AGL Upstream Investments Pty Ltd

2014-2015 Groundwater and Surface Water Monitoring Status Report

Camden Gas Project

8 October 2015





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Glossary

| Acidity | Base neutralising capacity. |
|----------------------------------|--|
| Alkalinity | Acid neutralising capacity. |
| Alluvium | Unconsolidated sediments (clays, sands, gravels and other materials) deposited by flowing water. Deposits can be made by streams on river beds, floodplains, and alluvial fans. |
| Alluvial aquifer | Permeable zones that store and produce groundwater from unconsolidated alluvial sediments. Shallow alluvial aquifers are generally unconfined aquifers. |
| Ammonia | A compound of nitrogen and hydrogen (NH ₃) that is a common by-product of animal waste and landfills but is also found naturally in reduced environments. Ammonia readily converts to nitrate in soils and streams. |
| Anion | An ion with a negative charge – usually non-metal ions when disassociated and dissolved in water. |
| Aquatic ecosystem | The stream channel, lake or estuary bed, water, and (or) biotic communities and the habitat features that occur therein. |
| Aquiclude | An impermeable unit that acts as a barrier to the flow of groundwater from one formation to another. |
| Aquifer | Rock or sediment in a formation, group of formations, or part of a formation that is saturated and sufficiently permeable to transmit economic quantities of water. |
| Aquifer properties | The characteristics of an aquifer that determine its hydraulic behaviour and its response to abstraction. |
| Aquifer, confined | An aquifer that is overlain by low permeability strata. The hydraulic conductivity of the confining bed is significantly lower than that of the aquifer. |
| Aquifer, semi-confined | An aquifer overlain by a low-permeability layer that permits water to slowly flow through it. During pumping, recharge to the aquifer can occur across the leaky confining layer – also known as a leaky artesian or leaky confined aquifer. |
| Aquifer, unconfined | Also known as a water table aquifer. An aquifer in which there are no confining beds between the zone of saturation and the surface. The water table is the upper boundary of an unconfined aquifer. |
| Aquitard | A low permeability unit that can store groundwater and also transmit it slowly from one formation to another. Aquitards retard but do not prevent the movement of water to or from adjacent aquifers. |
| Australian Height Datum (AHD) | The reference point (very close to mean sea level) for all elevation measurements, and used for correlating depths of aquifers and water levels in bores. |

| Beneficial aquifer | An aquifer with a water resource of sufficient quality and quantity to provide either ecosystem protection, raw water for drinking water supply, and agricultural or industrial water. |
|------------------------------|--|
| Bore | A structure drilled below the surface to obtain water from an aquifer or series of aquifers. |
| Boundary | A lateral discontinuity or change in the aquifer resulting in a significant change in hydraulic conductivity, storativity or recharge. |
| Cation | An ion with a positive charge – usually metal ions when disassociated and dissolved in water. |
| Claystone | A non-fissile rock of sedimentary origin composed primarily of clay-sized particles (less than 0.004 mm). |
| Coal | A sedimentary rock derived from the compaction and consolidation of vegetation or swamp deposits to form a fossilised carbonaceous rock. |
| Coal seam | A layer of coal within a sedimentary rock sequence. |
| Coal seam gas (CSG) | Coal seam gas is a form of natural gas (predominantly methane) that is extracted from coal seams. |
| Concentration | The amount or mass of a substance present in a given volume or mass of sample, usually expressed as milligram per litre (water sample) or micrograms per kilogram (sediment sample). |
| Conceptual model | A simplified and idealised representation (usually graphical) of the physical hydrogeologic setting and the hydrogeological understanding of the essential flow processes of the system. This includes the identification and description of the geologic and hydrologic framework, media type, hydraulic properties, sources and sinks, and important aquifer flow and surface-groundwater interaction processes. |
| Confining layer | Low permeability strata that may be saturated but will not allow water to move through it under natural hydraulic gradients. |
| Datalogger | A digital recording instrument that is inserted in monitoring and pumping bores to record pressure measurements and water level variations. |
| Dual permeability aquifer | An aquifer in which groundwater flow is through both the primary porosity of the rock matrix and the secondary porosity of fractures and fissures. |
| Electrical conductivity (EC) | A measure of a fluid's ability to conduct an electrical current and is an estimation of the total ions dissolved. It is often used as a measure of water salinity. |
| Facies | An assemblage or association of mineral, rock, or fossil features reflecting the environment and conditions of origin of the rock. It refers to the appearance and peculiarities that distinguish a rock unit from associated or adjacent units. |
| Fault | A fracture in rock along which there has been an observable amount of displacement. Faults are rarely single planar units; normally they occur as |

| | parallel to sub-parallel sets of planes along which movement has taken place to a greater or lesser extent. Such sets are called fault or fracture zones. |
|----------------------------|--|
| Groundwater | The water contained in interconnected pores or fractures located below the water table in the saturated zone. |
| Groundwater level | The water level measured in a bore; this may be at or close to the water table in unconfined aquifers, or represent the average piezometric level across the screened interval in confined aquifers. |
| Groundwater flow | The movement of water through openings in sediment and rock within the zone of saturation. |
| Groundwater system | A system that is hydrogeologically more similar than different in regard to geological province, hydraulic characteristics and water quality, and may consist of one or more geological formations. |
| Hydraulic conductivity | The rate at which water of a specified density and kinematic viscosity can move through a permeable medium (notionally equivalent to the permeability of an aquifer to fresh water). |
| Hydraulic gradient | The change in total hydraulic head with a change in distance in a given direction. |
| Hydraulic head | A specific measurement of water pressure above a datum. It is usually measured as a water surface elevation, expressed in units of length. In an aquifer, it can be calculated from the depth to water in a monitoring bore. The hydraulic head can be used to determine a hydraulic gradient between two or more points. |
| Hydrochemistry | Chemical characterisation of water (both surface water and groundwater). |
| Hydrogeology | The study of the interrelationships of geologic materials and processes with water, especially groundwater. |
| Hydrology | The study of the occurrence, distribution, and chemistry of all surface waters. |
| lon | An ion is an atom or molecule where the total number of electrons is not equal to the total number of protons, giving it a net positive or negative electrical charge. |
| Limit or reporting (LOR) | The concentration below which a particular analytical method cannot determine, with a high degree of certainty, a concentration. |
| Lithology | The study of rocks and their depositional or formational environment on a large specimen or outcrop scale. |
| Major ions | Constituents commonly present in concentrations exceeding 10 milligram per litre. Dissolved cations generally are calcium, magnesium, sodium, and potassium; the major anions are sulphate, chloride, fluoride, nitrate, and those contributing to alkalinity, most generally assumed to be bicarbonate and carbonate. |
| Methane (CH ₄) | An odourless, colourless, flammable gas, which is the major constituent of natural gas. It is used as a fuel and is an important source of hydrogen and a wide variety of organic compounds. |

| MicroSiemens per centimetre (µS/cm) | A measure of water salinity commonly referred to as EC (see also electrical conductivity). Most commonly measured in the field with calibrated field meters. |
|--|---|
| Monitoring bore | A non-pumping bore, is generally of small diameter that is used to measure the elevation of the water table and/or water quality. Bores generally have a short well screen against a single aquifer through which water can enter. |
| Monitoring period | Refers to data collected between October 2011 and June 2015. |
| Monitoring year | Refers to data collected between July 2014 and June 2015. |
| Normal faulting | Where the fault plane is vertical or dips towards the downthrow side of a fault. |
| Oxidising conditions | Conditions in which a species loses electrons and is present in oxidised form. |
| Permeability | The property or capacity of a porous rock, sediment, clay or soil to transmit a fluid. It is a measure of the relative ease of fluid flow under unequal pressure. The hydraulic conductivity is the permeability of a material for water at the prevailing temperature. |
| Permeable material | Material that permits water to move through it at perceptible rates under the hydraulic gradients normally present. |
| Permian | The last period of the Palaeozoic era that finished approximately 252 million years before present. |
| рН | Potential of Hydrogen; the logarithm of the reciprocal of hydrogen-ion concentration in gram atoms per litre; provides a measure on a scale from 0 to 14 of the acidity or alkalinity of a solution (where 7 is neutral, greater than 7 is alkaline and less than 7 is acidic). |
| Porosity | The proportion of open space within an aquifer, comprised of intergranular space, pores, vesicles and fractures. |
| Porosity, primary | The porosity that represents the original pore openings when a rock or sediment formed. |
| Porosity, secondary | The porosity caused by fractures or weathering in a rock or sediment after it has been formed. |
| Quaternary | The most recent geological period extending from approximately 2.6 million years ago to the present day. |
| Quality assurance | Evaluation of quality-control data to allow quantitative determination of the quality of chemical data collected during a study. Techniques used to collect, process, and analyse water samples are evaluated. |
| Recharge | The process which replenishes groundwater, usually by rainfall infiltrating from the ground surface to the water table and by river water reaching the water table or exposed aquifers. The addition of water to an aquifer. |
| Recharge area | A geographic area that directly receives infiltrated water from surface and in which there are downward components of hydraulic head in the aquifer. Recharge generally moves downward from the water table into the deeper |

| | parts of an aquifer then moves laterally and vertically to recharge other parts of the aquifer or deeper aquifer zones. |
|-----------------------------|---|
| Recovery | The difference between the observed water level during the recovery period after cessation of pumping and the water level measured immediately before pumping stopped. |
| Redox potential (ORP or Eh) | The redox potential is a measure (in volts) of the affinity of a substance for electrons – its electronegativity – compared with hydrogen (which is set at 0). Substances more strongly electronegative than (i.e. capable of oxidising) hydrogen have positive redox potentials. Substances less electronegative than (i.e. capable of reducing) hydrogen have negative redox potentials. Also known as oxidation-reduction potential and Eh. |
| Redox reaction | Redox reactions, or oxidation-reduction reactions, are a family of reactions that are concerned with the transfer of electrons between species, and are mediated by bacterial catalysis. Reduction and oxidation processes exert an important control on the distribution of species like O ₂ , Fe ²⁺ , H ₂ S and CH ₄ etc. in groundwater. |
| Salinity | The concentration of dissolved salts in water, usually expressed in EC units or milligrams of total dissolved solids per litre (mg/L TDS). |
| Salinity classification | Fresh water quality – water with a salinity <800 μ S/cm. |
| | Marginal water quality – water that is more saline than freshwater and generally waters between 800 and 1,600 μ S/cm. |
| | Brackish quality – water that is more saline than freshwater and generally waters between 1,600 and 4,800 $\mu\text{S/cm}.$ |
| | Slightly saline quality – water that is more saline than brackish water and generally waters with a salinity between 4,800 and 10,000 μ S/cm. |
| | Moderately saline quality – water that is more saline than slightly saline water and generally waters between 10,000 and 20,000 μ S/cm. |
| | Saline quality – water that is almost as saline as seawater and generally waters with a salinity greater than 20,000 μ S/cm. |
| | Seawater quality – water that is generally around 55,000 μ S/cm. |
| | (Australian Water Resources Council 1988) |
| Sandstone | Sandstone is a sedimentary rock composed mainly of sand-sized minerals or rock grains (predominantly quartz). |
| Screen | A type of bore lining or casing of special construction, with apertures designed to permit the flow of water into a bore while preventing the entry of aquifer or filter pack material. |
| Sedimentary rock aquifer | These occur in consolidated sediments such as porous sandstones and conglomerates, in which water is stored in the intergranular pores, and limestone, in which water is stored in solution cavities and joints. These aquifers are generally located in sedimentary basins that are continuous over large areas and may be tens or hundreds of metres thick. In terms of quantity, they contain the largest volumes of groundwater. |
| Shale | A laminated sedimentary rock in which the constituent particles are predominantly of clay size. |

| Siltstone | A fine-grained rock of sedimentary origin composed mainly of silt-sized particles (0.004 to 0.06 mm). |
|---|---|
| Standing water level (SWL) | The height to which groundwater rises in a bore after it is drilled and completed, and after a period of pumping when levels return to natural atmospheric or confined pressure levels. |
| Stratigraphy | The depositional order of sedimentary rocks in layers. |
| Surface water- groundwater interaction | This occurs in two ways: (1) streams gain water from groundwater through the streambed when the elevation of the water table adjacent to the streambed is greater than the water level in the stream; and (2) streams lose water to groundwater through streambeds when the elevation of the water table is lower than the water level in the stream. |
| Tertiary | Geologic time at the beginning of the Cainozoic era, 65 to 2.6 million years ago, after the Cretaceous and before the Quaternary. |
| Total Dissolved Solids (TDS) | A measure of the salinity of water, usually expressed in milligrams per litre (mg/L). See also EC. |
| Water quality | Term used to describe the chemical, physical, and biological characteristics of water, usually in respect to its suitability for a particular purpose. |
| Water quality data | Chemical, biological, and physical measurements or observations of the characteristics of surface and ground waters, atmospheric deposition, potable water, treated effluents, and waste water and of the immediate environment in which the water exists. |
| Well | Pertaining to a gas exploration well or gas production well. |

