric Iron (183.3%).
- **MW134\_10/02/11** and QC03\_10/02/11 for Manganese (34.2%).

Ferric iron and ferrous iron are used as one of a number of indicators, as relative concentrations of these analytes remains clear for interpretative purposes, this apparent lack of precision is not considered to affect the interpretation of results.

However, the apparent lack of precision for manganese should be taken into consideration when evaluating individual results, particularly where concentrations are reported close to the adopted investigation level.

### Interlaboratory Replicate (Field triplicate) RPDs

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

- **MW108\_9/02/11** and QC02\_09/02/11 for Methane (47.8%).
- **MW108\_9/02/11** and QC02\_09/02/11 for Potassium (38.8%).
- **MW108\_9/02/11** and QC02\_09/02/11 for Ferrous Iron (68.6%).
- MW118\_18/2/11 and **QC06\_18/02/11** for Potassium (30.8%).
- MW134\_10/02/11 and **QC04\_10/02/11** for Calcium (66.7%).
- MW134\_10/02/11 and **QC04\_10/02/11** for Sulphur as S (47.1%).

Methane, ferrous iron, potassium, calcium and sulphur (as S) are used as indicator species, as relative concentrations of these analytes remains clear for interpretative purposes, this apparent lack of precision is not considered to affect the interpretation of results.

### Laboratory duplicate RPDs

All laboratory duplicate RPDs were within control limits.

## Data accuracy

### Laboratory control spike recovery

Laboratory control spike recoveries outside the laboratory control limits were reported for at least one dissolved metal analyte and/or PAH in one or more batches, listed below.

Batch	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
ES1102970	Vanadium (dissolved)	87	91	109	Recovery less than lower control limit
ES1103571	Vanadium (dissolved)	79.8	91	109	Recovery less than lower control limit
ES1102970	Mercury (dissolved)	119	86	116	Recovery greater than upper control limit
ES1103571	Mercury (dissolved)	119	86	116	Recovery greater than upper control limit
ES1102970	Fluorene (dissolved)	62.7	63.9	115	Recovery less than lower control limit
ES1103571	Fluorene (dissolved)	62.7	63.9	115	Recovery less than lower control limit
ES1103572	Cadmium (dissolved)	109	89	107	Recovery greater than upper control limit
ES1103572	Chromium (dissolved)	114	91	111	Recovery greater than upper control limit
ES1103572	Lead (dissolved)	111	90	110	Recovery greater than upper control limit
ES1103572	Manganese (dissolved)	116	87	113	Recovery greater than upper control limit
ES1103572	Vanadium (dissolved)	85.5	91	109	Recovery less than lower control limit

For LCS recoveries reported greater than the upper control limit, the potential exists for reported concentrations to be biased high. Care should be taken with interpreting results for analytes close to



the adopted guidelines.

For LCS recoveries reported less than the lower control limit, the potential exists for reported concentrations to be biased low. Care should be taken with interpreting results for analytes close to the adopted guidelines.

LCS recoveries were not reported for total anions and cations from all batches, and alkalinity in batch ES1103571. 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, matrix LCS recoveries and surrogate recoveries (where applicable).

#### Matrix spike recovery

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

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
MW116_9/02/11	Manganese	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW103_10/02/11	Sulphate as SO <sub>4</sub>	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW113_14/02/11	Sulphate as SO <sub>4</sub>	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW113_14/02/11	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW113_14/02/11	Ferrous Iron	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

As the matrix spike issue was 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).

#### **Batch ES1102970**

Matrix spike recoveries were not reported for calcium, magnesium, potassium, sodium, alkalinity, total dissolved solids, trihalomethanes, aliphatic chlorinated compounds, monocyclic aromatic, oxygenated compounds 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 surrogate recoveries (where applicable).

#### **Batch ES1103571**

Matrix spike recoveries were not reported for PAHs, TPH C<sub>10</sub>-C<sub>36</sub>, calcium, magnesium, potassium, sodium, alkalinity, total dissolved solids, trihalomethanes, aliphatic halogenated compounds, monocyclic aromatic oxygenated compounds, chloride, total anions and cations, sulphur as S, sulphite as SO<sub>3</sub><sup>2-</sup> 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 surrogate recoveries (where applicable).

#### **Batch ES1103571**

Matrix spike recoveries were not reported for PAHs, TPH C<sub>10</sub>-C<sub>36</sub>, calcium, magnesium, potassium, sodium, alkalinity, total dissolved solids, trihalomethanes, aliphatic halogenated compounds, monocyclic aromatic and oxygenated compounds, total anions and cations, sulphur as S, sulphite as SO<sub>3</sub><sup>2-</sup> 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 surrogate recoveries (where applicable).

#### Surrogate spike recovery

A surrogate recovery of 78.9 % was reported for volatile TPH/BTEX compounds surrogate for sample QCC03\_14/02/11 in laboratory batch ES1103571. Therefore the reported concentrations of volatile TPH/BTEX compounds may be biased low by approximately one third. As concentrations of these compounds were not detected above the laboratory LOR this is not expected to affect interpretation of results.

#### **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.

**Other observations****Batch ES1102970**

- EG020: LCS recovery for Vanadium falls outside ALS Dynamic Control Limit. However, it is within the acceptance criteria based on ALS DQO. No further action is required.
- EG020: Samples ES1102970 #007 was diluted and rerun due to matrix interference and LOR's have been raised accordingly. (High sample salinity)
- IEG035F: LCS recovery for Mercury falls outside ALS Dynamic Control Limit. However, it is within the acceptance criteria based on ALS DQO. No further action is required
- EP074/EP080: Level of Reporting was raised for toluene due to ambient background levels in the laboratory.
- The entire bottle is required for PAH, phenol and semi-volatile TPH determinations. Additional sample bottles are required for laboratory reporting of duplicates.

**Batch ES1103571**

- ED041G: LOR raised for SO4 analysis on sample ID: L5 due to sample matrix
- EG020: LCS recoveries for some elements fall outside ALS Dynamic Control Limit. However, they are within the acceptance criteria based on ALS DQO. No further action is required.
- EG020: Some samples were diluted and rerun due to matrix interference and LOR's have been raised accordingly. (High sample salinity)
- EP074: Level of Reporting was raised for toluene due to ambient background levels in the laboratory.

**Batch ES1103571**

- EG020: LCS recoveries for some elements fall outside ALS Dynamic Control Limit. However, they are within the acceptance criteria based on ALS DQO. No further action is required.
- EG020A-T: An unpreserved aliquot from the natural bottle was used for analysis.
- EP074: Level of Reporting was raised for toluene due to ambient background levels in the laboratory.

Site Name Waterfront Precinct  
Project No. 42213719  
Project Manager Tim Smith  
Matrix Water  
Laboratory ALS / Labmark  
Batch File Number ESI102970 / 290561

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Number of Primary Samples	Holding Times (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)		Intr-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/SEDOF	Arsenic	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Cadmium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Chromium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Copper	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Lead	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Manganese	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Nickel	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Selenium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Sodium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
	Zinc	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	4	1	2	1	2	-
VOLATILES ANALYSIS/SEDOF	Benzene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	Ethylbenzene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	m,p-xylene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	o-xylene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	Toluene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	Chloroform	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Diethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Styrene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Triethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Phenol	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
SEMI-VOLATILES ANALYSIS/SEDOF	Chloroform	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Diethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Styrene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Triethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Phenol	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Chlorobenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Diethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Styrene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Triethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Phenol	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
METALS ANALYSIS/SEDOF	Arsenic	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Cadmium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Chromium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Copper	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Lead	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Manganese	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Nickel	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Selenium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Sodium	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
	Zinc	21	21	13	✓	✓	2	2	2	2	0	0	1	2	1	2	2	2	1	1	1	1	✓
VOLATILES ANALYSIS/SEDOF	Benzene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	Ethylbenzene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	m,p-xylene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	o-xylene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	Toluene	21	21	13	✓	✓	4	4	4	4	0	0	1	4	1	4	2	2	1	1	1	1	✓
	Chloroform	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Diethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Styrene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Triethylbenzene	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓
	Phenol	21	21	13	✓	✓	2	2	2	2	0	0	1	2	2	2	2	2	1	1	1	1	✓

Site Name Waterfront Precinct  
Project No. 42213719  
Project Manager Tim Smith  
Matrix Water  
Laboratory ALS / Labmark  
Batch File Number ESI102970 / 290561

Analytical Method	Analytical Parameters	Number of Tests Requested	Number of Tests Reported	Holding Times (h)	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)		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	
VOLATILES ANALYSISALSEP04/E	1,1,1-Trichloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,1,2-Trichloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,1,2,2-Tetrachloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,2-Dichloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,1-Dichloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,1,1-Trichloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,1-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,2,3-Trichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,2,3-Dibromo-4-chloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
VOLATILES ANALYSISALSEP04/F	1,2-Dichloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,3-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Benzonitrile	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon Tetrachloride	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chloroethane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
VOLATILES ANALYSISALSEP04/G	1,2-Dichloroethane (ED)	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2,2-Dichloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
VOLATILES ANALYSISALSEP04/H	1,2,3-Trichlorobenzene	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	1,2,4-Trichlorobenzene	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	1,2,5-Trichlorobenzene	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	1,3-Dichlorobenzene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,4-Dichlorobenzene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	2-Chlorodurene	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	4-Chlorodurene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Benzobenzene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chlorobenzene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Chlorobenzene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
VOLATILES ANALYSISALSEP04/I	1,2,4-Trinitrobenzene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	1,3,5-Trinitrobenzene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Benzoylethene	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Phenylacetone	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Phenylacetone	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Phenylacetone	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Phenylacetone	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Phenylacetone	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Phenylacetone	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Phenylacetone	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
VOLATILES ANALYSISALSEP04/J	2-Bromo-4-chloropropane	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	2-Bromo-4-chloropropane	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
VOLATILES ANALYSISALSEP04/K	4-Methyl-2-pentanone (MEK)	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Vinyl Acetate	19	19	✓	✓	2	2	2	2	0	0	1	1	1	2	1	0	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓
	Carbon disulfide	21	21	✓	✓	2	2	2	2	0	0	1	1	1	2	1	2	2	1	0	1	✓

NOTES:  
(a) Holding times are within project guidelines.  
(b) ✓ Indicates successful project guidelines.  
(c) ✓ Limits of reporting (LORs) comply with project specifications.  
(d) ✓ LORs do not comply with project specifications.  
NA Not Applicable

Site Name Waterfront Precinct  
Project No. 42213719  
Project Manager Tim Smith  
Matrix Water  
Laboratory ALS  
Batch File Number ESI103571