Abbreviations

| AGL | AGL Upstream Investments Pty Ltd | | |
|------|---|--|--|
| ВоМ | Bureau of Meteorology | | |
| BTEX | Benzene, toluene, ethyl benzene and xylenes | | |
| CDFM | Cumulative deviation from mean | | |
| CGP | Camden Gas Project | | |
| CSG | Coal seam gas | | |
| DO | Dissolved oxygen | | |
| EC | Electrical conductivity | | |
| LOR | Limit of reporting | | |
| ORP | Oxidation reduction potential | | |
| РАН | Polycyclic aromatic hydrocarbons | | |
| SCA | Sydney Catchment Authority | | |
| TDS | Total dissolved solids | | |
| ТРН | Total petroleum hydrocarbons | | |
| VWP | Vibrating wire piezometer | | |

Units

| °C | degrees Celsius |
|-------|--------------------------------|
| L/s | litres per second |
| m | metres |
| mAHD | metres Australian Height Datum |
| mbgl | metres below ground level |
| m/d | metres per day |
| mg/L | milligrams per litre |
| mV | millivolt |
| µg/L | micro grams per litre |
| µS/cm | microSiemens per centimetre |

Executive Summary

AGL Upstream Investments Pty Ltd (AGL) owns and operates the Camden Gas Project (CGP) located in the Macarthur region, 65 km southwest of Sydney, NSW. The CGP has been producing natural gas from coal seams for the Sydney region since 2001 and currently consists of 144 gas wells (of which, approximately 95 were operational on 30 June 2015). The target coal seams are the Bulli and Balgownie Coal Seams within the Illawarra Coal Measures at depths of approximately 550 – 700 metres below ground level (mbgl).

The CGP monitoring network comprises three nested groundwater monitoring sites (11 monitoring bores): Denham Court (monitored since 2011), Menangle Park (monitored since June 2013) and Glenlee (monitored since February 2014). One monitoring bore is installed in the alluvium near the Nepean River, two in the Ashfield Shale and eight monitoring bores at different depths in the Hawkesbury Sandstone. Groundwater levels have been recorded at six-hourly intervals and water quality data has been collected on a quarterly basis since installation. Surface water is monitored annually at one monitoring location along the Nepean River next to the Menangle Park site. This report presents an assessment of water level and water quality data from the groundwater monitoring network and from the Nepean River for the period up to 30 June 2015, with an emphasis on data obtained during the past 12 months.

The groundwater level in the Nepean River alluvium is shallow and shows a direct response to rainfall and flood events during the monitoring period. Groundwater levels in the Ashfield Shale are deep (~80 mbgl) and show no apparent response to rainfall over the monitoring period. Groundwater levels appear to follow similar trends in each of the screened Hawkesbury Sandstone water bearing zones (defined as upper, middle and lower). There is no apparent response to individual rainfall events over the monitoring period at the Denham Court and Glenlee sites, while a clear response to rainfall and flood events can be observed at the three uppermost monitoring bores at the Menangle Park site. Groundwater levels during the 2014/15 monitoring year were comparable to groundwater levels as recorded during previous monitoring years.

Groundwater sampled from the alluvium at the Menangle Park site is characterised as fresh to marginally saline. Dissolved metal concentrations were generally low and no dissolved gases were detected. Hydrocarbons (polycyclic aromatic hydrocarbons, total petroleum hydrocarbons and toluene) were detected at low levels. Groundwater sampled from the Hawkesbury Sandstone is fresh to marginal at the Menangle Park site and slightly saline at the Denham Court and Glenlee sites. Salinity decreases with depth at the Denham Court and Glenlee sites. Dissolved metal concentrations are generally low and minor detections of hydrocarbons were present at the three monitoring sites. Dissolved methane was detected at all Hawkesbury Sandstone bores, although concentrations at the Menangle Park and Glenlee sites were comparable to the control site at Denham Court. Low concentrations of dissolved ethane was detected at the Denham Court and Glenlee sites and low concentrations of propane and butane were detected only at the Glenlee site.

Groundwater quality during the 2014/15 monitoring year was overall comparable to groundwater quality as measured during the previous monitoring years.

The Nepean River water is characterised as fresh, neutral pH and dissolved metal concentrations were comparably lower than occurring in groundwater in the alluvium and underlying Hawkesbury Sandstone. Ammonia concentrations in Nepean River water exceeded the ANZECC (2000) guideline value and minor dissolved methane concentrations and no hydrocarbons were detected.

From assessment of the available data, there are no observable impacts to groundwater levels or quality that could be attributable to the CSG operations. There is also no evidence of connectivity between the shallower monitored zones and the coal seams and this is in agreement with the conceptual model (Parsons Brinckerhoff 2011), that indicates the presence of extensive and thick claystone formations (aquitards and aquicludes) between the Hawkesbury Sandstone and coal seams restricting upward depressurisation and impeding the vertical flow of groundwater.

1. Introduction

1.1 Background

AGL Upstream Investments Pty Ltd (AGL) owns and operates the Camden Gas Project (CGP) located in the Macarthur region, 65 km southwest of Sydney, NSW. The CGP has been producing natural gas from coal seams for the Sydney region since 2001 and currently consists of 144 gas wells (of which, approximately 95 were operational on 30 June 2015) within the Stage 1 and Stage 2 areas (Figure 1.1). A proposal for the expansion of the project into Stage 3 (Northern Expansion) was suspended in 2013 and was officially withdrawn in July 2015. The target coal seams are the Bulli and Balgownie Coal Seams within the Illawarra Coal Measures at depths of approximately 550 –700 metres below ground level (mbgl).

Parsons Brinckerhoff was engaged by AGL to characterise the hydrogeological environment and conduct routine groundwater monitoring and interpretation in relation to the CGP activities. Installation of a dedicated water monitoring network of 11 monitoring bores occurred between October 2011 and February 2014. The groundwater monitoring network comprises dedicated monitoring bores in the alluvium, the Ashfield Shale, and the Hawkesbury Sandstone as well as one surface water monitoring location that is monitored annually. The collection of groundwater level and groundwater quality data commenced in October 2011. Groundwater levels have been recorded at six-hourly intervals and following one initial baseline sample in November 2011, water quality data was collected on a quarterly basis between May 2013 and April 2015. From April 2015 onwards, water quality data is collected at six-monthly intervals. This reporting provides an emphasis on evaluation of data obtained during the past 12 months (July 2014 to June 2015 monitoring year) with comparison to the data obtained during the past monitoring years (Parsons Brinckerhoff 2012, 2013a, 2014a and 2014b).

Monitoring is undertaken at three sites located within the Camden-Campbelltown area, NSW (Figure 1.1). The Denham Court site is located to the north, outside of the CGP area, and can be considered as a control site given the distance from the operating CGP. The Menangle Park and Glenlee sites are located within the existing CGP wellfield (Stage 1 and 2, Figure 1.1).

The objective of the groundwater monitoring program is to provide water levels and water quality attributes for each of the monitored groundwater systems of the region, in areas within and distant from the currently operating CGP.

1.2 Scope of works

This report presents and interprets groundwater level data collected at six-hourly intervals and groundwater quality data collected since monitoring began at each of the established sites up to 30 June 2015, with emphasis on the data obtained during the past 12 months.

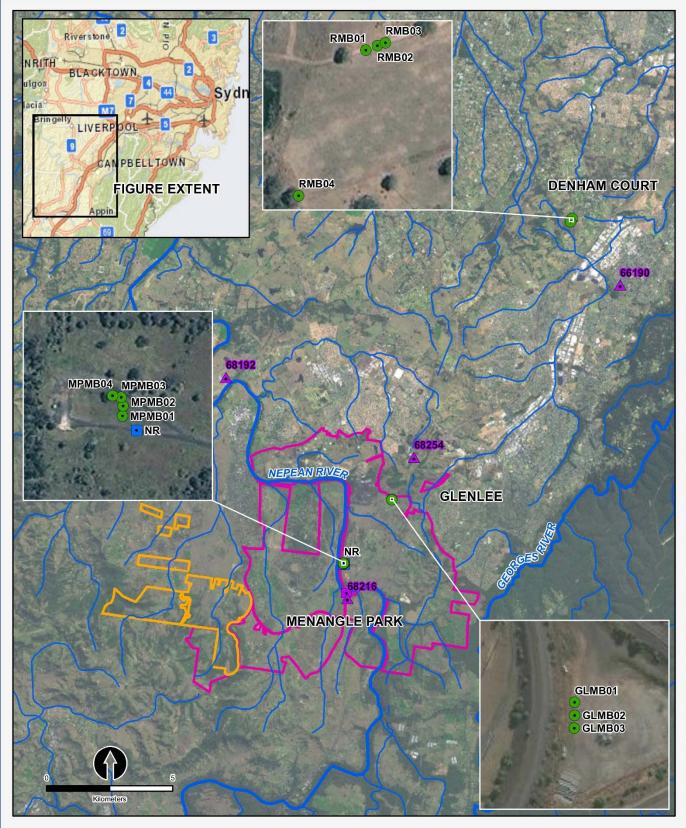
The scope of works was to:

- Conduct groundwater monitoring, including six-hourly groundwater level measurements and three groundwater quality sampling events testing for field parameters, major cations and anions, dissolved metals, nutrients, dissolved methane and hydrocarbons.
- Conduct surface water quality sampling on one occasion (September 2014) at one location (the Nepean River near the Menangle Park site as shown on Figure 1.1).
- Analyse and interpret water level and hydrochemical results with reference to the conceptual model where relevant.

• Establish whether there are any observable impacts from CSG activities within the shallow aquifers.

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CAMDEN GAS PROJECT AGL UPSTREAM INVESTMENTS PTY LTD



- C Stage 1 C Stage 2
- Camden Gas Project areas Groundwater monitoring bore ٠ Surface water sampling location
 - . BoM surface water gauging station
 - A BoM weather station
 - Rivers
 - Streams

Figure 1.1: Groundwater monitoring locations Camden Gas Project

2. Site characterisation

2.1 Rainfall

The nearest Bureau of Meteorology (BoM) weather station with consistent historical climate measurements is located at Camden airport (BoM site number 68192), approximately 2.5 km northwest of the Stage 2 area (Figure 1.1). Mean temperatures at Camden airport range from 17.3°C in July to 29.5°C in January (based on records from 1971 to 2015). The average annual rainfall is 788 mm (based on records from 1972 to 2015). On average, July receives the least rain, with a mean rainfall of 36.8 mm, while February receives the most rain, with a mean of 98.8 mm (BoM 2015).

The long-term, annual cumulative deviation from mean (CDFM) rainfall for Camden airport is plotted in Figure 2.1. Annual rainfall data for BoM site number 68216 (located approximately 10 km southeast from the Camden airport 68192 station) is presented from 2009 onwards as rainfall records at Camden airport were incomplete for that period. The long-term CDFM is generated by subtracting the long-term average annual rainfall for the recorded period from the actual annual rainfall and then accumulating these residuals over the assessment period. Periods of below average rainfall are represented as downward trending slopes while periods of above average rainfall are represented as upward trending slopes.

The cumulative deviation plot for Camden airport (Figure 2.1) shows a relatively wet period between 1972 and 1992, followed by a relatively dry period between 1998 and 2007. From 2007 onwards the rainfall has been typically close to the mean rainfall of 776 mm per year.

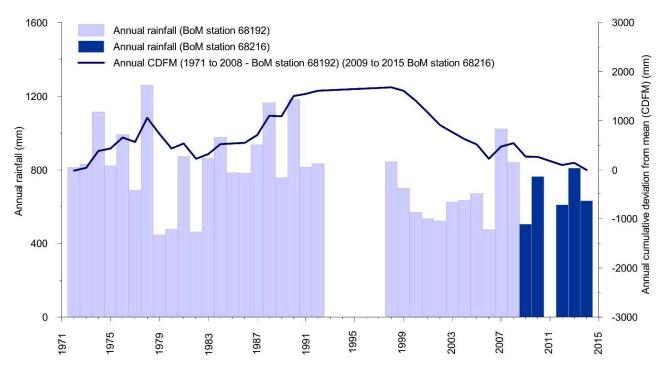


Figure 2.1 Cumulative deviation from annual mean rainfall at Camden airport (BoM station 068192)

The monitoring bore sites are separated by up to a distance of 16 km and therefore groundwater level data for each site have been compared with rainfall data from the closest BoM station (Figure 1.1) as follows:

- Denham Court: 66190 Ingleburn Station.
- Menangle Park: 68216 Menangle Bridge.