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Primary Samples	Holding Time (h)	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 esky)		Inter-Laboratory Duplicate Sample (1 in 20)		Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)	Surrogates (GCMS 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			
METALS ANALYSIS/ELSEP924F	Arsenic	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Chromium	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Copper	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Lead	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Manganese	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Nickel	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Vanadium	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	Zinc	9	9	7	✓	✓	1	1	1	1	0	0	1	2	0	0	1	4	1	2	1	2
	OS - GP Fraction	10	10	7	✓	✓	1	1	1	1	1	1	1	1	0	0	1	2	1	1	1	✓
VOLATILES ANALYSIS/ELSEP924F	Acenaphthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Acenaphthylene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Anthracene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benz[a]fluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benz[a]pyrene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
METALS ANALYSIS/ELSEP924F	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
INSTRUMENT LABORATORY NON-METALS ANALYSIS/ELSEP924F	Arsenic	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Chromium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Copper	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Lead	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Manganese	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Nickel	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Vanadium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Zinc	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	OS - GP Fraction	10	10	7	✓	✓	1	1	1	1	1	1	1	1	0	0	1	2	1	1	1	✓
VOLATILES ANALYSIS/ELSEP924F	Acenaphthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Acenaphthylene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Anthracene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benz[a]fluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benz[a]pyrene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
METALS ANALYSIS/ELSEP924F	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
VOLATILES ANALYSIS/ELSEP924F	Acenaphthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Acenaphthylene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Anthracene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benz[a]fluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benz[a]pyrene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Benzofluoranthene	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
METALS ANALYSIS/ELSEP924F	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	0	0	1	0	1	0	1	✓
	Cadmium	9	9																			



Site Name Waterfront Precinct  
Project No. 42213719  
Project Manager Tim Smith  
Matrix Water  
Laboratory ALS / SGS  
Batch File Number ESI103572 / 291299

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Primary Samples	Holding Time (h)	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)		Pre-Laboratory Duplicate Sample (1 in 20)		In-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 REF02F	Arsenic	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Cadmium	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Chromium	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Copper	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Lead	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Manganese	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Nickel	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Vanadium	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS REF03F	Zinc	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Cr: Cr Fraction	8	9	4	✓	✓	1	1	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
	Acenaphthene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Acenaphthylene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Anthracene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[a]anthracene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[a]fluorene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[a]pyrene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[b]fluoranthene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[b]fluorene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[k]fluoranthene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[k]pyrene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Chrysene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Diethylanthracene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Fluoranthene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Fluorene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Indeno[1,2,3-cd]pyrene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Naphthalene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Phenanthrene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Pyrene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
SEMIVOLATILES ANALYSIS REF04F	Cr: Cr Fraction	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Cr: Cr Fraction (a/r)	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Cr: Cr Fraction	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Cr: Cr Fraction	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
METALS ANALYSIS REF05F	Cadmium	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Magnesium	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Potassium	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Sodium	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
INSTRUMENT LABORATORY NON-METAL REF06F	Nitrate as N	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Nitrite as N	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Nitrate + Nitrite as N	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Nitrate as N	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
INSTRUMENT LABORATORY NON-METAL REF07F	Sulfate as S	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Sulfate as S	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Sulfate as S	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Sulfate as S	6	6	4	✓	✓	0	0	0	0	0	0	1	1	1	1	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS REF08F	Benzo[a]anthracene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[a]pyrene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[b]fluoranthene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[k]fluoranthene	8	8	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	0	1	1	1	1	✓	✓
	Benzo[e]pyrene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[g]herylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[h]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[i]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[j]fluoranthene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[k]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[l]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[m]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[n]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[o]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[p]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[q]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[r]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[s]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[t]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Benzo[u]perylene	7	7	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓





Sample ID
Date Sampled
Sample Type

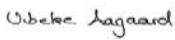

Analyte	LOR	Units	ANZECC 2000 - Marine Water - 85%	ANZECC 2000 - Irrigation - LTU
<b>Total Petroleum Hydrocarbons</b>				
C6-C8 fraction	20	µg/L	<20	<20
C10-C14 fraction	0.05	µg/L	<0.05	<0.05
C15-C28 fraction	0.1	mg/L	<0.1	<0.1
C29-C36 fraction	0.1	mg/L	<0.05	<0.05
Sum of TPH C10 - C36	50	µg/L	<50	<50
<b>BTEX Compounds</b>				
Benzene	0.001	mg/L	<0.001	<0.001
Toluene	0.001	mg/L	<0.005	<0.005
Ethylbenzene	0.001	mg/L	<0.002	<0.002
m,p-Xylene	2	µg/L	<2	<2
o-Xylene	0.001	mg/L	<0.002	<0.002
<b>Metals (Dissolved)</b>				
Asenic	0.001	mg/L	<0.001	<0.001
Cadmium	0.0001	mg/L	<0.0001	<0.0001
Chromium	0.001	mg/L	<0.001	<0.001
Copper	0.001	mg/L	<0.001	<0.001
Iron	0.05	mg/L	<2	<2
Lead	0.001	mg/L	<0.001	<0.001
Manganese	0.001	mg/L	<0.001	<0.001
Mercury	0.0001	mg/L	<0.0001	<0.0001
Nickel	0.001	mg/L	<0.001	<0.001
Vanadium	0.005	mg/L	<0.001	<0.001
Zinc	0.005	mg/L	<0.005	<0.005
<b>Polynuclear Aromatic Hydrocarbons</b>				
Naphthalene	0.001	mg/L	<0.001	<0.001
Acenaphthylene	0.001	mg/L	<0.001	<0.001
Acenaphthene	0.001	mg/L	<0.001	<0.001
Fluorene	0.001	mg/L	<0.001	<0.001
Phenanthrene	0.001	mg/L	<0.001	<0.001
Anthracene	0.001	mg/L	<0.001	<0.001
Fluoranthene	0.001	mg/L	<0.001	<0.001
Pyrene	0.001	mg/L	<0.001	<0.001
Benz[a]anthracene	0.001	mg/L	<0.001	<0.001
Chrysene	0.001	mg/L	<0.001	<0.001
Benz[b]fluoranthene	1	µg/L	<1	<1
Benz[k]fluoranthene	1	µg/L	<1	<1
Benz[a]pyrene	0.001	mg/L	<0.0005	<0.0005
Indene[1,2,3-c]pyrene	0.001	mg/L	<0.001	<0.001
Dibenz[a,h]anthracene	0.001	mg/L	<0.001	<0.001
Benz[g,h,i]perylene	0.001	mg/L	<0.001	<0.001
Benz[b]fluoranthene & Benz[c]fluoranthene	0.002	mg/L	<2	<2
<b>Chlorinated Aromatic Compounds</b>				
1,2,3-Trichlorobenzene	5	µg/L	<5	<5
1,2,4-Trichlorobenzene	5	µg/L	<5	<5
1,2,4-Trimethylbenzene	0.001	mg/L	<0.005	<0.005
1,2-dichlorobenzene	0.001	mg/L	<0.005	<0.005
1,3-Dichlorobenzene	0.001	mg/L	<0.005	<0.005
1,4-Dichlorobenzene	0.001	mg/L	<0.005	<0.005
2-Chlorotoluene	5	µg/L	<5	<5
4-Chlorotoluene	0.001	mg/L	<0.005	<0.005
Bromobenzene	0.001	mg/L	<0.005	<0.005
Chlorobenzene	0.001	mg/L	<0.005	<0.005







# DATA VALIDATION REPORT\_June 2011

<b>URS Project number:</b>	42213719	<b>Data verified by:</b>	Bek Aagaard	<b>Date:</b> 24/04/2013
<b>Client:</b>	Darwin Waterfront Corporation	<b>Signed:</b>		
<b>Site:</b>	Waterfront Precinct	<b>Validation by:</b>	Tim Smith	<b>Date:</b> 3/5/2013
<b>URS Project Manager:</b>	Jacques van Rensburg	<b>Signed:</b>		
<b>Matrix type:</b>	Groundwater	<b>Project Manager:</b>		<b>Date:</b>
<b>No Primary samples:</b>	10	<b>Signed:</b>		
<b>Laboratory:</b>	ALS; Labmark			
<b>Lab reference:</b>	ES1113641 304025			

## 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.
Frequency of laboratory QC	<p>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:</p> <ul style="list-style-type: none"> <li>Matrix spikes (MS) were reported on anonymous samples for all analytes. Where MS were conducted on anonymous samples, the reported MS recovery may not be considered representative of the URS field sample.</li> <li>Laboratory duplicate were not reported for Total Petroleum Hydrocarbons (TPH)/Total Recoverable Hydrocarbons (TRH). The precision of TPH and TRH has been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable.</li> <li>The inter-laboratory duplicate was not analysed for TRH, calcium, magnesium, potassium, sodium and alkalinity. The precision of this data can be assessed as acceptable based on the presence of intra-laboratory field duplicates for all analytes and laboratory duplicates except for TRH.</li> </ul>
Tests requested/reported	Samples were analysed and reported as requested on the COC.
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"> <li>Protocol LOR for anthracene (1 µg/L) above the adopted ILs (0.4 µg/L)</li> <li>Protocol LOR for benzo(a)pyrene (0.5 µg/L) above the adopted ILs (0.2 µg/L)</li> </ul> <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"> <li>Protocol LOR exceeds the guideline trigger value for chromium (0.0044 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for copper (0.0013 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for lead (0.0044 mg/L)</li> </ul> <p>The limits of reporting exceeded the guideline trigger levels for chromium copper and lead. 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	A random 10% check of the laboratory results identified no anomalies within the electronic data, the laboratory reports, and tables generated by URS.
<b>Sample management</b>	
Chain of Custody	Chain of custody documents completed.
Handling and preservation	<p>Samples were correctly preserved, chilled with ice and received at the laboratory at 7.6°C.</p> <p>Samples from batch ES1113641 were received above the recommended temperature range; therefore, some losses through volatilisation may have occurred and sample concentrations for BTEXN and TPH C<sub>6</sub> – C<sub>9</sub> may be may be biased low.</p>



It should be noted that the groundwater temperature across the site throughout the year and at the time of sampling is generally >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 24hr 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 are considered to be limited.

Holding time compliance

Samples were extracted and analysed within recommended holding times.

#### Data precision

Field duplicate RPDs

RPDs were within control limits.

Interlaboratory Replicate (Field triplicate) RPDs

RPDs were within control limits.

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
ES1113641	Mercury	126	86	116	Recovery greater than upper control limit
ES1113641	Mercury	119	86	116	Recovery greater than upper control limit

For LCS recoveries reported greater than the upper control limit, the potential exists for reported concentrations to be biased high by up to 10% for mercury. Care should be taken with interpreting results for mercury close to the adopted guidelines. As results for mercury were reported below LOR, 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
MW116_27/06/11	Manganese	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

As the matrix spike issue was 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).

Matrix spike recoveries were not reported for TPH C10-C36, calcium, magnesium, potassium, sodium, total dissolved solids 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 surrogate spike recoveries were within control limits.

#### 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.

#### Chromatograms

N/A

#### Other observations

- EG020: MW115\_23/06/11 was diluted and reanalysed due to matrix interference(s) caused by high salinity. LORs have been raised accordingly.
- EP080: Level of reporting raised for toluene due to ambient background levels in the laboratory.
- LCS recovery for Mercury falls outside ALS Dynamic Control Limit. However, it is within the acceptance criteria based on ALS DQO. No further action is required.
- The entire bottle is required for PAH, phenol and semi-volatile TPH determinations. Additional sample bottles are required for laboratory reporting of duplicates and matrix spikes.