Glenlee: 68254 Mount Annan Botanic Garden.

The rainfall characteristics are broadly similar between these BoM stations during the monitoring period, as presented in Figure 2.2. Total monthly rainfall for this monitoring period was overall higher than the long-term average.

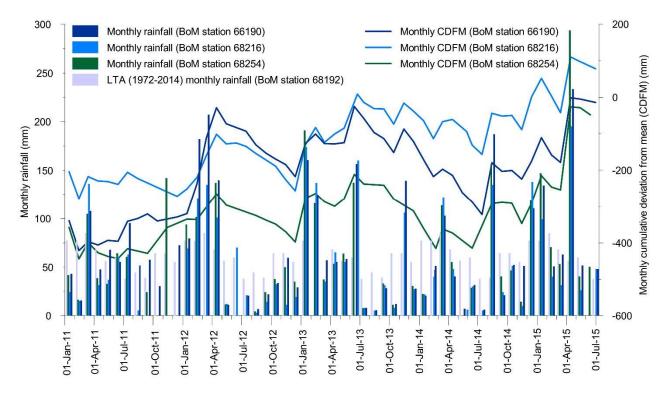


Figure 2.2 Monthly rainfall for the monitoring period (2011-2015)

2.2 Surface hydrology

The CGP includes two catchment areas: the Hawkesbury Nepean Catchment and the Sydney Metropolitan Catchment. The major surface hydrology features in the CGP are the Nepean River and its tributaries, which meander in a south to north direction within the project area; and the Georges River, which flows in a northerly direction, in the south-east of the project area.

Small farm dams are common in rural areas and provide water for stock, limited garden and irrigation purposes. Dams are replenished by rainfall and runoff, although some seepage flow through weathered soil profiles occurs after long wet periods. Dams and seepage flows are not related to the regional groundwater systems. There are no known springs in the CGP area.

2.3 Geological setting

The CGP is part of the Southern Coalfield of the Sydney Geological Basin. The Basin is primarily a Permo-Triassic sedimentary rock sequence (Parkin 2002) and is underlain by undifferentiated sediments of Carboniferous and Devonian age. The stratigraphy of the CGP in the Camden-Campbelltown area is summarised in Table 2.1. The geology and structure of the CGP is shown on Figure 2.3.

| Period | Group | Sub- group | Formation | Description | Average thickness (m) ^a |
|------------|-------------------------------------|--|---|--|--|
| Quaternary | | | Alluvium | Quartz and lithic 'fluvial' sand, silt and clay. | |
| Tertiary | a ta | | Alluvium | High level alluvium. | <20 |
| | | | Bringelly Shale | Shale, carbonaceous claystone, laminate, lithic sandstone, rare coal. | 80 (top eroded) |
| | Wianamatta Group | | Minchinbury Shale | Fine to medium-grained lithic sandstone. | - |
| | Wia | | Ashfield Shale ^b | Black to light grey shale and laminate (Bembrick <i>et al.</i> 1987). | - |
| | | | Mittagong Formation | Dark grey to grey alternating beds of shale laminate, siltstone and quartzose sandstone (Alder <i>et al.</i> 1991). | 11 |
| | | | Hawkesbury Sandstone | Massive or thickly bedded quartzose sandstone with siltstone, claystone and grey shale lenses up to several metres thick (Bowman 1974; Moffitt 2000). | 173 |
| <u>.</u> | | Subgroup | Newport Formation | Fine-grained sandstone (less than 3 m thick) interbedded with light to dark grey, fine-grained sandstones, siltstones and minor claystones (Bowman 1974). | 35 |
| Triassic | Narrabeen Group Clifton Subgroup | Garie Formation | Cream, massive, kaolinite-rich pelletal claystone, which grades upwards to grey, slightly carbonaceous claystone containing plant fossils at the base of the Newport Formation (Moffitt 2000). | 8 | |
| | | | Bald Hill Claystone ^b | Massive chocolate coloured and cream pelletal claystones and mudstones, and occasional fine-grained channel sand units (Moffitt 2000). | 34 |
| | | roup | Bulgo Sandstone | Thickly bedded sandstone with intercalated siltstone and claystone bands up to 3 m thick (Moffitt 2000). | 251 |
| | | on Subç | Stanwell Park Claystone ^b | Red-green-grey shale and quartz sandstone (Moffitt 1999). | 36 |
| | | Clift | Scarborough Sandstone | Quartz-lithic sandstone, pebbly in part (Moffitt 1999). | 20 |
| | | | Wombarra Claystone ^b | Grey shale and minor quartz-lithic sandstone (Moffitt 1999). | 32 |
| | | | Bulli Coal Seam | | 4 |
| | res | Illawarra Coal Measures Sydney Subgroup | Loddon Sandstone | | 12 |
| | Measu | | Balmain Coal Member | Coal interbedded with shale, quartz-lithic sandstone, | 24 |
| Permian | Illawarra Coal | | Balgownie Coal Seam | conglomerate, chert, torbanite seams and occasionally carbonaceous mudstone (Moffitt 2000). | 2 |
| Å | | | (Remaining Sydney Subgroup) | | - |
| | | Cumberla | nd Subgroup | | - |
| | Shoalhaven Group | | | Sandstone, siltstone, shale, polymictic conglomerate, claystone; rare tuff, carbonate, evaporate. | - |
| Palaeozic | Lachlan F | old Belt | | Intensely folded and faulted slates, phyllites, quartzite sandstones and minor limestones of Ordovician to Silurian age (Moffitt 2000). | - |

Notes:

(a) Average thickness from available well data within CGP (AGL 2013); (b) Aquitard or aquiclude

The Illawarra Coal Measures is the economic sequence of interest for CSG development in the area, and consists of interbedded sandstone, shale and coal seams, with a thickness of approximately 300 m. The upper sections of the Permian Illawarra Coal Measures (Sydney Subgroup) contain the major coal seams: Bulli Coal Seam, Balgownie Coal Seam, Wongawilli Coal Seam, and Tongarra Coal Seam. The seams targeted for CSG production within the CGP are the Bulli and Balgownie coal seams, both of which are 2 to 5 m thick within the CGP.

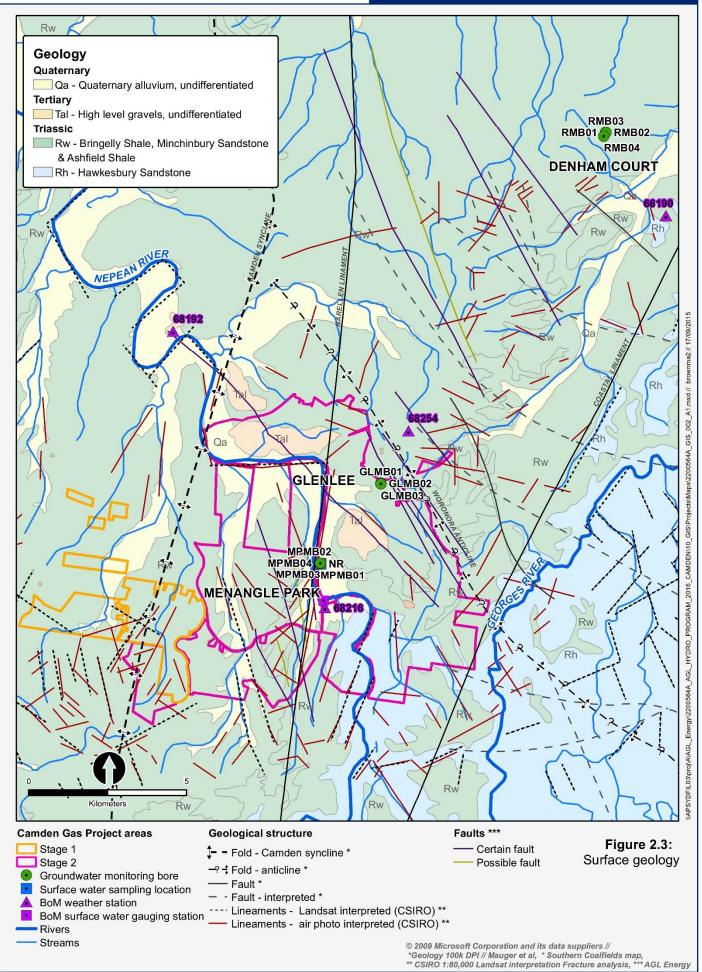
The Illawarra Coal Measures is overlain by Triassic sandstones, siltstones and claystones of the Narrabeen Group and the Hawkesbury Sandstone. Overlying the Hawkesbury Sandstone is the Triassic Wianamatta Group shales which comprises the majority of the surficial geology (where thin alluvial deposits are not present).

Structurally, the CGP area and surrounds are dominated by the north-northeast plunging Camden Syncline, which is a broad and gentle warp structure (Alder *et al.* 1991 and Bray *et al.* 2010). The Camden Syncline is bounded in the west and truncated in the southwest by the north-south trending Nepean Structural Zone, part of the Lapstone Structural Complex.

The CGP is relatively unaffected by major faulting apart from a set of NW-NNW trending faults associated with the Lapstone Monocline Structure (Alder *et al.* 1991 and Blevin *et al.* 2007). These faults have been identified from exploration and 2D seismic studies and they have been identified as high-angle, low to moderate displacement normal faults (Blevin *et al.* 2007). Many of these features intersect coal seams however very few, if any, affect the entire stratigraphic sequence and display no expression at surface.

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2.4 Hydrogeological setting

The Southern Coalfield is located within the Sydney Basin sedimentary rock groundwater system. The recognised hydrogeological units within the CGP are shown in Table 2.2.

| Table 2.2 | Hydrogeological units within the CGP area |
|-----------|--|
| | Tryanogeological antis within the ool area |

| Hydrogeological unit | Aquifer type | Hydraulic conductivity – horizontal (m/d) | Hydraulic conductivity – vertical (m/d) | TDS (mg/L) |
|--|--------------------------------------|---|---|---------------|
| Alluvium | Unconfined aquifer | 0.15 - 10 | 1 - 10 | <3,000 |
| Ashfield Shale (Wianamatta Group) | Aquitard or unconfined/perched | Very low | 0.05 | >3,000 |
| Hawkesbury Sandstone | Unconfined/semi- confined aquifer | 0.0009 - 0.69 | 0.05 - 6 x 10 ⁻⁴ | <500 - 10,000 |
| Bald Hill Claystone (Narrabeen Group) | Aquitard/aquiclude | 1 x 10 ⁻⁵ | 2 x 10 ⁻⁶ | - |
| Bulgo Sandstone (Narrabeen Group) | Confined aquifer | 5 x 10 ⁻⁴ | 1 x 10 ⁻⁴ | 1,500 - 5,000 |
| Stanwell Park Claystone (Narrabeen Group) | Aquitard/aquiclude | 3 x 10⁻⁵ | 6 x 10 ⁻⁶ | - |
| Scarborough Sandstone (Narrabeen Group) | Confined aquifer | 0.01 | 5 x 10 ⁻³ | - |
| Wombarra Claystone (Narrabeen Group) | Aquitard/aquiclude | 3 x 10 ⁻⁵ | 6 x 10 ⁻⁶ | - |
| Illawarra Coal Measures | Confined water bearing zones | 5 x 10 ⁻² (Bulli Coal Seam) | 2.5 x 10 ⁻² (Bulli Coal Seam) | >2,000 |

Note: table summarises data from a number of investigations including SCA (2005); GHD (2007); Broadstock (2011); Parsons Brinckerhoff (2011); AGL (2013); Parsons Brinckerhoff (2014a).

Alluvium occurs along the floodplain of the Nepean River and its tributaries. Alluvial deposits are generally thin, discontinuous (except along the Nepean River) and relatively permeable. The unconfined aquifers within the alluvium are responsive to rainfall and stream flow and form a minor beneficial aquifer. There are also small terrace areas of Tertiary alluvium within the CGP area that contain localised aquifers of variable quality (Figure 2.3).

The Ashfield Shale (which outcrops across the majority of the CGP) is generally of low permeability and yield; however small water bearing zones are sometimes present. Water is typically brackish to saline, especially in low relief areas of western Sydney (due to the marine depositional environment of the shales) (Old 1942). Average bore yields are 1.3 litres per second (L/s) (AGL 2013).

The Hawkesbury Sandstone and Narrabeen Group form part of an extensive generally semi-confined regional aquifer system within the Sydney Basin sequence. The Hawkesbury Sandstone is more widely exploited for groundwater than the overlying and underlying formations, being of generally higher yield, better water quality and either outcropping or buried to shallow depths over the basin. Groundwater flow within the Hawkesbury Sandstone and Narrabeen Group aquifers at a regional scale has a major horizontal

component, due to the alternation of sheet and massive facies, with some vertical leakage. The Hawkesbury Sandstone and Narrabeen Group are characterised by dual porosity. The primary porosity is imparted by connected void space between sand grains and the secondary porosity is due to the interconnected rock defects such as joints, fractures, faults and bedding planes. Superior bore yield in the sandstone aquifers of the Hawkesbury Sandstone is often associated with major fractures or a high fracture zone density, and yields of up to 40 L/s have been recorded in bores intercepting these zones within deformed areas of the Sydney Basin (McLean and Ross 2009). Typically within the CGP area bore yields within the Hawkesbury Sandstone rarely exceed 2 L/s (SCA 2007 and Ross 2014). The Narrabeen Group aquifer is generally not used as a water source as it is considered to be poorer quality and lower permeability compared to the overlying Hawkesbury Sandstone aquifer (Madden 2009).

There is a lack of major fracturing and fault systems intersecting the Hawkesbury Sandstone within the CGP. Yields in the Hawkesbury Sandstone are highest and salinities are freshest south of the Nepean River due to the proximity to recharge areas. North of the Nepean River, the groundwater within the Hawkesbury Sandstone is characterised by higher salinity, becoming moderately saline. Groundwater is used for irrigation and domestic purposes to the south and immediately to the north of the Nepean River; however, further north of the river, groundwater quality is typically only suitable for stock (AGL 2013).

The coal seams present in the Illawarra Coal Measures are both regionally and locally minor water bearing zones. Due to the greater depth of burial of the coal measures and fine-grained nature of the sedimentary rocks, the permeability is generally lower than the overlying sandstone aquifers. Recharge to the Permian water bearing zones is likely to occur where formations are outcropping, which occurs a significant distance to the south of the CGP. Salinity of the water bearing zones is typically brackish to moderately saline.

Within the CGP, there is limited rainfall recharge to the Ashfield Shale with most rainfall generating runoff and overland flow. Some leakage through the Ashfield Shale into the Hawkesbury Sandstone is expected where there is adequate fracture spacing, however, it is anticipated that most recharge to the sandstone aquifers occurs via lateral groundwater through-flow from upgradient and updip areas to the south. There is insufficient data within the CGP to define local flow paths and natural discharge zones; however, regionally, groundwater flow is predominantly towards the north or northeast, eventually discharging via the Georges, Parramatta or Hawkesbury River systems, and ultimately offshore to the east. Locally, there may be a small base flow or interflow discharge component to local stream headwaters during wet periods; however groundwater-surface water interactions are not well defined within the area (Parsons Brinckerhoff 2010).

3. Monitoring program

3.1 Monitoring network

Table 3.1 and Figure 3.1 provide construction details for the 11 monitoring bores within the CGP area (Figure 1.1).

| Monitoring bore | Location | Total depth (mbgl) | Screened interval (mbgl) | Lithology | Formation |
|--------------------|---------------|-----------------------|--------------------------------|----------------|----------------------------------|
| RMB01 | Denham Court | 84.0 | 69.0 - 81.0 | Siltstone | Ashfield Shale |
| RMB02 | Denham Court | 150.0 | 135.0 – 147.0 | Sandstone | Hawkesbury Sandstone (upper) |
| RMB03 | Denham Court | 300.0 | 290.0 - 299.0 | Sandstone | Hawkesbury Sandstone (lower) |
| RMB04 | Denham Court | 8.5 | 4.5 – 7.5 | Clay/siltstone | Ashfield Shale (weathered) |
| MPMB01 | Menangle Park | 18.5 | 10.0 – 16.0 | Clay | Alluvium |
| MPMB02 | Menangle Park | 42.0 | 27.4 - 39.4 | Sandstone | Hawkesbury Sandstone (upper) |
| MPMB03 | Menangle Park | 108.5 | 97.0 - 106.0 | Sandstone | Hawkesbury Sandstone (middle) |
| MPMB04 | Menangle Park | 192.6 | 182.6 – 191.6 | Sandstone | Hawkesbury Sandstone (lower) |
| GLMB01 | Glenlee | 102.2 | 87.0 – 99.0ª | Sandstone | Hawkesbury Sandstone (upper) |
| GLMB02 | Glenlee | 190.3 | 168.0 – 180.0ª | Sandstone | Hawkesbury Sandstone (middle) |
| GLMB03 | Glenlee | 228.3 | 212.0 – 224.0 | Sandstone | Hawkesbury Sandstone (lower) |

Table 3.1 Groundwater monitoring bore details

Note:

(a) Monitoring bores GLMB01 and GLMB02 were converted to vibrating wire piezometers (VWP) on 12 March 2015 to maintain borehole integrity (Parsons Brinckerhoff 2015b); the VWP sensors are installed at 93.0 mbgl and 174 mbgl respectively.

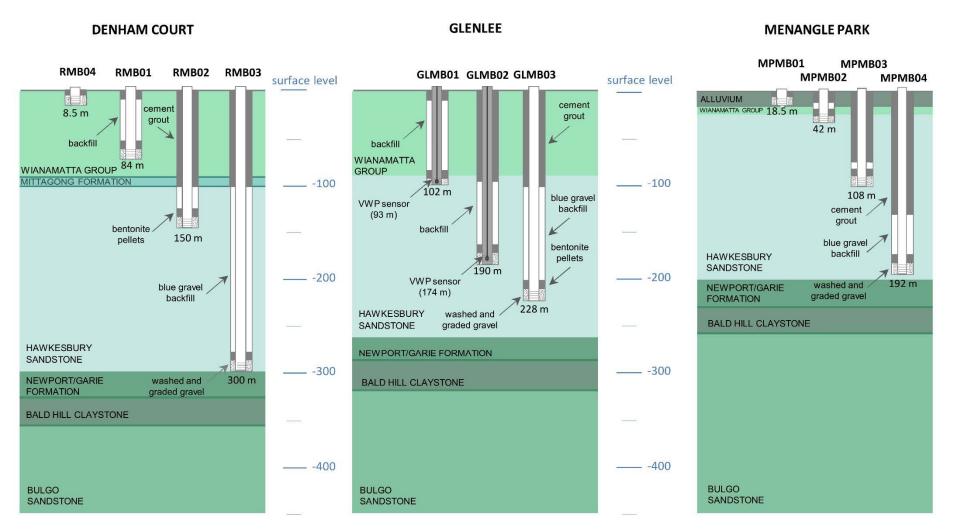


Figure 3.1 Nested groundwater monitoring bores at the Denham Court, Glenlee (following conversion to VWPs in March 2015) and Menangle Park sites

3.2 Water level monitoring

3.2.1 Groundwater levels

Following completion of each monitoring bore, pressure transducers (Solinist Levelogger (M30) dataloggers) were suspended from a galvanised steel wire in the water column and programmed to record a groundwater level every six hours. To verify the level recorded by the dataloggers, manual measurements are recorded periodically using an electronic dip meter. The monitoring start date of the datalogger data at each monitoring bore is shown in Table 3.2.

Barometric loggers installed above the water table at monitoring bores RMB01 and MPMB01 record changes in atmospheric pressure. Data from these loggers is used to correct for the effects of changing barometric pressure on water level loggers in the adjacent monitoring bores.

Table 3.2 Summary of current water level monitoring locations and data collection periods

| Monitoring location | Monitoring start date | |
|---|---|--|
| Denham Court (RMB01, RMB02, RMB03, RMB04) | November 2011 (and June 2013 for RMB04) | |
| Menangle Park (MPMB01, MPMB02, MPMB03, MPMB04)) | June 2013 | |
| Glenlee (GLMB01, GLMB02, GLMB03) | February 2014 | |

The VWP sensors at GLMB01 and GLMB02, which were installed in March 2015, have not stabilised since installation. As such, the data has not been presented in this report; however, groundwater level data that was collected from these bores prior to their conversion to VWPs is included in this report.

3.2.2 Surface water levels

Water levels in the Nepean River are monitored by the BoM (gauging station 68216) using automatic dataloggers close to the Menangle Park site (Figure 1.1). These water levels have been included in the hydrograph for the Menangle Park site for comparison (refer to Section 4 Groundwater levels; Figure 4.2). The river height data is real-time operational data from automated telemetry systems and has been processed to remove erroneous data.

3.3 Water quality monitoring

3.3.1 Overview

Groundwater sampling has been undertaken on nine occasions at Denham Court (since November 2011), seven occasions at Menangle Park (since August 2013) and five occasions at Glenlee (since February 2014) with details provide in Table 3.3. Three sampling events occurred in the 2014/15 monitoring year.

| Sampling event | Denham Court | Menangle Park | Glenlee | Reference report |
|----------------|----------------|---------------|--------------|------------------------------|
| November 2011 | \sqrt{a} | - | - | Parsons Brinckerhoff (2012) |
| May 2013 | √a | - | - | Parsons Brinckerhoff (2013a) |
| August 2013 | √b | νc | - | Parsons Brinckerhoff (2013c) |
| November 2013 | √d | \checkmark | - | Parsons Brinckerhoff (2014c) |
| February 2014 | √ b | \checkmark | \checkmark | Parsons Brinckerhoff (2014d) |
| May 2014 | √b | \checkmark | \checkmark | Parsons Brinckerhoff (2014e) |
| August 2014 | √ ^b | \checkmark | \checkmark | Parsons Brinckerhoff (2014f) |
| January 2015 | √ ^b | \checkmark | \checkmark | Parsons Brinckerhoff (2015a) |
| April 2015 | √b | \checkmark | √e | Parsons Brinckerhoff (2015b) |

Table 3.3 Groundwater quality sampling program

Notes:

(a) RMB01 not sampled due to insufficient water in monitoring bore.

(b) RMB01 and RMB04 not sampled due to insufficient water in monitoring bores.

(c) MPMB04 not sampled due to blockage in monitoring bore (Parsons Brinckerhoff 2013b).

(d) RMB04 not sampled due to insufficient water in monitoring bore.

(e) GLMB01 and GLMB02 not sampled as converted to vibrating wire piezometer in March 2015 (Parsons Brinckerhoff 2015b).

- = monitoring locations not yet installed.

Surface water quality sampling was undertaken on one occasion (9 September 2014) at the Nepean River site next to the Menangle Park groundwater monitoring site.

3.3.2 Sampling techniques

Two methods were used to obtain groundwater quality samples from the monitoring bores. The methods were selected based on the permeability of the screened formation of each bore, which was determined during hydraulic conductivity testing. Higher yielding bores were purged and sampled using a submersible pump. Lower yielding bores and selected deeper bores with high purge volumes were sampled using a low flow pump. In summary:

- A submersible 12V pump was used in monitoring bores MPMB01 and MPMB02 (higher yielding bores).
- A micro-purge[™] low flow sampling pump was used in monitoring bores RMB02, RMB03, MPMB03, MPMB04, GLMB01, GLMB02 and GLMB03 (lower yielding bores).
- RMB01 was not sampled during this monitoring year due to insufficient water in the monitoring bore at each sampling event.
- RMB04 has never been sampled because no perched water has ever been detected within the bore.

Where a submersible pump was used, a minimum of three well volumes was purged from the monitoring bore prior to sampling to allow a representative groundwater sample to be collected. Water quality parameters were measured during and immediately after purging to monitor water quality changes and to indicate representative groundwater suitable for sampling and analysis.

The micro-purge[™] system allows groundwater to be drawn into the pump intake directly from the screened portion of the aquifer, eliminating the need to purge relatively large volumes of groundwater from these bores. Water quality parameters were monitored during the micro-purge[™] pumping to ensure that a representative groundwater sample was collected.

Physiochemical parameters (pH, electrical conductivity (EC), temperature, total dissolved solids (TDS), dissolved oxygen (DO) and oxidation reduction potential (ORP)) were measured during and following purging using a calibrated hand-held water quality meter.

The surface water sample was taken at the river bank using a telescopic sampler. The sample was collected from just below the water surface and approximately 1 m away from the river bank.