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

Waterfront Precinct  
422137119  
Tim Smith  
Water  
ALS / Labmark  
ESI1113641 / 304025

Analytical Method	Analytical Parameter	Number of Tests Requested		Number of Primary Samples	Holding Times (a)	Unit 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)		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)		
		Requested	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	Reported	OK	
METALS ANALYSEALS E6002F	Arsenic	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Cadmium	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Chromium	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Copper	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Lead	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Zinc	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Vanadium	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Barium	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Zinc	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	4	1	2	1	2	-	-	
	Benzene	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	2	1	1	✓	✓	
	Ethylbenzene	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	2	1	1	✓	✓	
	m,p-xylene	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	2	1	1	✓	✓	
	p-xylene	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	2	1	1	✓	✓	
	o-xylene	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	2	1	1	✓	✓	
VOLATILES ANALYSEALS E6001	Hexane	14	14	10	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	2	2	1	2	1	1	✓	✓	
	2,4- DFB Fraction	15	15	10	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓	
	2,6- DFB Fraction	15	15	10	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	2	2	1	1	1	1	✓	✓	
	Semivolatiles ANALYSEALS E6003	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Acetophenone	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Acetophenone	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Acetophenone	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Benzofuran	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Benzofuran	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Benzofuran	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Benzofuran	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Benzofuran	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
	Benzofuran	2	2	0	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	0	0	0	0	0	0	1	✓	
SEMIVOLATILES ANALYSEALS E6003F	2,4- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,6- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,4- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,6- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,4- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,6- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,4- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,6- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,4- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,6- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,4- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,6- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,4- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
	2,6- DFB Fraction	13	13	10	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	2	0	1	0	1	1	✓	✓	
METALS ANALYSEALS E6002F	Cadmium	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Chromium	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Copper	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Sodium	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Vanadium	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Barium	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Zinc	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Iron	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Lead	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Aluminum	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Calcium	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Chlorine	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Fluorine	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Sulfur	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
INSTRUMENT LABORATORY NON-METALS E6002F	Wet Chemistry and Preparation of Samples	12	12	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14	14	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	0	2	4	1	0	1	2	-	-
	Carbonate Alkalinity as CaCO <sub>3</sub>	14																									

Groundwater Analytical Results - Field Blanks  
June 2011

Sample ID
Date Sampled
Sample Type

QCA01_27/06/11	QCB01_27/06/11	QCC01_27/06/11
27/06/2011	27/06/2011	27/06/2011
Trip Blank	Field Blank	Rinsate Blank

Analyte	LOR	Units	ANZECC 2000 - Marine Water - 95%	ANZECC 2000 - Irrigation - LTU			
<b>Total Petroleum Hydrocarbons</b>							
C6-C9 fraction	20	µg/L			<20	<20	<20
C10-C14 fraction	50	µg/L			-	<50	<50
C15-C28 fraction	100	µg/L			-	<100	<100
C29-C36 fraction	50	µg/L			-	<50	<50
C10 - C36 Fraction (sum)	50	µg/L			-	<50	<50
<b>BTEX Compounds</b>							
Benzene	1	µg/L	700		-	<1	<1
Toluene	2	µg/L			-	<5	<5
Ethylbenzene	2	µg/L			-	<2	<2
m&p-Xylene	2	µg/L			-	<2	<2
o-Xylene	2	µg/L			-	<2	<2
Total Xylenes	2	µg/L			-	<2	<2
Sum of BTEX	1	µg/L			-	<1	<1
<b>Metals (Dissolved)</b>							
Calcium	0.1	mg/L			-	-	-
Potassium	1	mg/L			-	-	-
Sodium	0.1	mg/L			-	-	-
Arsenic	0.001	mg/L		0.1	-	<0.001	<0.001
Cadmium	0.0001	mg/L	0.0055	0.01	-	<0.0001	<0.0001
Chromium	0.001	mg/L	0.0044	0.1	-	<0.001	<0.001
Copper	0.001	mg/L	0.0013	0.2	-	<0.001	<0.001
Lead	0.001	mg/L	0.0044	2	-	<0.001	<0.001
Magnesium	1	mg/L			-	-	-
Manganese	0.001	mg/L		0.2	-	<0.001	<0.001
Mercury	0.0001	mg/L	0.0004	0.002	-	<0.0001	<0.0001
Nickel	0.001	mg/L	0.07	0.2	-	<0.001	<0.001
Vanadium	0.01	mg/L	0.1	0.1	-	<0.01	<0.01
Zinc	0.005	mg/L	0.015	2	-	<0.005	<0.005
<b>Polynuclear Aromatic Hydrocarbons</b>							
Naphthalene	5	µg/L	70		-	<1	<1
Acenaphthylene	1	µg/L			-	<1	<1
Acenaphthene	1	µg/L			-	<1	<1
Fluorene	1	µg/L			-	<1	<1
Phenanthrene	1	µg/L			-	<1	<1
Anthracene	1	µg/L			-	<1	<1
Fluoranthene	1	µg/L			-	<1	<1
Pyrene	1	µg/L			-	<1	<1
Benz(a)anthracene	1	µg/L			-	<1	<1
Chrysene	1	µg/L			-	<1	<1
Benzo(b)fluoranthene	1	µg/L			-	<1	<1
Benzo(k)fluoranthene	1	µg/L			-	<1	<1
Benzo(a)pyrene	0.5	µg/L			-	<0.5	<0.5
Indeno(1,2,3-cd)pyrene	1	µg/L			-	<1	<1
Dibenz(a,h)anthracene	1	µg/L			-	<1	<1
Benzo(g,h,i)perylene	1	µg/L			-	<1	<1
Sum of polycyclic aromatic hydrocarbons	0.5	µg/L			-	<1	<0.9
<b>Total Recoverable Hydrocarbons</b>							
>C10 - C16 Fraction	100	µg/L			-	<100	<100
>C10 - C40 Fraction (sum)	100	µg/L			-	<100	<100
>C16 - C34 Fraction	100	µg/L			-	<100	<100
>C34 - C40 Fraction	100	µg/L			-	<100	<100
C6 - C10 Fraction	20	µg/L			<20	<20	<20
C6 - C10 Fraction minus BTEX (F1)	20	µg/L			-	<20	<20

Legend:

Exceeds the ANZECC/ARMCANZ, 2000, Trigger values for marine water ecosystems - Level of protection 95% species

Exceeds the ANZECC/ARMCANZ, 2000, Irrigation

- Not analysed / not calculated

\* LOR Exceeds Guideline Trigger Value

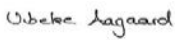



10% check  
18/4/2013 VBA

[illegible]

**Legend:**

# DATA VALIDATION REPORT September 2011

<b>URS Project number:</b>	42213719	<b>Data verified by:</b>	Bek Aagaard	<b>Date:</b> 24/04/2013
<b>Client:</b>	Darwin Waterfront Corporation			
<b>Site:</b>	Waterfront Precinct	<b>Signed:</b>		
<b>URS Project Manager:</b>	Jacques van Rensburg	<b>Validation by:</b>	Tim Smith	<b>Date:</b> 3/5/2013
<b>Matrix type:</b>	Groundwater			
<b>No Primary samples:</b>	22			
<b>Laboratory:</b>	ALS; SGS	<b>Signed:</b>		
<b>Lab reference:</b>	ES1121061, EM1110994, EM1111058, SE89800, SW102312 and SW102311	<b>Project Manager:</b>		<b>Date:</b>
		<b>Signed:</b>		

## 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.
Frequency of laboratory QC	<p>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:</p> <ul style="list-style-type: none"> <li>Matrix spikes (MS) were reported on anonymous samples for several analytes. Where MS were conducted on anonymous samples, the reported MS recovery may not be considered representative of the URS field sample.</li> <li>Laboratory duplicate were not reported for ferric iron, total anions and cations in all batches, PAHs and Total Petroleum Hydrocarbons (TPH)/Total Recoverable Hydrocarbons (TRH) in batches EM1110994 and EM1111058 and trihalomethanes in batch EM1111058. The precision of these analytes have been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable.</li> <li>The inter-laboratory duplicate was reported less than the required frequency for analysed for pentachloroethene, total anions and cations and alkalinity. The precision of this data can be assessed as acceptable based on the presence of intra-laboratory field duplicates for all analytes and laboratory duplicates.</li> </ul>
Tests requested/reported	Samples were analysed and reported as requested on the COC.
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"> <li>Protocol LOR for anthracene (1 µg/L) above the adopted ILs (0.4 µg/L)</li> <li>Protocol LOR for benzo(a)pyrene (0.5 µg/L) above the adopted ILs (0.2 µg/L)</li> </ul> <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"> <li>Protocol LOR exceeds the guideline trigger value for chromium (0.0044 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for copper (0.0013 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for lead (0.0044 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for zinc (0.015 mg/L)</li> </ul> <p>The limits of reporting exceeded the guideline trigger levels for 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 below 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.
<b>Sample management</b>	
Chain of Custody	Chain of custody documents completed.
Handling and preservation	Samples were correctly preserved, chilled with ice and received at the laboratory at 3.0°C, 1.0-4.5°C and 4.5-6.0°C for batches ES1121061, EM1110994 and EM1111058, respectively.



Holding time compliance

Samples were extracted and analysed within recommended holding times.

#### Data precision

Field duplicate RPDs

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

- **MW119\_27/09/11** and QC01\_27/09/11 for Ionic Balance (56.8%).
- **MW116\_28/09/11** and QC05\_28/09/11 for Bicarbonate Alkalinity (36.9%).
- **MW116\_28/09/11** and QC05\_28/09/11 for Ionic Balance (30.1%).
- **MW108\_29/09/11** and QC09\_29/09/11 for Ionic Balance (152.2%).
- **MW108\_29/09/11** and QC09\_29/09/11 for Methane (42%).
- **MW108\_29/09/11** and QC09\_29/09/11 for Methane (42%).

Inter-laboratory Replicate (Field triplicate) RPDs

RPDs were within control limits.

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
ES1121061	Arsenic	115	88	110	Recovery greater than upper control limit
ES1121061	Cadmium	111	89	107	Recovery greater than upper control limit
ES1121061	Manganese	118	87	113	Recovery greater than upper control limit
ES1121061	Nickel	116	89	109	Recovery greater than upper control limit
ES1121061	Vanadium	89.0	91	109	Recovery less than lower control limit
ES1121061	Acenaphthylene	62.7	63.6	114	Recovery less than lower control limit
ES1121061	Fluorene	62.3	63.9	115	Recovery less than lower control limit
EM1110994	Bromobenzene	72.9	75	119	Recovery less than lower control limit
EM1110994	1,4-Dichlorobenzene	73.1	74	120	Recovery less than lower control limit
EM1110994	1,2-Dichlorobenzene	73.5	73.5	118	Recovery less than lower control limit

For LCS recoveries reported greater than the upper control limit, the potential exists for reported concentrations to be biased high. Care should be taken with interpreting results for analytes close to the adopted guidelines.

For LCS recoveries reported less than the lower control limit, the potential exists for reported concentrations to be biased low. Care should be taken with interpreting results for analytes close to the adopted guidelines.

LCS recoveries were not reported for total anions and cations in all batches, and ferric iron in batches EM1110998 and EM1111058. 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, matrix spike recoveries and surrogate recoveries (where applicable).

Matrix spike recovery

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

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
MW124_27/09/11	Sulphate as SO4 - Turbidimetric	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW124_27/09/11	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW117_27/09/11	Manganese	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW124_27/09/11	Ferrous Iron	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Batch EM1110994 Anonymous	Sulphate as SO4 - Turbidimetric	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

MW136_28/09/11	Manganese	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW116_28/09/11	Ferrous Iron	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW136_28/09/11	Methane	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Batch EM1111058 Anonymous	Sulphate as SO4 - Turbidimetric	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Batch EM1111058 Anonymous	Ferrous Iron	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

As the matrix spike issue was 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).

#### **All Batches**

Matrix spike recoveries were not reported for calcium, magnesium, potassium, sodium, trihalomethanes, oxygenated compounds, monocyclic aromatic compounds, carbon disulphide, halogenated aliphatics, sulphite, sulphur, total dissolved solids, total anions and cations 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).

#### **EM1110998**

Matrix spike recoveries were not reported for PAHs, sulphide, TPH C<sub>10</sub>-C<sub>36</sub> and TRH C<sub>10</sub>-C<sub>40</sub>. 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 surrogate spike recoveries were within control limits.

#### **Blank monitoring**

Equipment rinsate blank

Concentrations of all analytes were reported below the LOR, except for QC07\_28/09/2011, which reported a chromium concentration equal to the LOR. Chromium was detected in two primary samples during this sampling round on the previous day of sampling (MW123\_27/09/2011 and MW124\_27/09/2011), given that chromium was not detected in any other primary or QC samples this is not considered to be indicative of potential for cross contamination and is not considered to affect interpretation of analytical 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.

#### **Chromatograms**

N/A

#### **Other observations**

##### **Batch ES1121061**

- EG020-F: LCS recovery for some elements fall outside ALS dynamic control limit, however they are within the acceptance criteria based on ALS DQO. No further action is required.
- EP080: Level of reporting raised for toluene due to ambient background levels in the laboratory.

##### **Batch EM1110998**

- EG020F : Chromium result for EM1110994 #9 has been confirmed by re-preparation and re-analysis.
- Ionic balances were calculated using: major anions - chloride, alkalinity and sulphate; and major cations - calcium, magnesium, potassium and sodium.
- Methane analysis conducted by ALS Sydney, NATA accreditation no. 825, site no 10911.

##### **Batch EM1111058**

- Ionic Balance out of acceptable limits for sample EM1111058 #6 & #7 due to analytes not quantified in this report.
- Ionic balances were calculated using: major anions - chloride, alkalinity and sulphate; and major cations - calcium, magnesium, potassium and sodium.
- Methane conducted by ALS Sydney, NATA accreditation no. 825, site no 10911











Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Number of Primary Samples	Holding Time (h)	Units of Reporting (b)	Field Blank (1 per day)		Finsate Blank (1 per day)		Trip Blank (1 per csk with VOCs)		Method Blank (1 per batch)		Reference 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	Reported	OK
METALS ANALYSIS REF002#	Arsenic	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Barium	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Boron	11	11	7	✓	✓	1	1	1	1	0	0	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	2	1	1	1	1	-	-
	Lead	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Manganese	11	11	7	✓	✓	1	1	1	1	0	0	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	2	1	1	1	1	-	-
	Vanadium	11	11	7	✓	✓	1	1	1	1	0	0	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	2	1	1	1	1	-	-
	11	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
VOLATILES ANALYSIS REF001	Benzene	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Ethylbenzene	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	m,p- & para-Xylene	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Napthalene	11	11	4	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	ortho-Xylene	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	11	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	11	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	11	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	11	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	11	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
METALS ANALYSIS REF006#	Calcium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Magnesium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Potassium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Sodium	9	9	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	-	-
	Boron	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Boron	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Boron	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Boron	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Boron	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Boron	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS REF004	1,1,1-Trichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1,1,2-Tetrachloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1,2,2-Tetrachloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1,2-Trichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,3-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS REF007	Bromomethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Carbon Tetrachloride	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	Chloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1,1-Trichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1,2,2-Tetrachloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1,2-Trichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,1-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS REF008	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
VOLATILES ANALYSIS REF009	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓
	1,2-Dichloroethane	11	11	7	✓	✓	1	1	1	1	0	0	1	1	1	1	1	2	1	1	1	1	✓	✓

Surrogate Sample ID	Lab Duplicate (1 in 10)		Matrix Spike (1 in 20)		LCS (1 per batch)		Surrogates (6-CAS organics)	
	Number Reported	Number Required	Number Reported	Number Required	Number Reported	Number Required	Reported	OK
1	1	2	1	0	1	1	✓	✓
2	1	2	1	0	1	1	✓	✓
3	1	2	1	0	1	1	✓	✓
4	1	2	1	0	1	1	✓	✓
5	1	2	1	0	1	1	✓	✓
6	1	2	1	0	1	1	✓	✓
7	1	2	1	0	1	1	✓	✓
8	1	2	1	0	1	1	✓	✓
9	1	2	1	0	1	1	✓	✓
10	1	2	1	0	1	1	✓	✓
11	1	2	1	0	1	1	✓	✓
12	1	2	1	0	1	1	✓	✓
13	1	2	1	0	1	1	✓	✓
14	1	2	1	0	1	1	✓	✓
15	1	2	1	0	1	1	✓	✓
16	1	2	1	0	1	1	✓	✓
17	1	2	1	0	1	1	✓	✓
18	1	2	1	0	1	1	✓	✓
19	1	2	1	0	1	1	✓	✓
20	1	2	1	0	1	1	✓	✓
21	1	2	1	0	1	1	✓	✓
22	1	2	1	0	1	1	✓	✓
23	1	2	1	0	1	1	✓	✓
24	1	2	1	0	1	1	✓	✓
25	1	2	1	0	1	1	✓	✓
26	1	2	1	0	1	1	✓	✓
27	1	2	1	0	1	1	✓	✓
28	1	2	1	0	1	1	✓	✓
29	1	2	1	0	1	1	✓	✓
30	1	2	1	0	1	1	✓	✓
31	1	2	1	0	1	1	✓	✓
32	1	2	1	0	1	1	✓	✓
33	1	2	1	0	1	1	✓	✓
34	1	2	1	0	1	1	✓	✓
35	1	2	1	0	1	1	✓	✓
36	1	2	1	0	1	1	✓	✓
37	1	2	1	0	1	1	✓	✓
38	1	2	1	0	1	1	✓	✓
39	1	2	1	0	1	1	✓	✓
40	1	2	1	0	1	1	✓	✓
41	1	2	1	0	1	1	✓	✓
42	1	2	1	0	1	1	✓	✓
43	1	2	1	0	1	1	✓	✓
44	1	2	1	0	1	1	✓	✓
45	1	2	1	0	1	1	✓	✓
46	1	2	1	0	1	1	✓	✓
47	1	2	1	0	1	1	✓	✓
48	1	2	1	0	1	1	✓	✓
49	1	2	1	0	1	1	✓	✓
50	1	2	1	0	1	1	✓	✓
51	1	2	1	0	1	1	✓	✓
52	1	2	1	0	1	1	✓	✓
53	1	2	1	0	1	1	✓	✓
54	1	2	1	0	1	1	✓	✓
55	1	2	1	0	1	1	✓	✓
56	1	2	1	0	1	1	✓	✓
57	1	2	1	0	1	1	✓	✓
58	1	2	1	0	1	1	✓	✓
59	1	2	1	0	1	1	✓	✓
60	1	2	1	0	1	1	✓	✓
61	1	2	1	0	1	1	✓	✓
62	1	2	1	0	1	1	✓	✓
63	1	2	1	0	1	1	✓	✓
64	1	2	1	0	1	1	✓	✓
65	1	2	1	0	1	1	✓	✓
66	1	2	1	0	1	1	✓	✓
67	1	2	1	0	1	1	✓	✓
68	1	2	1	0	1	1	✓	✓
69	1	2	1	0	1	1	✓	✓
70	1	2	1	0	1	1	✓	✓
71	1	2	1	0	1	1	✓	✓
72	1	2	1	0	1	1	✓	✓
73	1	2	1	0	1	1	✓	✓
74	1	2	1	0	1	1	✓	✓
75	1	2	1	0	1	1	✓	✓
76	1	2	1	0	1	1	✓	✓
77	1	2	1	0	1	1	✓	✓
78	1	2	1	0	1	1	✓	✓
79	1	2	1	0	1	1	✓	✓
80	1	2	1	0	1	1	✓	✓
81	1	2	1	0	1	1	✓	✓
82	1	2	1	0	1	1	✓	✓
83	1	2	1	0	1	1	✓	✓
84	1	2	1	0	1	1	✓	✓
85	1	2	1	0	1	1	✓	✓
86	1	2	1	0	1	1	✓	✓
87	1	2	1	0	1	1	✓	✓
88	1	2	1	0	1	1	✓	✓
89	1	2	1	0	1	1	✓	✓
90	1	2	1	0	1	1	✓	✓
91	1	2	1	0	1	1	✓	✓
92	1	2	1	0	1	1	✓	✓
93	1	2	1	0	1	1	✓	✓
94	1	2	1	0	1	1	✓	✓
95	1	2	1	0	1	1	✓	✓
96	1	2	1	0	1	1	✓	✓
97	1	2	1	0	1	1	✓	✓
98	1	2	1	0	1	1	✓	✓
99	1	2	1	0	1	1	✓	✓
100	1	2	1	0	1	1	✓	✓

[illegible]

Sample ID		ANZECC 2007 - Marine Value - 4%		ANZECC 2007 - Tribution -110		Sample Type									
Date Sampled	Base Sampled	LOR	Units	ANZECC 2007 - Marine Value - 4%	ANZECC 2007 - Tribution -110	Q404	TRP BLANK-1	TRP BLANK-2	Q408	TRP BLANK-3	TRP BLANK-4	Q412 28-9-11	TRPBLANK-5	TRPBLANK-6	
Sample Type				Field Blank	Trips Blank	Field Blank	Trips Blank	Trips Blank	Field Blank	Trips Blank	Trips Blank	Field Blank	Trips Blank	Trips Blank	
Analyte Fertilizants	1-2-Ethionomethane	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	1-2-Ethionomethane	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
Analyte Fertilizants	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
Analyte Fertilizants	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
	2,4-Dichlorophenol	5	µg/L	<5	-	<5	-	<5	<5	-	-	<5	-	-	
Analyte Fertilizants	2,4-Dichlorophenol	5													

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[illegible][illegible]







# DATA VALIDATION REPORT December 2011

<b>URS Project number:</b>	42213719	<b>Data verified by:</b>	Bek Agaard	<b>Date:</b> 28/04/2013
<b>Client:</b>	Darwin Waterfront Corporation	<b>Signed:</b>		
<b>Site:</b>	Waterfront Precinct	<b>Validation by:</b>	Tim Smith	<b>Date:</b> 2/5/2013
<b>URS Project Manager:</b>	Jacques van Rensburg	<b>Signed:</b>		
<b>Matrix type:</b>	Groundwater	<b>Project Manager:</b>		<b>Date:</b>
<b>No Primary samples:</b>	10	<b>Signed:</b>		
<b>Laboratory:</b>	ALS; SGS			
<b>Lab reference:</b>	EP1108921 PE063730			

## 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.
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:

- Matrix spikes (MS) were reported on anonymous samples for dissolved metals. As MS was conducted on anonymous samples, the reported MS recovery may not be considered representative of the URS field sample.
- Laboratory duplicate were not reported for total anions and cations and Total Petroleum Hydrocarbons (TPH)/Total Recoverable Hydrocarbons (TRH). The precision is considered acceptable based on the data quality assessment of intra- and inter-laboratory duplicates for TPH and TRH, and inter-laboratory for total anions and cations.
- The inter-laboratory duplicate was not analysed total anions and cations. The precision of this data is considered acceptable based on the data quality assessment of intra-laboratory field duplicates.