3.3.3 Chemical analysis of water

Groundwater samples collected in the field were analysed for a broad chemical suite designed specifically to assess the chemical characteristics of the different water bearing zones at the monitoring sites. Table 3.4 details the groundwater analytical suite.

| Table 3.4 | Groundwater a | nalytical suite |
|-----------|---------------|-----------------|
|-----------|---------------|-----------------|

| Category | Parameters | |
|---|---|---|
| Physiochemical parameters (measured in the field) | EC Temperature DO | pH ORP TDS |
| General parameters | EC ^a TDS | pH ^{a,c} |
| Major ions | Calcium Magnesium Sodium Potassium | Chloride Bicarbonate Sulphate Fluoride Silica |
| Metals and minor/trace elements | Aluminium Antimony ^a Arsenic Barium Boron Bromine Beryllium Cadmium Cobalt Copper Cyanide ^a | Iron Lead Manganese Molybdenum Mercury ^b Nickel Selenium Strontium Zinc Uranium Vanadium |
| Nutrients | Ammonia Nitrite Nitrate | Total organic carbon (TOC) Phosphorus (total) Phosphorus (reactive) |
| Hydrocarbons | Phenol compounds Polycyclic aromatic hydrocarbons (PAH) | Total petroleum hydrocarbons (TPH) Benzene, toluene, ethyl benzene and xylenes (BTEX) |
| Dissolved gases | Methane Ethene Ethane Propene | Propane Butene Butane |

Notes:

(a) For samples collected since the May 2013 sampling event.

(b) For samples collected since the August 2013 sampling event.

(c) Generally analysed outside of recommended holding times.

Groundwater samples for laboratory analysis were collected in sample bottles specified by the laboratory, with appropriate preservation where required. Samples undergoing dissolved metal analysis were filtered through 0.45 µm filters in the field prior to collection.

3.3.4 Quality assurance

3.3.4.1 Field QA/QC

The field sampling procedures conformed to Parsons Brinckerhoff's quality assurance/quality control protocols to prevent cross-contamination and preserve sample integrity. The following QA/QC procedures were applied:

- samples were collected in clearly labelled bottles with appropriate preservation solutions
- samples were delivered to the laboratories within the specified holding times (except for pH).
- unstable parameters were analysed in the field (physiochemical parameters).

3.3.4.2 Laboratory QA/QC

The laboratories conduct their own internal QA/QC program to assess the repeatability of the analytical procedures and instrument accuracy. These programs include analysis of laboratory sample duplicates, spike samples, certified reference standards, surrogate standards/spikes and laboratory blanks. In addition, a duplicate sample is collected in the field for every ten samples collected to assess sampling and laboratory analysis accuracy.

4. Groundwater levels

Hydrographs showing groundwater levels and rainfall from the start of monitoring until the end of June 2015 are presented for Denham Court in Figure 4.1, Menangle Park in Figure 4.2 and Glenlee in Figure 4.3. The Menangle Park site is located close to the Nepean River and river levels from BoM gauging station 68216 have also been included in the hydrograph for comparison (Figure 4.2). Individual hydrographs for each monitoring bore are included in Appendix A.

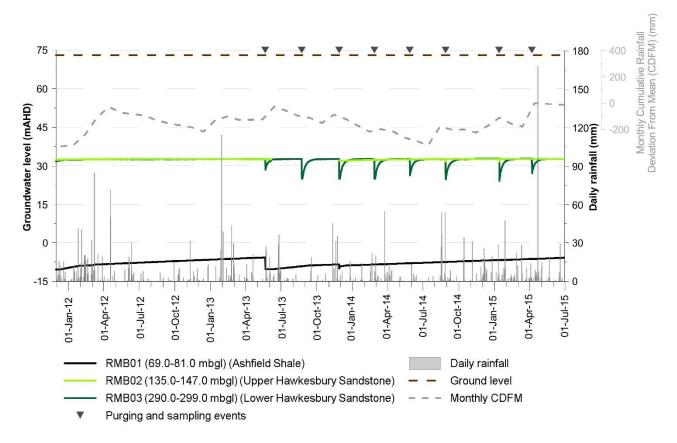
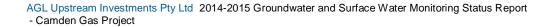


Figure 4.1 Groundwater levels at the Denham Court site



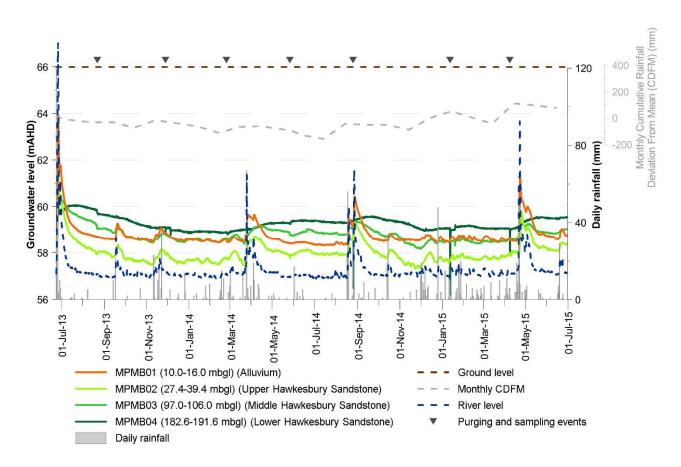


Figure 4.2 Groundwater levels and Nepean River level at the Menangle Park site

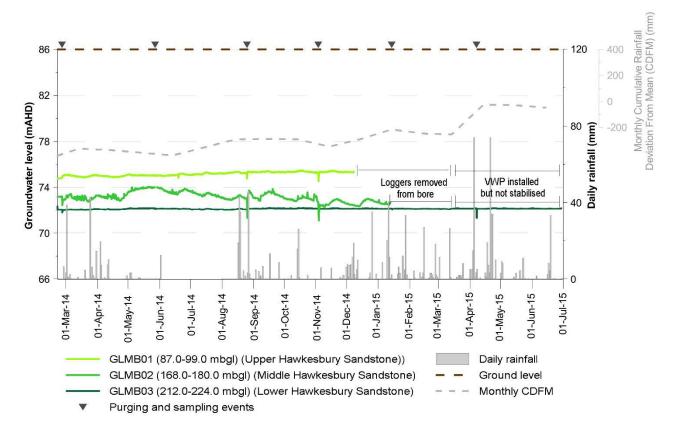


Figure 4.3 Groundwater levels at the Glenlee site

4.1 Temporal trends

4.1.1 Alluvium

The groundwater level in the alluvium (MPMB01) is shallow (less than 10 mbgl) and shows a direct response to rainfall and flood events during the monitoring year (Figure 4.1). The groundwater level increased in response to high rainfall events in August 2014 and April 2015, after which it receded to pre-rainfall level, approximately 1.5 m above river height. The time taken for the water level to recede to pre-rainfall events was over a period of up to two months.

4.1.2 Ashfield Shale

Groundwater levels in RMB01 located in the Ashfield Shale are deep (~80 mbgl) and show no apparent response to rainfall over the monitoring year as shown in Figure 4.1.

The groundwater level in RMB01 shows very slow recovery (longer than 1 year) after installation and after purging and sampling. This slow recovery is due to the very low permeability of the formation. Because of the slow recovery in RMB01 the observed water levels do not reflect the regional groundwater level or trends during the monitoring year.

Since its installation in June 2013, there has been insufficient perched water present in shallow bore RMB04 to allow for water level data to be collected. This monitoring bore was installed to monitor the presence of perched water that may sustain the adjacent Cumberland Plain Woodlands' vegetation.

4.1.3 Hawkesbury Sandstone

Groundwater levels in the Hawkesbury Sandstone show no apparent response to individual rainfall events over the monitoring year at the Denham Court and Glenlee sites (Figure 4.1 and Figure 4.3 respectively), while at the Menangle Park site, located next to the Nepean River, a definite response to rainfall and flood events is observed at the upper and middle Hawkesbury Sandstone monitoring bores (MPMB02 and MPMB03) and a subdued and delayed response in the lower Hawkesbury Sandstone monitoring bore (MPMB04) (Figure 4.2).

Groundwater levels are deep (~40 mbgl) at the Denham Court site and are above the base of the Ashfield Shale indicating confining conditions. The groundwater levels remained fairly constant with fluctuations of less than ~0.5 m throughout the monitoring year. Sudden decreases and slow recovery of groundwater level in the lower Hawkesbury Sandstone (RMB03) are related to sampling events and are consistent with the low permeability of the Hawkesbury Sandstone at this location.

Groundwater levels are shallow (less than 15 mbgl) in the Hawkesbury Sandstone at the Glenlee site. Groundwater levels have been stable over the monitoring year in the upper and lower Hawkesbury Sandstone (GLMB01 and GLMB03) and show a slight decrease in the middle Hawkesbury Sandstone (GLMB02) since the start of this monitoring year. The erratic nature of the data from GLMB02 was investigated and appears to be a natural (Parsons Brinckerhoff 2015a).

At Menangle Park the groundwater level for each of the water bearing zones monitored in the Hawkesbury Sandstone is shallow (less than 10 mbgl). Groundwater levels at the upper and middle Hawkesbury Sandstone monitoring bores (MPMB02 and MPMB03) increased in response to rainfall events in August 2014 and April 2015, followed by a decrease in groundwater levels over a period of two months. The response to rainfall in the lower Hawkesbury Sandstone (MPMB04) indicates slow and delayed recharge after high rainfall periods.

4.2 Spatial trends in the Hawkesbury Sandstone

The conceptual model (AGL 2013) and hydrogeological setting (Section 2.4) suggest that regional groundwater flow within the Hawkesbury Sandstone is from south to north towards the incised river systems of the Sydney Basin.

The groundwater level elevations in the Hawkesbury Sandstone aquifer can be compared between each of the three monitoring sites. The data collected at the CGP suggests that groundwater flow (in the Hawkesbury Sandstone at least) is more complex than the regional conceptual model. The data suggests that:

- The Nepean River in the vicinity of the Menangle Park site is a probable groundwater discharge area (as there is upward groundwater flow within the Hawkesbury Sandstone and there is no Ashfield Shale to act as a cap rock) - groundwater elevations here are between 58 and 61 mAHD and the Nepean River height is typically between 57 and 58 mAHD;
- The Glenlee site may be close to a groundwater divide as groundwater elevations are between 72 and 75 mAHD; and
- The Denham Court site may be on the other side of the groundwater divide as groundwater elevations are lower at around 31 mAHD.

Further studies and data from a larger number of monitoring sites would be required to further understand the regional flow patterns in the Hawkesbury Sandstone.

4.3 Groundwater-surface water interaction

Hydraulic connection between surface water and groundwater exists where the river is in direct contact with the underlying aquifer (Bouwer and Maddock 1997). A 'gaining' stream exists where the water table or groundwater level in a connected aquifer is higher than the running level in a stream and groundwater will flow or discharge to the stream (Land and Water Australia 2007).

The Nepean River level shows a clear response to rainfall (Figure 4.2). The river level is usually lower than the level in the alluvium and Hawkesbury Sandstone units, indicating the river is a gaining river at the Menangle Park site during most of the monitoring period, except for short periods during extremely high rainfall events, when recharge to the alluvial groundwater system occurs.

It is noted that the location of the Nepean River level monitoring station is approximately 1.5 km south of the MPMB bores. This distance, and potential differences in elevation at each site, has not been taken into consideration during this analysis.

4.4 Vertical gradients

Vertical gradients provide an indication of the potential for groundwater to flow vertically upward or downward at that particular location. A downward hydraulic gradient indicates a potential for downward flow from the shallower unit to the deeper unit, while an upward gradient indicates the opposite. It is noted that the actual flow direction and velocity is also governed by permeability.

Potential vertical gradients between the various hydrogeological units were assessed and vary between sites:

- Groundwater levels are comparable in the upper and lower Hawkesbury Sandstone at the Denham Court site indicating no vertical gradient between these monitored zones. Although the sandstone is confined, there is an apparent upward gradient between the upper Hawkesbury Sandstone and the Ashfield Shale.
- There is an apparent upward hydraulic gradient at the Menangle Park site within the monitored zones of the Hawkesbury Sandstone; however a downward gradient exists between the alluvium and the upper

Hawkesbury Sandstone. The similar response to rainfall and flooding events between the alluvial monitoring bore and the Hawkesbury Sandstone monitoring bores indicates connectivity between the two formations at this location, which is expected given the lack of a substantial confining layer (for example shale) between the formations.

 There is an apparent downward hydraulic gradient within the Hawkesbury Sandstone at the Glenlee site.