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.
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	<p>Samples were correctly preserved, chilled with ice and received at the laboratory at 12.4°C.</p> <p>While the recommended temperature for water samples on receipt is 4°C (<math>\pm 2^\circ\text{C}</math>) batch EP1108921 recorded on receipt above the recommended temperature; therefore, some losses through volatilisation may have occurred and sample concentrations for BTEXN and TPH C<sub>6</sub> – C<sub>9</sub> may be may be biased low.</p> <p>It should be noted that the groundwater temperature across the site throughout the year and at the time of sampling is generally <math>&gt;30^\circ\text{C}</math>. 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 24hr 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 are considered to be limited.</p>
Holding time compliance	<p>All samples were extracted and analysed within recommended holding times, with the following exceptions:</p> <p>Calcium, magnesium, potassium and sodium analysis was exceeded by 2 days. These analytes are used as one of a number of indicators of natural attenuation and sea water intrusion and as such will not affect the overall interpretation of results.</p>

## Data precision

Field duplicate RPDs	RPDs were within control limits.
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Interlaboratory Replicate (Field triplicate) RPDs	RPDs exceeded control limits for the following sample analysis. (Sample with higher reported concentrations is in bold). <ul style="list-style-type: none"><li>MW136_13/12/12 and <b>QC02_13/12/12</b> for Total Alkalinity (63%).</li></ul>												
Laboratory duplicate RPDs	RPDs were within control limits.												
<b>Data accuracy</b>													
Laboratory control spike recovery	<p>The laboratory control spike recoveries were within control limits.</p> <p>Laboratory control spike recoveries were not reported for total anions and cations. The accuracy of the data for total anions and cations is considered acceptable based on the presence of surrogate recoveries.</p>												
Matrix spike recovery	<p>The following recoveries were outside control limits and may affect data interpretation:</p> <table><tr><th>Sample</th><th>Analyte</th><th>Recovery (%)</th><th>LCL (%)</th><th>UCL (%)</th><th>Comment</th></tr><tr><td>MW138_13/12/11</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></table> <p>As the matrix spike issue was 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).</p> <p>Matrix spike recoveries were not reported for TPH C<sub>10</sub>-C<sub>36</sub>, total anions and cations, calcium, magnesium, potassium, sodium, total dissolved solids 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).</p>	Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment	MW138_13/12/11	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment								
MW138_13/12/11	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.								
Surrogate spike recovery	The surrogate spike recoveries were within control limits.												
<b>Blank monitoring</b>													
Equipment rinsate blank	<p>Concentrations were not detected above the LOR for all analytes tested, with the following exceptions:</p> <p>Concentrations of cadmium (0.0002 mg/L), lead (0.001 mg/L), manganese (0.001 mg/L) and zinc (0.006 mg/L) were detected above the LOR in the rinsate sample (collected on 14<sup>th</sup> of December 2011 (QCC01_14/12/12)).</p> <p>It should be noted that dedicated well sampling equipment was used at each location. The manganese concentration detected in the rinsate blank was well below manganese reported in primary samples and is not considered to affect interpretation of analytical results.</p> <p>Concentrations of cadmium, lead and zinc were not detected in the field blank; potential for cross-contamination to have occurred between the five locations sampled on the 14<sup>th</sup> of December 2011 is not considered likely. Rinsate water used may have been from a glass bottle and may have contained trace metals impurities. Analytical results from rinsate blank analysis are close to laboratory LOR and are not considered to affect primary results or interpretation of analytical data.</p>												
Field blank	Manganese (0.002 mg/L) was detected above the LOR in the field blank sample (collected on 14 <sup>th</sup> of December 2011 (QCB01_14/12/12). Rinsate water used may have been from a glass bottle and may have contained trace metals impurities. Analytical results from field blank analysis are close to laboratory LOR and are not considered to affect primary results or interpretation of analytical data.												
Trip blank	Concentrations of all analytes were reported below the LOR.												
Method blank	Concentrations of all analytes were reported below the LOR.												
<b>Chromatograms</b>	N/A												
Other observations	None reported												

Site Name: Wascott Precinct  
Project No. 4221719  
Project Manager Tim Smith  
Water ALS/SGS  
Batch File Number EP110821/PE063730

Analytical Parameter	Number of Tests Requested	Number of Tests Reported	Holding Times (a)	Limits of Reporting (b)	Field Blank (1 per day)		Pinstate 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
Arsenic	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Barium	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Benzene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Boron	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Bromine	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Chlorine	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Cadmium	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Copper	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Lead	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Manganese	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Nickel	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Nitrogen	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Vanadium	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Zinc	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Mercury	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	-	-
Calcium	11	11	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	2	1	0	1	2	-	-
Magnesium	11	11	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	2	1	0	1	2	-	-
Potassium	11	11	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	2	1	0	1	2	-	-
Sodium	11	11	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	2	1	0	1	2	-	-
Total Arsenic	10	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	0	1	0	1	0	1	0	-	-
Total Cadmium	10	10	✓	✓	0	0	0	0	0	0	1	0	1	1	1	0	1	0	1	0	1	0	-	-
Total Chromium	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	0	1	1	-	-
Total Copper	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	0	1	1	-	-
Total Lead	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	0	1	1	-	-
Total Manganese	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	0	1	1	-	-
Total Nickel	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	0	1	1	-	-
Total Nitrogen	13	13	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
Total Vanadium	13	13	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
Total Zinc	13	13	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
AsO4 - C41 Fraction	13	13	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
AsO4 - C41 Fraction	13	13	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
AsO4 - C41 Fraction	14	14	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
AsO4 - C28 Fraction	14	14	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
AsO4 - C28 Fraction	14	14	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
AsO4 - C28 Fraction	14	14	✓	✓	1	1	1	1	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
Chloride	11	11	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
Chloride	11	11	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	0	1	0	1	2	-	-
Chloride	15	15	✓	✓	1	1	1	1	2	2	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
Chloride	15	15	✓	✓	1	1	1	1	2	2	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
Chloride	16	16	✓	✓	1	1	1	1	1	1	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
Benzene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
Ethylbenzene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
m,p-Xylene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
p-Xylene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
Naphthalene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
o-Xylene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
Toluene	14	14	✓	✓	1	1	1	1	0	0	1	1	1	1	1	1	1	2	1	1	1	1	✓	✓
Toluene	10	10	✓	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	3	1	1	1	2	✓	✓
Surrogate AsO4 - Tridecenoic	11	11	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	2	1	0	1	1	-	-

Groundwater Analytical Results - Field Blanks  
December 2011

Sample ID
Date Sampled
Sample Type

QCA01	QCA02	QCB01	QCC01
13/12/2011	14/12/2011	14/12/2011	14/12/2011
Trip Blank	Trip Blank	Field Blank	Rinsate Blank

Analyte	LOR	Units	ANZECC 2000 - Marine Water - 95%	ANZECC 2000 - Irrigation - LTU				
<b>Total Petroleum Hydrocarbons</b>								
C6-C9 fraction	20	µg/L			<20	<20	<20	<20
C10-C14 fraction	50	µg/L			-	-	<50	<50
C15-C28 fraction	100	µg/L			-	-	<100	<100
C29-C36 fraction	50	µg/L			-	-	<50	<50
C10 - C36 Fraction (sum)	50	µg/L			-	-	<50	<50
<b>BTEX Compounds</b>								
Benzene	1	µg/L	700		-	-	<1	<1
Toluene	2	µg/L			-	-	<2	<2
Ethylbenzene	2	µg/L			-	-	<2	<2
m&p-Xylene	2	µg/L			-	-	<2	<2
o-Xylene	2	µg/L			-	-	<2	<2
Total Xylenes	2	µg/L			-	-	<2	<2
Sum of BTEX	1	µg/L			-	-	<1	<1
Naphthalene	5	µg/L	70		-	-	<5	<5
<b>Metals (Dissolved)</b>								
Arsenic	0.001	mg/L		0.1	-	-	<0.001	<0.001
Cadmium	0.0001	mg/L	0.0055	0.01	-	-	<0.0001	<b>0.0002</b>
Chromium	0.001	mg/L	0.0044	0.1	-	-	<0.001	<0.001
Copper	0.001	mg/L	0.0013	0.2	-	-	<0.001	<0.001
Lead	0.001	mg/L	0.0044	2	-	-	<0.001	<b>0.001</b>
Manganese	0.001	mg/L		0.2	-	-	<b>0.002</b>	<b>0.001</b>
Mercury	0.0001	mg/L	0.0004	0.002	-	-	-	-
Nickel	0.001	mg/L	0.07	0.2	-	-	<0.001	<0.001
Vanadium	0.01	mg/L	0.1	0.1	-	-	<0.01	<0.01
Zinc	0.005	mg/L	0.015	2	-	-	<0.005	<b>0.006</b>
<b>Major Ions</b>								
Calcium	1	mg/L			-	-	-	<1
Magnesium	1	mg/L			-	-	-	<1
Potassium	1	mg/L			-	-	-	<1
Sodium	1	mg/L			-	-	-	<1
Sulfate as SO4 - Turbidimetric	1	mg/L			-	-	-	<1
<b>Miscellaneous Compounds</b>								
Chloride	1	mg/L			-	-	-	<1
Ionic Balance	0.01	%			-	-	-	<0.01
Total Anions	0.01	meq/l			-	-	-	<0.01
Total Cations	0.01	meq/l			-	-	-	<0.01
Total Dissolved Solids	5	mg/L			-	-	-	<5
<b>Alkalinity</b>								
Hydroxide Alkalinity as CaCO3	1	mg/L			-	-	-	<1
Carbonate Alkalinity as CaCO3	1	mg/L			-	-	-	<1
Total Alkalinity	1	mg/L			-	-	-	<1
Bicarbonate Alkalinity as CaCO3	1	mg/L			-	-	-	<1
<b>Total Recoverable Hydrocarbons</b>								
>C10 - C16 Fraction	100	µg/L			-	-	<100	<100
>C10 - C40 Fraction (sum)	100	µg/L			-	-	<100	<100
>C16 - C34 Fraction	100	µg/L			-	-	<100	<100
>C34 - C40 Fraction	100	µg/L			-	-	<100	<100
C6 - C10 Fraction	20	µg/L			<20	<20	<20	<20
C6 - C10 Fraction minus BTEX (F1)	20	µg/L			-	-	<20	<20

Legend:

Exceeds the ANZECC/ARMCANZ, 2000, Trigger values for marine water ecosystems - Level of protection 95% species

Exceeds the ANZECC/ARMCANZ, 2000, Irrigation

- Not analysed / not calculated

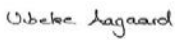

\* LOR Exceeds Guideline Trigger Value

10% checked  
1/5/2013 BA

Table 777  
Groundwater Analytical Results - Analysis Results  
Waterfront

Location		MW108	MW115	MW116	MW121	MW125	MW127	MW128	MW134	MW135	MW136	MW136	MW136
Sample ID		MW108	MW115	MW116	MW121	MW125	MW127	MW128	MW134	MW135	MW136	MW136	MW136
Date Sampled		13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011	13/12/2011
Sample Type		Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample	Primary Sample
Primary Sample ID		MW108	MW115	MW116	MW121	MW125	MW127	MW128	MW134	MW135	MW136	MW136	MW136
Area		-	-	-	-	-	-	-	-	-	-	-	-
Analyte		LOR	Units	ANZECC 2000 - Marine Life	ANZECC 2000 - Ingestion - L10								
Total Petroleum Hydrocarbons		20	µg/L										
C6-C9 fraction		<20		<20	<20	<20	<20	<20	<20	<20	<20	<20	<20
C10-C14 fraction		130	µg/L	<50	<50	<50	80	50	50	<50	<50	<50	<50
C15-C28 fraction		1480	µg/L	<200	<100	<100	890	160	890	<100	480	890	890
C29-C36 fraction		300	µg/L	<50	<50	<50	160	<50	160	<50	160	160	160
C37-C44 fraction		1910	µg/L	<50	<50	<50	1120	260	1040	<50	640	950	950
BTEX Compounds													
Benzene		1	µg/L	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Toluene		2	µg/L	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2
Ethylbenzene		2	µg/L	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2
m,p-Xylene		2	µg/L	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2
o-Xylene		2	µg/L	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2
Total Xylenes		2	µg/L	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2
Sum of BTEX		1	µg/L	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Naphthalene		5	µg/L	<5	<5	<5	<5	<5	<5	<5	<5	<5	<5
Metals (Dissolved)													
Cadmium		0.001	µg/L	<0.001	<0.001	<0.001	0.002	<0.001	0.002	<0.001	<0.001	<0.001	<0.001
Chromium		0.001	µg/L	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Chromium		0.001	µg/L	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Copper		0.001	µg/L	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Lead		0.001	µg/L	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Manganese		0.001	µg/L	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Mercury		0.001	µg/L	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Nickel		0.001	µg/L	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001
Vanadium		0.01	µg/L	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01	<0.01
Zinc		0.005	µg/L	0.01	0.009	0.007	0.012	0.01	<0.005	<0.005	<0.005	<0.005	<0.005
Major Ions													
Calcium		1	mg/L	164	72	29	45	-	-	155	70	70	70
Magnesium		1	mg/L	92	81	29	20	-	-	49	52	52	52
Potassium		1	mg/L	48	415	26	17	-	-	28	31	31	37
Sodium		1	mg/L	892	10000	200	455	-	-	271	812	837	890
Sulfate as SO <sub>4</sub> - Tetrathionate		1	mg/L	4	633	16	4	-	-	200	63	63	69
Miscellaneous Compounds													
Carbonate Alkalinity as CO <sub>3</sub>		1	mg/L	-	-	-	-	-	-	-	-	-	<1
Chloride		1	mg/L	1490	784	328	344	-	-	432	170	1160	1360
Total Hardness		0.01	%	4.63	2.76	1.81	2.85	-	-	0.89	1.89	2.35	-
Total Anions		0.01	mg/L	46	253	133	142	-	-	21.8	42.8	45.1	-
Total Cations		0.01	mg/L	54.5	636	31	13.6	-	-	21.3	42.8	45.1	-
Total Dissolved Solids		5	mg/L	2970	32600	828	783	-	-	1490	2010	2550	2460
Alkalinity													
Hydroxide Alkalinity as CaCO <sub>3</sub>		1	mg/L	<1	<1	<1	<1	-	-	<1	<1	<1	-
Carbonate Alkalinity as CaCO <sub>3</sub>		1	mg/L	<1	<1	<1	<1	-	-	<1	<1	<1	-
Total Alkalinity		1	mg/L	376	737	247	241	-	-	374	308	401	760
Bicarbonate Alkalinity as CaCO <sub>3</sub>		1	mg/L	376	737	247	241	-	-	374	308	401	-
Total Recoverable Hydrocarbons													
C6-C9 Fraction		100	µg/L	<100	<100	<100	240	<100	250	<100	<100	110	-
C10-C14 Fraction (sum)		100	µg/L	<100	<100	<100	1480	160	1070	<100	890	1320	-
C15-C28 Fraction		100	µg/L	<100	<100	<100	1020	160	820	<100	700	1060	-
C29-C36 Fraction		100	µg/L	<100	<100	<100	230	<100	200	<100	200	360	-
C37-C44 Fraction		100	µg/L	<100	<100	<100	<20	<100	<20	<20	<20	<20	-
C6-C10 Fraction minus BTEX (F1)		20	µg/L	<20	<20	<20	<20	<20	<20	<20	<20	<20	-
F10 C10-C14		50	µg/L	-	-	-	-	-	-	-	-	140	-
F10 C15-C28		200	µg/L	-	-	-	-	-	-	-	-	480	-
F10 C29-C36		200	µg/L	-	-	-	-	-	-	-	-	<200	-
F10 C4-C9		40	µg/L	-	-	-	-	-	-	-	-	<40	-

# DATA VALIDATION REPORT\_March 2012

URS Project number:	42213719	Data verified by:	Bek Aagaard	Date: 26/04/2013
Client:	Darwin Waterfront Corporation			
Site:	Waterfront Precinct	Signed:		
URS Project Manager:	Jacques Van Rensburg	Validation by:	Tim Smith	Date: 3/5/2013
Matrix type:	Groundwater			
No Primary samples:	23			
Laboratory:	ALS; SGS	Signed:		
Lab reference:	ES1205485 & SE106212; EP1201888; ES1205892 & SE106317	Project Manager:		Date:
		Signed:		

## 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, except for no trip blank, field blank and rinsate blank was not taken for batch ES1205892; hence the potential for cross-contamination has not been assessed for that batch, this is not considered to affect data quality for interpretive purposes. Potential data considerations are discussed below.
Frequency of laboratory QC	<p>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:</p> <ul style="list-style-type: none"> <li>Matrix spikes (MS) were reported on anonymous samples for all analytes. Where MS were conducted on anonymous samples, the reported MS recovery may not be considered representative of the URS field sample.</li> </ul> <p><b>Batch ES1205485</b></p> <ul style="list-style-type: none"> <li>Laboratory duplicate were not reported for total anions and cations and ferric iron. The precision of these analytes have been considered acceptable based on the presence intra-laboratory duplicates were available and acceptable.</li> <li>The inter-laboratory duplicate was not analysed for total anions and cations. The precision of this data can be assessed as acceptable based on the presence of intra-laboratory field duplicates for these analytes.</li> </ul> <p><b>Batch EP1201888</b></p> <ul style="list-style-type: none"> <li>Laboratory duplicate were not reported for total anions and cations, TPH/Total Recoverable Hydrocarbons (TRH), PAHs and ferric iron. The precision of these analytes been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable from other batches within the complete data set.</li> </ul> <p><b>Batch ES1205892</b></p> <ul style="list-style-type: none"> <li>Laboratory duplicate were not reported for total anions and cations, TPH/TRH, PAHs and ferric iron. The precision of these analytes have been considered acceptable based on the presence intra- and inter-laboratory duplicates were available and acceptable (where applicable).</li> <li>The inter-laboratory duplicate was not analysed for total anions and cations, volatiles and ferric iron. The precision of this data can be assessed as acceptable based on the presence of intra-laboratory field duplicates for these analytes.</li> </ul>
Tests requested/reported	Samples were analysed and reported as requested on the COC.
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"> <li>Protocol LOR for anthracene (1 µg/L) above the adopted ILs (0.4 µg/L)</li> <li>Protocol LOR for benzo(a)pyrene (0.5 µg/L) above the adopted ILs (0.2 µg/L)</li> </ul> <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"> <li>Protocol LOR exceeds the guideline trigger value for chromium (0.0044 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for copper (0.0013 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for lead (0.0044 mg/L)</li> <li>Protocol LOR exceeds the guideline trigger value for zinc (0.015 mg/L)</li> </ul> <p>The limits of reporting exceeded the guideline trigger levels for chromium copper, lead and zinc. As a</p>

	<p>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>
<b>Sample management</b>	
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 4.1°C, 10.0°C and 9.1°C for batches ES1205485, EP1201888 and ES1205892, respectively.</p> <p>Samples from batches EP1201888 and ES1205892 were received above the recommended temperature range; therefore, some losses through volatilisation may have occurred and sample concentrations for BTEXN and TPH C<sub>6</sub> – C<sub>9</sub> may be biased low.</p> <p>It should be noted that the groundwater temperature across the site throughout the year and at the time of sampling is generally &gt;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 24hr 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 are considered to be limited.</p>
Holding time compliance	<p>All samples were extracted and analysed within recommended holding times, with the following exceptions:</p> <p><b><u>Batch ES1205485</u></b>            Ferrous iron analysis was exceeded by 2 days, nitrite as N by 1 day, sulphide as S<sup>2-</sup> by 7 days and sulphite as SO<sub>3</sub><sup>2-</sup> by 3 days. These analytes are used as indicators of natural attenuation and as such will not affect the overall interpretation of results.</p> <p><b><u>Batch EP1201888</u></b>            PAHs extraction was exceed by 1 day, TPH by 1 day and TRH by 1 day. Total dissolved solids analysis was exceeded by 4 days and sulphite as SO<sub>3</sub><sup>2-</sup> by 3 days,. These analytes are used as indicators of natural attenuation and as such will not affect the overall interpretation of results.</p> <p><b><u>Batch ES1205892</u></b>            Sulphide as S<sup>2-</sup> analysis was exceeded by 2 days. This analyte is used as one of a number of indicators of natural attenuation and as such will not affect the overall interpretation of results.</p>
<b>Data precision</b>	
Field duplicate RPDs	<p>RPDs exceeded control limits for the following sample analysis. (Sample with higher reported concentrations is in bold).</p> <ul style="list-style-type: none"> <li>• <b>MW103_12/03/12</b> and QC03_12/03/12 for Copper (32%).</li> <li>• <b>MW103_12/03/12</b> and QC03_12/03/12 for Manganese (49%).</li> <li>• <b>MW103_12/03/12</b> and QC03_12/03/12 for Zinc (31%).</li> </ul> <p>For most samples, reported concentrations were either well above or below the adopted ILs; however, the apparent lack of precision should be taken into consideration when evaluating individual results, particularly where concentrations are reported close to the adopted investigation level.</p>
Inter-laboratory Replicate (Field triplicate) RPDs	<p>RPDs exceeded control limits for the following sample analysis. (Samples with higher reported concentrations is in bold).</p> <ul style="list-style-type: none"> <li>• MW128_07/03/12 and <b>QC02_07/03/12</b> for Total Alkalinity (31%).</li> <li>• MW128_07/03/12 and <b>QC02_07/03/12</b> for Sodium (67%).</li> <li>• <b>MW103_12/03/12</b> and QC04_12/03/12 for Total Alkalinity (179%).</li> <li>• <b>MW103_12/03/12</b> and QC04_12/03/12 for Sodium (59%).</li> <li>• <b>MW103_12/03/12</b> and QC04_12/03/12 for Manganese (179%).</li> <li>• <b>MW103_12/03/12</b> and QC04_12/03/12 for Zinc (39%).</li> </ul> <p>For most samples, reported concentrations were either well above or below the adopted ILs; however, the apparent lack of precision should be taken into consideration when evaluating individual results, particularly where concentrations are reported close to the adopted investigation level.</p>
Laboratory duplicate RPDs	<p>RPDs were within control limits.</p>

## Data accuracy

### Laboratory control spike recovery

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

Batch	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
ES1205485	Chromium	80.8	91	111	Recovery less than lower control limit
ES1205485	Copper	80.8	87	111	Recovery less than lower control limit
ES1205485	Lead	77.9	90	110	Recovery less than lower control limit
ES1205485	Manganese	80.4	87	113	Recovery less than lower control limit
ES1205485	Nickel	82.2	89	109	Recovery less than lower control limit
ES1205485	Vanadium	73.6	91	109	Recovery less than lower control limit
ES1205485	Zinc	82.7	85	116	Recovery less than lower control limit
ES1205485	Mercury	81.6	86	116	Recovery less than lower control limit
ES1205892	Chromium	90.3	91	111	Recovery less than lower control limit
ES1205892	Vanadium	88.9	91	109	Recovery less than lower control limit

For LCS recoveries reported less than the lower control limit, the potential exists for reported concentrations to be biased low. Care should be taken with interpreting results for these analytes close to the adopted guidelines.