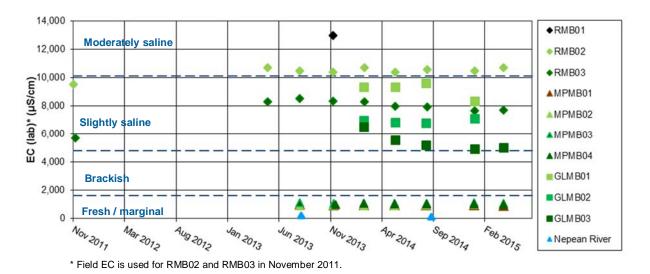
Vertical gradients can be influenced by structural geological features (*i.e.* faults, folds and lineaments) and differences in local porosity, as described in sections 2.3 and 2.4 respectively.

5. Water quality

Water quality monitoring was undertaken between November 2011 and April 2015. Water quality results for the 2014/15 monitoring year are summarised in this chapter and compared to previous monitoring years (Parsons Brinckerhoff 2014a and 2014b). The 2014/15 monitoring year full water quality results are presented in Appendix B and laboratory results in Appendix C.

5.1 Groundwater quality

5.1.1 Field parameters



Time series of field EC and pH for the CGP monitoring bores are presented in Figure 5.1 and Figure 5.2.

Figure 5.1 EC time series for CGP monitoring bores and Nepean River sample

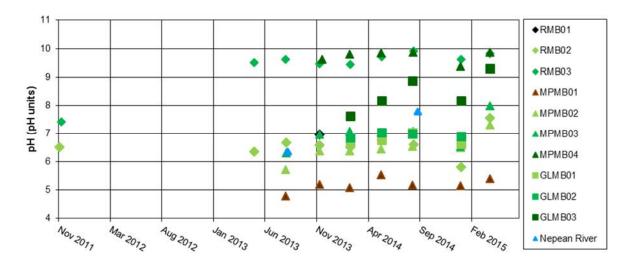


Figure 5.2 pH time series for CGP monitoring bores and Nepean River sample

Groundwater sampled from the alluvium at Menangle Park (MPMB01) is classified as fresh to marginally saline and has a pH of approximately 4.5 to 5.5. The Menangle Park site is a former sand and gravel quarry

that has been subsequently rehabilitated. The observed low pH may be related to these previous land use activities.

No groundwater sample was collected from the Ashfield Shale monitoring bore at Denham Court (RMB01) due to insufficient water in the monitoring bore during the 2014/15 monitoring year.

Groundwater in the Hawkesbury Sandstone at the Menangle Park site (MPMB02-04) is classified as fresh to marginal, while slightly saline conditions are observed at the Glenlee site (GLMB01-03) and slightly saline to moderately saline conditions are observed at the Denham Court site (RMB02-03). The fresh to marginal conditions at the Menangle Park site are likely due to the influence of rainfall recharge and connectivity with the Nepean River.

Salinity within the Hawkesbury Sandstone does not show a clear depth related trend at Menangle Park, however, salinity decreases with depth at the Denham Court and Glenlee sites. This decrease is likely a result of saline groundwater within the Ashfield Shale migrating into the underlying aquifer as a result of the leakage. The pH generally increases with depth within the Hawkesbury Sandstone. The pH has increased in the Hawkesbury Sandstone at the Menangle Park and Glenlee sites over the monitoring period.

5.1.2 Major ions

The major ion characteristics of groundwater samples for this monitoring year are shown in a piper diagram and representative bivariate plots in Figure 5.3. A piper diagram is a graphical representation of the relative concentrations of major ions (Ca^{2+} , Mg^{2+} , Na^+ , K^+ , Cl^- , HCO_3^- , CO_3^{2-} and SO_4^{2-}). The ratios of sodium/chloride and magnesium/chloride versus chloride concentrations are also presented in two bivariate plots. Chloride is typically assumed to be a conservative (non-reactive) ion in groundwater systems. Evapotranspiration of the initial water with low chloride concentration would therefore be expected to result in a horizontal trend in a major ion/chloride versus chloride plot.

Groundwater at monitoring bores is typically dominated by sodium and chloride, although, groundwater at the Menangle Park site is a mixed ion, Na-Mg-Ca-HCO₃ type water. Interestingly the water type at each of the four nested monitoring bores at the Menangle Park site is different.

Major ion/chloride ratios of groundwater at the Denham Court and Glenlee sites are similar and generally close to those of typical sea water (Turekian 1968). The Hawkesbury Sandstone at the Menangle Park site has a different geochemical signature to the Denham Court and Glenlee sites, which may be related to local recharge and connectivity with the overlying alluvium at this site.

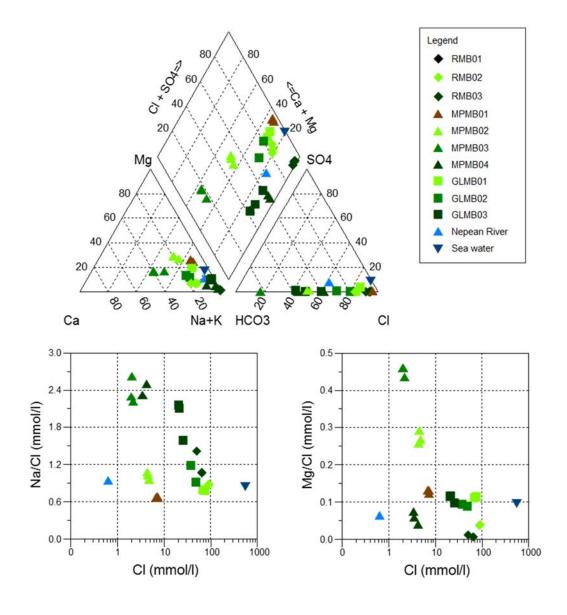


Figure 5.3 Major ion chemistry of groundwater for CGP monitoring bores and Nepean River sample (2014/15 monitoring year)

5.1.3 Dissolved metals

Concentrations of dissolved metals in groundwater are presented in Figure 5.4. The major findings for dissolved metals for this monitoring year are as follows:

- Dissolved metal concentrations are generally similar in the alluvium and the Hawkesbury Sandstone. The exceptions are copper, cobalt, nickel and lead concentrations, which are higher in the alluvium compared to the other hydrogeological units. Dissolved metal concentrations in the alluvium were generally comparable to previous monitoring years.
- Slightly elevated dissolved metal concentrations were detected in the Hawkesbury Sandstone at the Denham Court and Glenlee sites (*e.g.* barium, bromine and strontium) compared with other locations. This is not unexpected given the higher salinity at these locations and the influence from the overlying Ashfield Shale. Elevated concentrations of barium, cadmium, molybdenum, strontium and zinc are not uncommon for groundwater in the Hawkesbury Sandstone (Parsons Brinckerhoff 2006 and 2013b). Dissolved metal concentrations in the Hawkesbury Sandstone were generally comparable to previous monitoring years, with the exception of decreasing iron concentrations at RMB02, decreasing molybdenum concentrations at MPMB04 and GLMB03. Mercury was also detected at MPMB04 just above the laboratory LOR in April 2015 at 0.0003 mg/L.

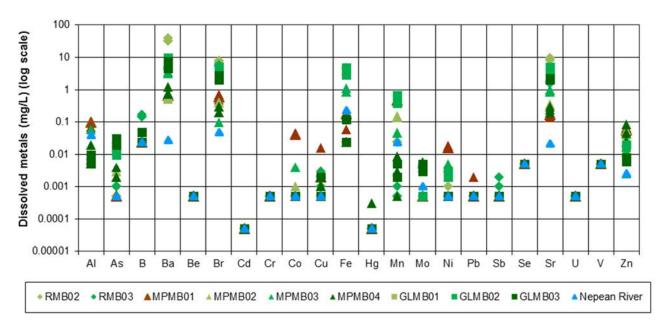


Figure 5.4 Dissolved metal concentrations in groundwater for CGP monitoring bores (2014/15 monitoring year)

5.1.4 Nutrients

A plot showing ammonia versus nitrate in groundwater is presented in Figure 5.5.

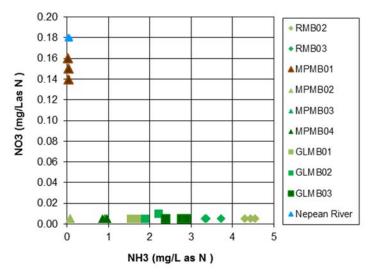


Figure 5.5 Ammonia versus nitrate concentrations in groundwater for CGP monitoring bores and the Nepean River sample (2014/15 monitoring year)

The major findings for nutrients are as follows:

- Nitrate (NO₃) concentrations in groundwater remained low at all CGP monitoring bores (<0.2 mg/L as N) compared to previous monitoring years. The highest concentrations continued to be recorded in the alluvial aquifer (MPMB01) (Figure 5.5).
- Ammonia (NH₃) concentrations were the lowest at Menangle Park and the highest at Denham Court, which is consistent with previous monitoring years (Figure 5.5).
- Nitrite concentrations remained below the laboratory LOR at all monitoring bores with the exception of a detection in the Hawkesbury Sandston at Denham Court (RMB02-03) in August 2014.

 Total organic carbon (TOC) concentrations differ between hydrogeological units, being generally the highest in the lower Hawkesbury Sandstone at all sites and comparable in the middle and upper Hawkesbury Sandstone and alluvium.

5.1.5 Dissolved gases

A time series plot of dissolved methane concentrations in groundwater is presented in Figure 5.6.

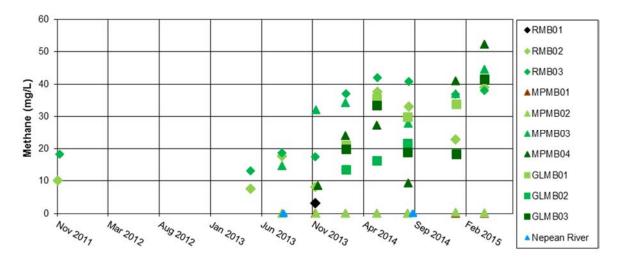


Figure 5.6 Dissolved methane (CH₄) in groundwater time series for CGP monitoring bores

The major findings for dissolved gases are as follows:

- Dissolved gases were not detected in the alluvium (MPMB01).
- Dissolved methane was detected in the Hawkesbury Sandstone at all monitoring sites, with increased concentrations at depth. Dissolved methane was also detected at the control site (RMB) remote from development activities. Dissolved methane concentrations continued to be the lowest at MPMB02. Dissolved methane concentrations at the majority of sites in the Hawkesbury Sandstone continued to increase during this monitoring year and the April 2015 concentrations were the highest concentrations recorded at most monitoring bores since monitoring commenced (Figure 5.6).
- Ethane continued to be detected at low concentrations in groundwater from the Hawkesbury Sandstone at the Denham Court and Glenlee sites and propane and butane continued to be detected at low concentrations at the Glenlee site only. Concentrations of these dissolved gases overall decreased at the Glenlee site since monitoring commenced in February 2014.

Dissolved methane is shown to be of mostly thermogenic origin (Parsons Brinckerhoff 2014). The presence of dissolved hydrocarbons observed in the groundwater within the Hawkesbury Sandstone and Ashfield Shale is assessed to be naturally occurring, based on the values present within the groundwater at the control site (Denham Court, RMB) located at significant distance from the CGP gas production wells. It is likely that the increasing dissolved methane trends are related to degassing of naturally occurring methane after purging during groundwater sampling events. Future trends will be monitored.

5.1.6 Dissolved hydrocarbons

The Menangle Park site is a former sand and gravel quarry that has been subsequently rehabilitated, and hydrocarbon detections (PAHs and TPH/TRH) at this site in the alluvium may be related to these previous land use activities.

Toluene continued to be present in most sandstone units at all sites in variable concentrations. No other BTEX compounds (*i.e.* benzene, xylenes and ethyl benzene) were detected during this monitoring year. Toluene is assessed to be naturally occurring, given that it has been detected in groundwater at all

monitoring sites since monitoring began, including the control site (Denham Court, RMB) located at a significant distance from development activities.

Dissolved hydrocarbons can occur naturally in groundwater, with concentrations derived from any carbonaceous material (CSIRO 2011). Detections of PAHs, phenols and heavy chain hydrocarbons in both upper and lower Hawkesbury Sandstone across all sites are most likely natural, however possible residues from monitoring bore drilling (*e.g.* lubricating oils) cannot be excluded. Future trends will be monitored.