LCS recoveries were not reported for total anions and cations and ferric for all batches. And alkalinity for batches ES1205485 and ES125892. 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).

### Matrix spike recovery

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

Sample	Analyte	Recovery (%)	LCL (%)	UCL (%)	Comment
Anonymous	Sulphate as SO <sub>4</sub> - Turbidimetric	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW118_6/3/12	Manganese	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.
MW135_7/3/12	Manganese	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW102_8/3/12	Sulphate as SO <sub>4</sub> - Turbidimetric	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.
MW102_8/3/12	Chloride	Not determined	-	-	MS recovery not determined, background level greater than or equal to 4x spike level.

As the matrix spike issue was 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).

#### **Batch ES1205485**

Matrix spike recoveries were not reported for alkalinity, total anions and cations, calcium, magnesium, potassium, sodium, total dissolved solids, ferric iron, sulphur as S, sulphite as SO<sub>3</sub><sup>2-</sup>, trihalogenated methanes, oxygenated compounds, monocyclic aromatic hydrocarbons (MAHs) and halogenated aliphatics. 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 EP1201888**

Matrix spike recoveries were not reported for TPH C<sub>9</sub>-C<sub>36</sub>/TRH C<sub>10</sub>-C<sub>40</sub>, PAHs, alkalinity, total anions and cations, calcium, magnesium, potassium, sodium, total dissolved solids, ferric iron, sulphur as S, sulphide as S<sup>2-</sup>, sulphite as SO<sub>3</sub><sup>2-</sup>, carbon disulphide, fumigants, trihalogenated methanes and oxygenated compounds. 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><b>Batch ES1205892</b></p> <p>Matrix spike recoveries were not reported for TPH C<sub>9</sub>-C<sub>36</sub>/TRH C<sub>10</sub>-C<sub>40</sub>, PAHs, alkalinity, total anions and cations, calcium, magnesium, potassium, sodium, total dissolved solids, ferric iron, sulphur as S, sulphite as SO<sub>3</sub><sup>2-</sup>, trihalogenated methanes, fumigants, oxygenated compounds, monocyclic aromatic hydrocarbons (MAHs) and halogenated aliphatics. 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	The surrogate spike recoveries were within control limits.
<b>Blank monitoring</b>	
Equipment rinsate blank	<p>Concentrations were not detected above the LOR for all analytes tested, with the following exceptions:</p> <p>Concentrations of chromium (0.051 mg/L), manganese (0.006 mg/L), and nickel (0.025 mg/L) were detected above the LOR in the rinsate sample (collected on 8th of March 2012 (QCC01_80312) in batch EP1201888. Concentrations of chromium, manganese and nickel were not detected in the field blank; hence, potential exists for cross-contamination to have occurred between the eight locations sampled on the 8<sup>th</sup> of March 2012.</p> <p>Rinsate blanks were not collected for batch ES1205892 12<sup>th</sup> March 2012. Given that rinsate blanks were collected for all other sampling days, analytical results except those listed above were &lt;LOR and dedicated sampling equipment was used at each location this is not considered to affect interpretation of analytical results.</p> <p>It should be noted that dedicated well sampling equipment and disposable gloves were used at each location. This is not considered to affect interpretation of analytical results</p>
Field blank	<p>Concentrations of all analytes were reported below the LOR, with the following exceptions:</p> <p>Field blanks were not collected for batch ES1205892 12<sup>th</sup> March 2012. Given that field blanks were collected for all other sampling days, analytical results were &lt;LOR and dedicated sampling equipment was used at each location this is not considered to affect the interpretation of results.</p>
Trip blank	<p>Concentrations of all analytes were reported below the LOR, with the following exceptions:</p> <p>As no volatile or semi-volatile COPCs were reported above LOR in primary samples from batch ES1205892 12<sup>th</sup> March 2012, it is not considered that cross contamination during sample transport is likely.</p>
Method blank	Concentrations of all analytes were reported below the LOR.
<b>Chromatograms</b>	N/A
<b>Other observations</b>	<p><b>Batch ES1205485</b></p> <ul style="list-style-type: none"> <li>EG020: LCS recoveries for some elements fall outside ALS Dynamic Control Limit. However, they are within the acceptance criteria based on ALS DQO. No further action is required.</li> <li>EG020: Samples were diluted and rerun due to matrix interference and LOR's have been raised accordingly. (High sample salinity).</li> <li>EG035T: LCS recovery for Mercury falls outside ALS Dynamic Control Limit. However, it is within the acceptance criteria based on ALS DQO. No further action is required.</li> <li>EK085: Sulphide ;:sample IDs "MW118" and "MW117" invalidated due to insufficient sample.</li> <li>EP080: Level of reporting raised for toluene due to ambient background levels in the laboratory.</li> </ul> <p><b>Batch ES1205892</b></p> <ul style="list-style-type: none"> <li>EP080: Level of reporting raised for toluene due to ambient background levels in the laboratory.</li> </ul>

[illegible][illegible]

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Primary Samples	Holding Times (a)	Limits of Reporting (b)	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	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number	Number
						Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported	Required	Reported
VOLATILES ANALYSIS/ALS/EP04E	1,1,2-Tetrachloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,1,1-Trichloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,1,2,2-Tetrachloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,1,2-Trichloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,1-Dichloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,1-Dichloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,1-Dichloropropane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,2,3-Trichloropropane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,2-Dibromo-3-chloropropane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,2-Dichloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
VOLATILES ANALYSIS/ALS/EP04C	1,3-Dichloropropane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Bromomethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Carbon Tetrachloride	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Chloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Chloromethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Chloromethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	cis-1,2-Dichloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	dis-1,4-Dichloro-2-butene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Dibromomethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Dichlorodifluoromethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
VOLATILES ANALYSIS/ALS/EP04G	Hexachlorobutadiene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Iodomethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Pentachloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Pentachloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Tetrachloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	trans-1,2-Dichloroethene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	trans-1,4-Dichloro-2-butene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Trichloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Trichlorofluoromethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Vinyl chloride	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
VOLATILES ANALYSIS/ALS/EP04B	Carbon disulfide	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Bromodichloromethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Bromomethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Chloroform	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Dibromodichloromethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Dibromodichloromethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	2-Butanone (MEK)	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	2-Hexanone (MIBK)	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	4-Methyl-2-pentanone (MIBK)	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Vinyl Acetate	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
VOLATILES ANALYSIS/ALS/EP04F	1,2,3-Trichlorobenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,2,4-Trichlorobenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,2-Dichlorobenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,3-Dichlorobenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,4-Dichlorobenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	2-Chloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	4-Chloroethane	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	4-Chlorotoluene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Bromobenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Chlorobenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
VOLATILES ANALYSIS/ALS/EP04A	1,2,4-Trimethylbenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	1,3,5-Trimethylbenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Isopropylbenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	n-Butylbenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	n-Propylbenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	p-Isopropyltoluene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	sec-Butylbenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	Styrene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			
	tert-Butylbenzene	5	5	2	✓	1	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	1	1	1	✓	✓			

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Primary Samples	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
VOLATILES ANALYSIS/ALS/SEP0740	1,2-Dibromochloroethane (EDB)	5	5	2	✓	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	0	1	1	✓	✓
	2,2-Dichloropropane	5	5	2	✓	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	0	1	1	✓	✓
	2,2-Dichloropropane	5	5	2	✓	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	0	1	1	✓	✓
	98-13-Dichloropropylene	5	5	2	✓	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	0	1	1	✓	✓
	trans-1,3-Dichloropropylene*	5	5	2	✓	1	1	1	1	1	1	1	1	0	0	0	0	1	1	1	0	1	1	✓	✓
SEMIVOLATILES ANALYSIS/ALS/SEP060071	2-C10 - C14 Fraction	10	10	7	✓	1	1	1	1	1	0	1	1	1	1	0	0	1	1	1	1	1	1	-	-
	C16 - C24 Fraction	10	10	7	✓	1	1	1	1	1	0	1	1	1	1	0	0	1	1	1	1	1	1	-	-
	C24 - C40 Fraction	11	11	7	✓	1	1	1	1	1	0	1	1	1	1	1	1	1	1	1	1	1	1	-	-
	C18 - C28 Fraction	11	11	7	✓	1	1	1	1	1	0	1	1	1	1	1	1	1	1	1	1	1	1	-	-
	C26 - C36 Fraction	11	11	7	✓	1	1	1	1	1	0	1	1	1	1	1	1	1	1	1	1	1	1	-	-
METALS ANALYSIS/ALS/EG050F	Mercury	9	9	7	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	-	-
	Sulfate as SO3 2-	9	9	7	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	-	-
	Sulfate as SO3 2-	7	7	5	✗	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	1	1	1	-	-
	Calcium	9	9	7	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	1	1	0	1	2	-	-
	Magnesium	9	9	7	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	1	1	0	1	2	-	-
METALS ANALYSIS/ALS/EG050F	Potassium	9	9	7	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	1	1	0	1	2	-	-
	Sodium	9	9	7	✓	0	0	0	0	0	0	1	2	1	1	1	1	1	1	1	0	1	2	-	-

NOTES:

- (a) Holding times are either project guideline limits.  
✗ Holding times exceed project guideline limits.  
(b) ✓ Limiting values exceed project guideline limits.  
✓ Limiting values are within project specifications.  
\* LODs do not comply with project specifications.