5.2 Surface water quality

Surface water quality results from a single sample collected from the Nepean River on 9 September 2014 were compared to ANZECC (2000) guidelines for freshwater ecosystems (95% protection level) and it was found that:

- pH is neutral, within the ANZECC (2000) guideline range (6.5 8.0 pH units) and higher than the pH of groundwater in the alluvium.
- Salinity is fresh, and within the ANZECC (2000) guideline range (125 2,200 µS/cm) and lower than groundwater in the alluvium. The low salinity is likely the result of the rainfall that preceded the collection of the sample.
- Dominant major ions are sodium, chloride and bicarbonate (Figure 5.3).
- Dissolved metal concentrations were lower than those of groundwater in the alluvium and underlying Hawkesbury Sandstone units, except for iron and molybdenum concentrations. All dissolved metal concentrations were below the ANZECC (2000) guideline values.
- Ammonia concentrations were above the ANZECC (2000) guideline (0.02 mg/L) and are similar to the alluvial groundwater. Total phosphorus concentration was lower than the August 2013 concentration and was below the ANZECC (2000) guideline value (0.05 mg/L).
- Dissolved methane was detected at low concentrations (0.01 mg/L).
- No other dissolved hydrocarbons were detected.

The water quality of the Nepean River was overall comparable to that of the sample collected in August 2013.

6. Discussion and conclusions

Monitoring of groundwater levels at three nested monitoring bore sites was undertaken using dataloggers, allowing water level trends to be identified in the alluvium, Ashfield Shale and Hawkesbury Sandstone. Sampling of water quality at all sites also established useful trends.

The main findings for this monitoring year in regards to water levels are:

- The groundwater level in the alluvium at Menangle Park is less than 10 mbgl and shows a direct response to rainfall and flood events.
- Groundwater levels appear to follow similar trends in each of the Hawkesbury Sandstone units (upper, middle and lower) at each site. There is no apparent response to individual rainfall events at the Denham Court and Glenlee sites, while a clear response to rainfall and flood events can be observed at most monitoring bores at the Menangle Park site even though this is an apparent groundwater discharge area.
- Groundwater levels are deep (~40 mbgl) at the Denham Court site, although they are above the base of the Ashfield Shale indicating confining conditions. In contrast, groundwater levels are shallow (less than 15 mbgl) at the Menangle Park and Glenlee sites.
- For the regional Hawkesbury Sandstone aquifer, groundwater elevations are highest at the Glenlee site (~75 mAHD); lower at the Menangle Park site (~60 mAHD); and lowest at the Denham Court site (~31 mAHD).
- Vertical gradients vary between sites; no vertical gradient is apparent between the lower and upper Hawkesbury Sandstone at the Denham Court site, whereas upward and downward gradients are evident at Menangle Park and Glenlee respectively.
- The Nepean River elevation is usually lower than the water elevation in the alluvium and Hawkesbury Sandstone units, indicating the river is a gaining stream around the Menangle Park site, except for short periods during extremely high rainfall events when recharge to the underlying groundwater systems is observed.
- The groundwater level data collected in the alluvium and Hawkesbury Sandstone is indicative of natural systems in long-term equilibrium with occasional seasonal responses to recharge when there is a connection with surface features, as evident at the Menangle site within the alluvium and Hawkesbury Sandstone.

No downward temporal trends have been observed in the groundwater level data at any of the monitored locations and therefore it can be assumed that no depressurisation due to CSG operations (which involve dewatering of the coal seam) is occurring within the monitored zones.

The main findings for this monitoring year in regards to water quality are:

- Groundwater quality in the alluvium at the Menangle Park site is characterised as fresh to marginally saline and slightly acidic pH. Dissolved metal concentrations are typically low and no dissolved gases were detected. Minor dissolved hydrocarbons were detected and may be related to previous land use activities.
- No sample was collected from the Ashfield Shale at Denham Court due to insufficient water in the monitoring bore.
- Groundwater quality in the Hawkesbury Sandstone ranges from fresh to marginally saline at the Menangle Park site and is slightly to moderately saline at the Denham Court and Glenlee sites. Salinity decreases with depth at the Denham Court and Glenlee sites. Minor detections of dissolved hydrocarbons were present in the Hawkesbury Sandstone. Dissolved methane was detected at all

Hawkesbury Sandstone bores and it is likely related to degassing of naturally occurring methane after purging during groundwater sampling events. Dissolved ethane was detected at the Denham Court and Glenlee sites and propane and butane were detected at the Glenlee site only. These compounds are assessed to be naturally occurring given that methane has been observed to occur at all sites, including the control site (Denham Court) which is located at a significant distance from any development activities.

- Toluene was present at all sites (except the alluvium and the shallowest sandstone interval at Menangle Park) at variable concentrations. No other BTEX compounds were detected. Toluene is assessed to be naturally occurring given that it has been detected at all sites, including the control site (Denham Court) which is located at a significant distance from the CGP gas production wells.
- The Nepean River at the Menangle Park site is of neutral pH and fresh water quality. Dissolved metal concentrations were typically lower for the surface water than the alluvium and underlying Hawkesbury Sandstone. Ammonia concentrations were above the ANZECC (2000) guideline values and similar to alluvial groundwater. Minor dissolved methane concentrations were detected however no other hydrocarbons were detected.
- No significant change in water quality was detected during the 2014/15 monitoring year compared to the previous monitoring year (Parsons Brinckerhoff 2014b).

No adverse water quality impacts that can be attributed to CSG operations were observed at any of the monitored locations. Water quality results are not significantly different between the control site (Denham Court) and the Stage 2 area where CSG operations are occurring (Menangle Park and Glenlee).

From the available data, there are no observable impacts to groundwater levels or quality that could be attributable to the CSG operations. There is also no evidence of connectivity between the shallower monitored zones and the coal seams. This is in agreement with the conceptual model developed during the Phase 1 studies (Parsons Brinckerhoff 2011). The presence of extensive and thick claystone formations (aquitards and aquicludes) between the Hawkesbury Sandstone and the targeted coal seams restricts upward depressurisation and impedes the vertical flow of groundwater.

7. Statement of limitations

7.1 Scope of services

This report has been prepared in accordance with the scope of services set out in the contract, or as otherwise agreed, between the client (AGL) and Parsons Brinckerhoff (scope of services). In some circumstances the scope of services may have been limited by a range of factors such as time, budget, access and/or site disturbance constraints.

7.2 Reliance on data

In preparing the report, Parsons Brinckerhoff has relied upon data, surveys, plans and other information provided by the client and other individuals and organisations, most of which are referred to in the report (the data). Except as otherwise stated in the report, Parsons Brinckerhoff has not verified the accuracy or completeness of the data. To the extent that the statements, opinions, facts, information, conclusions and/or recommendations in the report (conclusions) are based in whole or part on the data, those conclusions are contingent upon the accuracy and completeness of the data. Parsons Brinckerhoff will not be liable in relation to incorrect conclusions should any data, information or condition be incorrect or have been concealed, withheld, misrepresented or otherwise not fully disclosed to Parsons Brinckerhoff.

7.3 Environmental conclusions

In accordance with the scope of services, Parsons Brinckerhoff has relied upon the data and has conducted environmental field monitoring and/or testing in the preparation of the report. The nature and extent of monitoring and/or testing conducted is described in the report.

On all sites, varying degrees of non-uniformity of the vertical and horizontal soil or groundwater conditions are encountered. Hence no monitoring, common testing or sampling technique can eliminate the possibility that monitoring or testing results/samples are not totally representative of soil and/or groundwater conditions encountered. The conclusions are based upon the data and the environmental field monitoring and/or testing and are therefore merely indicative of the environmental condition of the site at the time of preparing the report, including the presence or otherwise of contaminants or emissions.

Within the limitations imposed by the scope of services, the monitoring, testing, sampling and preparation of this report have been undertaken and performed in a professional manner, in accordance with generally accepted practices and using a degree of skill and care ordinarily exercised by reputable environmental consultants under similar circumstances. No other warranty, expressed or implied, is made.

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The report has been prepared for the benefit of the client (and no other party). Parsons Brinckerhoff assumes no responsibility and will not be liable to any other person or organisation for or in relation to any matter dealt with or conclusions expressed in the report, or for any loss or damage suffered by any other person or organisation arising from matters dealt with or conclusions expressed in the report (including without limitation matters arising from any negligent act or omission of Parsons Brinckerhoff or for any loss or damage suffered by any other party relying upon the matters dealt with or conclusions expressed in the report). Parties other than the client should not rely upon the report or the accuracy or completeness of any conclusions and should make their own enquiries and obtain independent advice in relation to such matters.

7.5 Other limitations

Parsons Brinckerhoff will not be liable to update or revise the report to take into account any events or emergent circumstances or facts occurring or becoming apparent after the date of the report.

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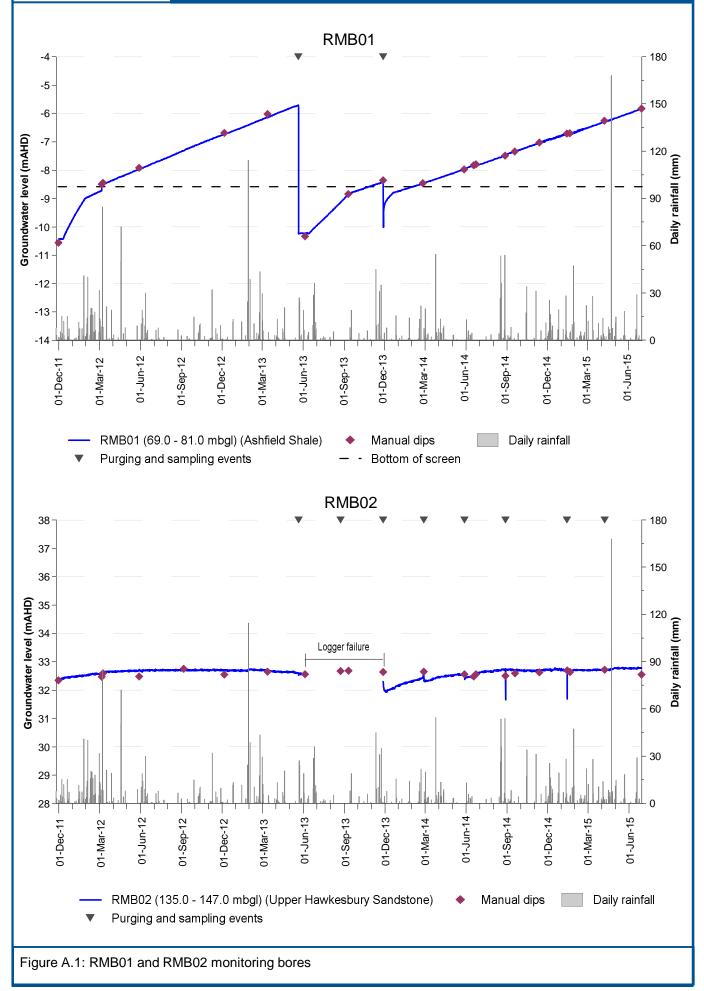
Appendix A