NA - Not Applicable

Analytical Method	Analytical Parameter	Number of Tests Requested	Number of 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)		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	Number Reported	Number Required	Number Reported	Number Required	Number Reported
METALS ANALYSIS/SE005F	Arsenic	10	10	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Sodium	10	10	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Chromium	10	10	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Copper	10	10	8	✓	✗	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Lead	10	10	8	✓	✗	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Manganese	10	10	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Nickel	10	10	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Vanadium	10	10	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Zinc	10	10	9	✓	✗	1	1	0	0	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Mercury	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	1	1	1	-	-
SEMI-VOLATILES ANALYSIS/SE003	Substrate as S	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	0	1	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	0	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	0	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	0	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	0	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	0	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	0	1	1	-	-
	METALS ANALYSIS/SE005F	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	0	1	1	-	-
NON-METALS INSTRUMENTALS/SE04S	Chloride	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	1	1	1	1	1	-	-
	Substrate as Se02.2	10	10	10	✗	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	1	-	-
	Substrate as Se04.1 - Cyanide-free	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	1	-	-
NON-METALS INSTRUMENTALS/SE05S	Total Arsenic	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	0	-	-
	Total Cadmium	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	0	1	0	1	0	-	-
	Se010 - C10 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
	Se016 - C16 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
	Se024 - C24 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
	Se030 - C30 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
	Se036 - C36 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
	Se042 - C42 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
	Se048 - C48 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
	Se054 - C54 Fraction	12	12	10	✗	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	0	1	0	1	1	-	-
NON-METALS INSTRUMENTALS/SE003P	Se059 - C59 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se065 - C65 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se071 - C71 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se077 - C77 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se083 - C83 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se089 - C89 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se095 - C95 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se101 - C101 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se107 - C107 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
	Se113 - C113 Fraction	10	10	10	✓	✓	0	0	0	0	0	0	1	1	0	0	0	0	1	2	1	0	1	2	✓	✓
VOLATILES/SE006	Dimethylsiloxane	12	12	10	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	2	1	1	2	✓	✓	
	Ethylbenzene	12	12	10	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	2	1	0	2	✓	✓	
	m,p-Xylene	12	12	10	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	2	1	0	2	✓	✓	
	o-Xylene	12	12	8	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	2	1	0	2	✓	✓	
	p-Xylene	12	12	10	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	2	1	0	2	✓	✓	
	Toluene	12	12	10	✓	✓	1	1	1	1	0	0	1	1	0	0	0	0	2	2	1	1	2	✓	✓	
	Se0 - C10 Fraction	12	12	10	✓	✓	1	1	1	1	1	1	1	1	0	0	0	0	2	2	1	1	1	✓	✓	
	Se0 - C10 Fraction	12	12	10	✓	✓	1	1	1	1	1	1	1	1	0	0	0	0	2	2	1	1	1	✓	✓	
	Se0 - C10 Fraction	12	12	10	✓	✓	1	1	1	1	1															

[illegible]



Analytical Method	Analytical Parameter	Number of Tests Requested	Number of Primary Samples	Holding Time(s)	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)		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	
INSTRUMENT LABORATORY (NON-METALS) SEC03D8G5 METALS ANALYSIS SEC03F	Chloride	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Calcium	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Magnesium	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Potassium	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Sodium	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D7P SEM/VOLATILES ANALYSIS SEC03D7P1	Barium/Arsenic Accuracy at GC03	7	7	✓	✓	0	0	0	0	0	0	0	1	1	1	1	1	1	1	0	-	
	Carbonate Accuracy at GC03	7	7	✓	✓	0	0	0	0	0	0	0	1	1	1	1	1	1	1	0	-	
	Peroxide Accuracy at GC03	7	7	✓	✓	0	0	0	0	0	0	0	1	1	1	1	1	1	1	0	-	
	Total Alkalinity at GC03	7	7	✓	✓	0	0	0	0	0	0	0	1	1	1	1	1	1	1	1	-	
	Total Alkalinity at GC03	7	7	✓	✓	0	0	0	0	0	0	0	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Total Alkalinity	8	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Total Calcium	8	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Total Chloride	8	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	8	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Total Organic Carbon	8	8	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
	Surfactant at SEC	7	7	✓	✓	0	0	0	0	0	0	1	1	1	1	1	1	1	1	1	-	
INSTRUMENT LABORATORY (NON-METALS) SEC03D5 SEM/VOLATILES ANALYSIS SEC03D5F	Surfactant at SEC	7	7	✓	✓	0	0	0														





Groundwater Analytical Results - Field Blanks  
March 2012

Sample ID
Date Sampled
Sample Type

QCA01	QCB01	QCC01	QCA01_080312	QCB01_080312	QCC01_080312
7/03/2012	7/03/2012	7/03/2012	9/03/2012	9/03/2012	9/03/2012
Trip Blank	Field Blank	Rinsate Blank	Trip Blank	Field Blank	Rinsate Blank

Analyte	LOR	Units	ANZECC 2000 - Marine Water - 95%	ANZECC 2000 - Irrigation - LTU
<b>Total Petroleum Hydrocarbons</b>				
C6-C9 fraction	20	µg/L		
C10-C14 fraction	50	µg/L		
C15-C28 fraction	100	µg/L		
C29-C36 fraction	50	µg/L		
C10 - C36 Fraction (sum)	50	µg/L		
C10-C36 fraction		µg/L		
C6-C36 fraction		µg/L		
<b>BTEX Compounds</b>				
Benzene	1	µg/L	700	
Toluene	2	µg/L		
Ethylbenzene	2	µg/L		
m&p-Xylene	2	µg/L		
o-Xylene	2	µg/L		
Total Xylenes		µg/L		
Sum of BTEX	1	µg/L		
Total BTEX		µg/L		
Naphthalene	1	µg/L	70	
<b>Metals (Dissolved)</b>				
Arsenic	0.001	mg/L		0.1
Cadmium	0.001	mg/L	0.0055	0.01
Chromium	0.001	mg/L	0.0044	0.1
Copper	0.001	mg/L	0.0013	0.2
Lead	0.001	mg/L	0.0044	2
Manganese	0.001	mg/L		0.2
Nickel	0.001	mg/L	0.07	0.2
Vanadium	0.01	mg/L	0.1	0.1
Zinc	0.005	mg/L	0.015	2
			<20	<20
			<50	<50
			<100	<100
			<50	<50
			<50	<50
			<200	<200
			<220	<220
			<1	<1
			<5	<2
			<2	<2
			<2	<2
			<2	<2
			<4	<4
			<1	<1
			<12	<9
			-	<1
			-	-
			-	<0.001
			-	<0.0001
			-	<0.001
			-	<0.001
			-	<0.001
			-	<0.001
			-	<0.01
			-	<0.005

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7/03/2012	7/03/2012	7/03/2012	9/03/2012	9/03/2012	9/03/2012
Trip Blank	Field Blank	Rinsate Blank	Trip Blank	Field Blank	Rinsate Blank

Analyte	LOR	Units	ANZECC 2000 - Marine Water - 95%	ANZECC 2000 - Irrigation - LTU
<b>Polynuclear Aromatic Hydrocarbons</b>				
Naphthalene	7	µg/L	70	
Acenaphthylene	1	µg/L		
Acenaphthene	1	µg/L		
Fluorene	1	µg/L		
Phenanthrene	1	µg/L		
Anthracene	1	µg/L		
Fluoranthene	1	µg/L		
Pyrene	1	µg/L		
Benz(a)anthracene	1	µg/L		
Chrysene	1	µg/L		
Benzo(b)fluoranthene	1	µg/L		
Benzo(k)fluoranthene	1	µg/L		
Benzo(a)pyrene	0.5	µg/L		
Indeno(1,2,3-cd)pyrene	1	µg/L		
Dibenz(a,h)anthracene	1	µg/L		
Benzo(g,h,i)perylene	1	µg/L		
Sum of polycyclic aromatic hydrocarbons	0.5	µg/L		
<b>Chlorinated Aromatic Compounds</b>				
1,2,3-Trichlorobenzene	5	µg/L		
1,2,4-Trichlorobenzene	5	µg/L	80	
1,2,4-Trimethylbenzene	5	µg/L		
1,2-Dichlorobenzene	5	µg/L		
1,3-Dichlorobenzene	5	µg/L		
1,4-Dichlorobenzene	5	µg/L		
2-Chlorotoluene	5	µg/L		
4-Chlorotoluene	5	µg/L		
Bromobenzene	5	µg/L		
Chlorobenzene	5	µg/L		

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QCA01	QCB01	QCC01	QCA01_080312	QCB01_080312	QCC01_080312
7/03/2012	7/03/2012	7/03/2012	9/03/2012	9/03/2012	9/03/2012
Tripp Blank	Field Blank	Rinsate Blank	Tripp Blank	Field Blank	Rinsate Blank

Analyte	LOR	Units	ANZECC 2000 -	ANZECC 2000 -
			Marine Water - 95%	Irrigation - LTU
Chlorinated Aliphatic Compounds				
1,1,1,2-Tetrachloroethane	5	µg/L		
1,1,1-Trichloroethane	5	µg/L		
1,1,2,2-Tetrachloroethane	5	µg/L		
1,1,2-Trichloroethane	5	µg/L	1900	
1,1-Dichloroethane	5	µg/L		
1,1-Dichloroethene	5	µg/L		
1,1-Dichloropropylene	5	µg/L		
1,2,3-Trichloropropane	5	µg/L		
1,2-Dibromo-3-chloropropane	5	µg/L		
1,2-Dichloroethane	5	µg/L		
1,3-Dichloropropane	5	µg/L		
Bromomethane	50	µg/L		
Carbon Tetrachloride	5	µg/L		
Chloroethane	50	µg/L		
Chloromethane	50	µg/L		
cis-1,2 Dichloroethene	5	µg/L		
cis-1,4-Dichloro-2-butene	5	µg/L		
Dibromomethane	5	µg/L		
Dichlorodifluoromethane	50	µg/L		
Hexachlorobutadiene	5	µg/L		
Iodomethane	5	µg/L		
Pentachloroethane	5	µg/L		
Tetrachloroethene	5	µg/L		
trans-1,2-Dichloroethene	5	µg/L		
trans-1,4-Dichloro-2-butene	5	µg/L		
Trichloroethane	5	µg/L		
Trichlorofluoromethane	50	µg/L		
Vinyl chloride	50	µg/L		
Fumigants				
1,2-Dibromomethane	5	µg/L		
1,2-Dichloropropane	5	µg/L		
2,2-Dichloropropane	5	µg/L		
cis-1,3-dichloropropene	5	µg/L		
trans-1,3-dichloropropane	5	µg/L		
Trihalomethanes				
Bromodichloromethane	5	µg/L		
Bromofom	5	µg/L		
Chlorofom	5	µg/L		
Dibromochloromethane	5	µg/L		

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7/03/2012	7/03/2012	7/03/2012	9/03/2012	9/03/2012	9/03/2012
Trip Blank	Field Blank	Rinsate Blank	Trip Blank	Field Blank	Rinsate Blank

Analyte	LOR	Units	ANZECC 2000 - Marine Water - 95%	ANZECC 2000 - Irrigation - LTU
<b>Monocyclic Aromatic Hydrocarbons</b>				
Styrene	5	µg/L	<5	-
Isopropylbenzene	5	µg/L	<5	-
n-Butylbenzene	5	µg/L	<5	-
n-Propylbenzene	5	µg/L	<5	-
p-Isopropyltoluene	5	µg/L	<5	-
sec-Butylbenzene	5	µg/L	<5	-
tert-Butylbenzene	5	µg/L	<5	-
1,3,5-Trimethylbenzene	5	µg/L	<5	-
<b>Oxygenated Compounds</b>				
4-Methyl-2-pentanone	50	µg/L	<50	-
Vinyl Acetate	50	µg/L	<50	-
2-Butanone (MEK)	50	µg/L	<50	-
2-Hexanone (MBK)	50	µg/L	<50	-
<b>Sulfonated Compounds</b>				
Carbon disulfide	5	µg/L	<5	-
<b>Total Recoverable Hydrocarbons</b>				
>C10 - C16 Fraction	100	µg/L	<100	<100
>C10 - C40 Fraction (sum)	100	µg/L	<100	<100
>C16 - C34 Fraction	100	µg/L	<100	<100
>C34 - C40 Fraction	100	µg/L	<100	<100
C6 - C10 Fraction	20	µg/L	<20	<20
C6 - C10 Fraction minus BTEX (F1)	20	µg/L	<20	<20

Legend:

Exceeds the ANZECC/ARMCANZ, 2000, Trigger values for marine water ecosystems - Level of protection 95% species

Exceeds the ANZECC/ARMCANZ, 2000, Irri

- Not analysed / not calculated

\* LOR Exceeds Guideline Trigger Value