Groundwater hydrographs



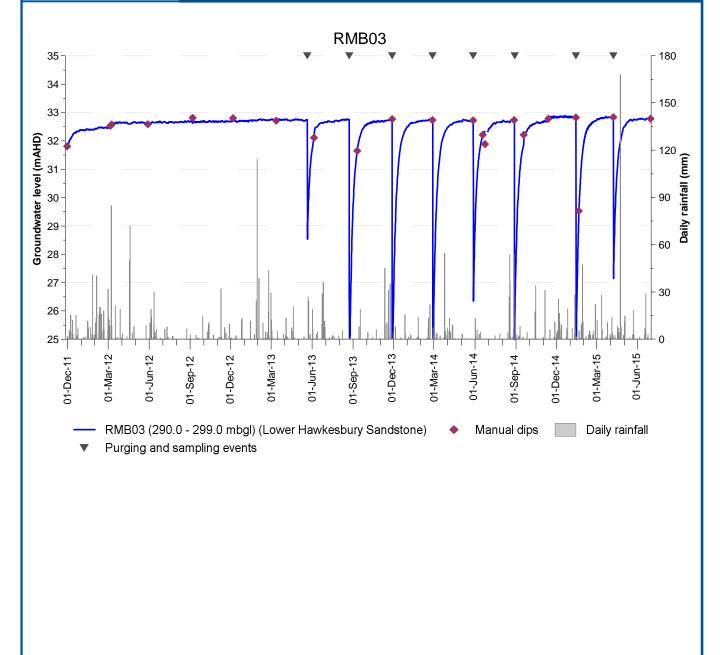
PARSONS BRINCKERHOFF

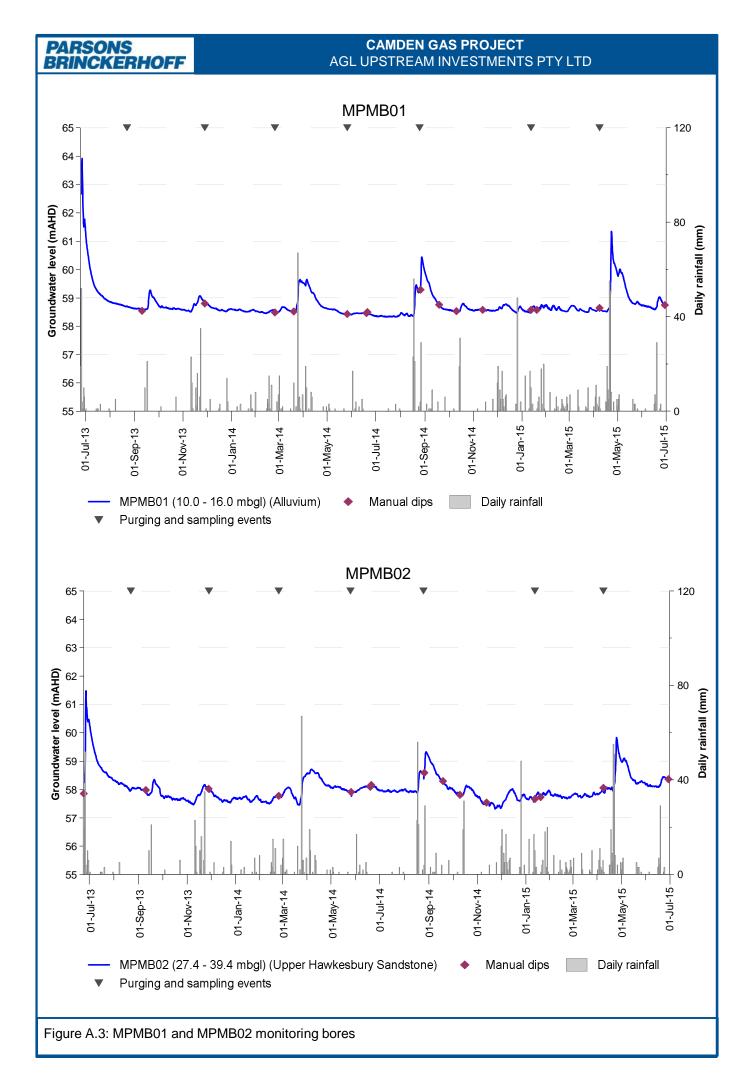
CAMDEN GAS PROJECT AGL UPSTREAM INVESTMENTS PTY LTD

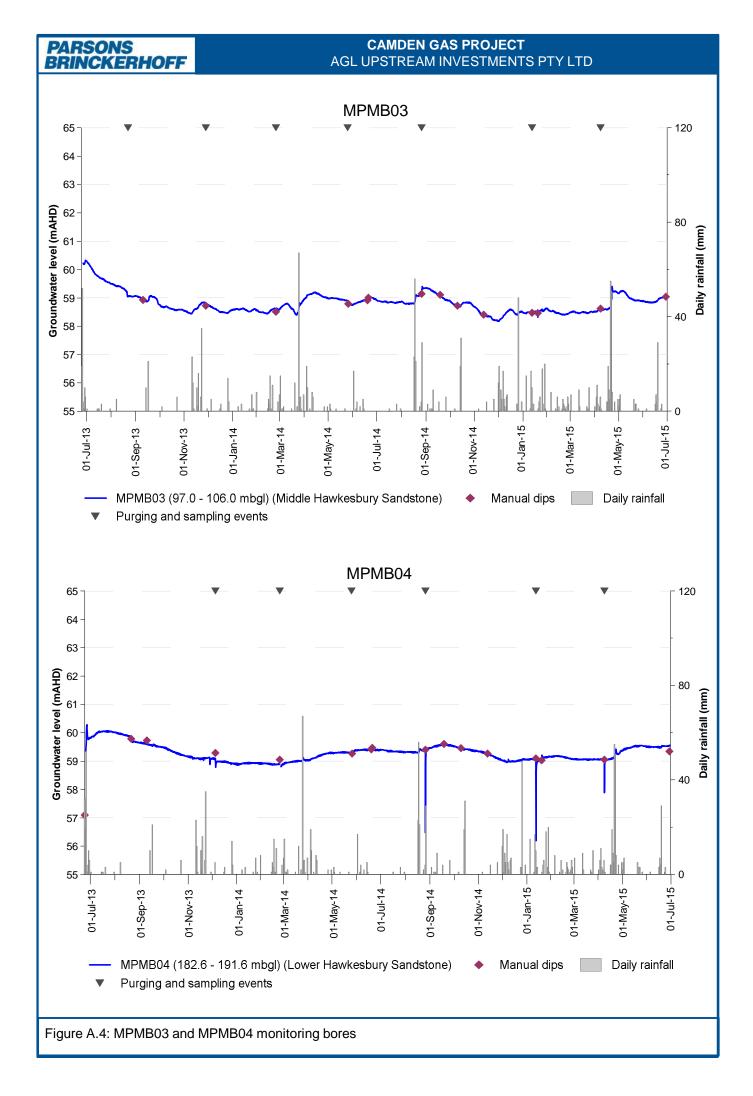


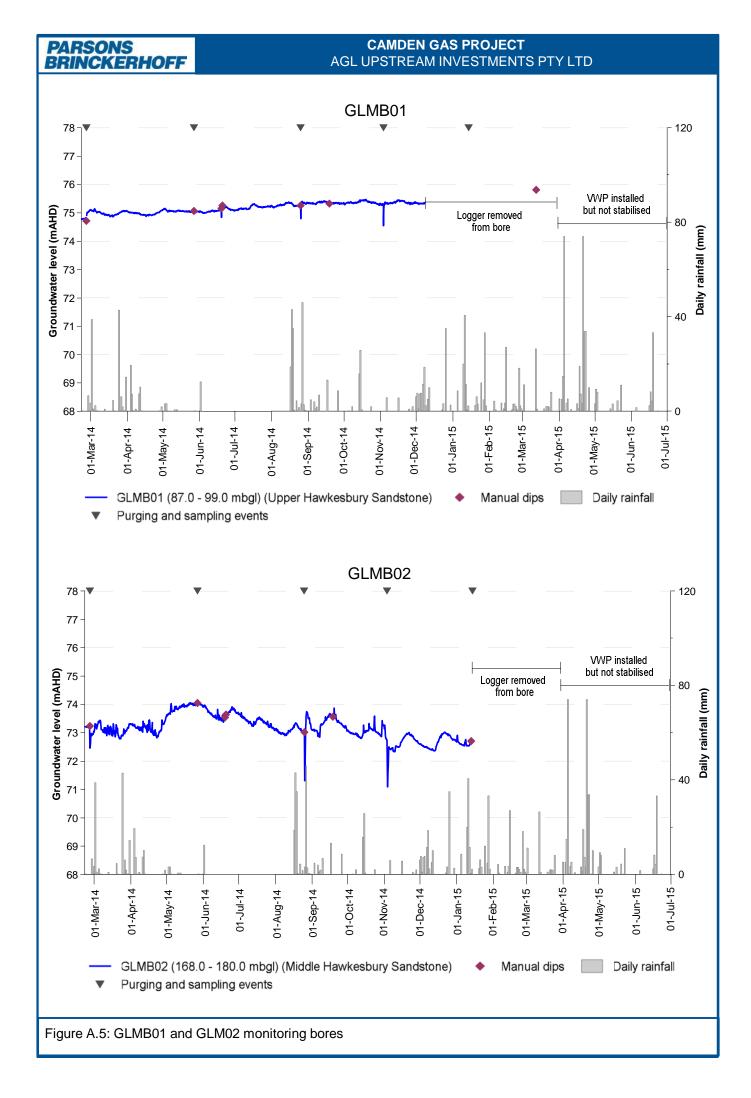
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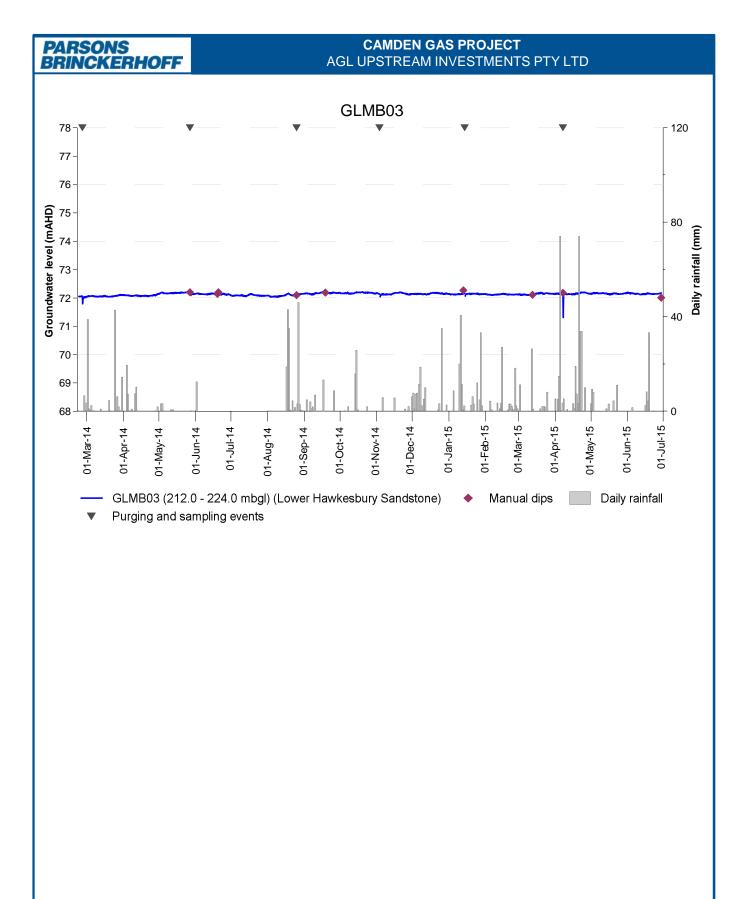
CAMDEN GAS PROJECT AGL UPSTREAM INVESTMENTS PTY LTD











Appendix B

Water quality summary tables



SUMMARY TABLE B.1: Water quality results Denham Court and Glenlee sites (2014-2015)

| Hydrogeological unit | | | Hawkes | RMB02 bury Sandstor | e (upper) | Hawkes | RMB03 bury Sandstor | ne (lower) | Hawkesbur | MB01 y Sandstone | Hawkesbur | VB02 Ty Sandstone | Indstone Hawkesbury Sandstone (lower) | | |
|---|--------------|----------------|-------------------|------------------------|-------------------|------------------|------------------------|------------------|-------------------|---------------------|----------------------------|----------------------|---------------------------------------|-----------------------------|----------------|
| Sample date | | Territori | 29/08/2014 | 13/01/2015 | 7/04/2015 | 29/08/2014 | 13/01/2015 | 7/04/2015 | (up 25/08/2014 | per) 14/01/2015 | (mi 25/08/2014 | ddle) 14/01/2015 | 25/08/2014 | 14/01/2015 | 7/04/2015 |
| General parameters | Units | Typical LOR | | | | | | | | | | | | | |
| H (field) H (lab) | pH units | 0.01 | 6.60 7.18 | 5.81 7.39 | 7.55 7.19 | 9.92 9.48 | 9.61 9.60 | 9.81 9.63 | 7.03 7.40 | 6.63 7.48 | 7.00 7.39 | 6.90 7.66 | 8.87 8.77 | 8.17 8.85 | 9.31 8.60 |
| lectrical conductivity (field) | µS/cm | 1 | 10,959 | 9,831 | 10,390 | 8,296 | 7,522 | 7,486 | 9,332 | 7,675 | 6,545 | 6,902 | 5,017 | 4,704 | 4,842 |
| enperature | °C | 0.01 | 10,600 20.55 | 10,500 22.30 | 10,700 19.46 | 7,940 20.46 | 7,690 20.96 | 7,730 19.45 | 9,640 19.73 | 8,360 24.53 | 6,810 20.82 | 7,140 27.41 | 5,200 20.25 | 4,930 24.53 | 5,010 20.47 |
| Dissolved oxygen Total dissolved solids (field) | % sat | 0.01 | 10.30 7,123 | 6.30 6,391 | - 6,753 | 20.10 5,389 | 2.40 4,889 | 5.10 4,865 | 42.90 6,066 | 3.60 4,986 | 1.40 4,253 | 5.00 4,484 | 2.70 3,260 | 1.40 3,057 | 2.40 3,151 |
| Fotal dissolved solids (lab) | mg/L | 10 | 5,440 | 5,980 | 6,170 | 3,830 | 4,130 | 4,140 | 5,320 | 5,110 | 3,360 | 4,320 | 2,460 | 2,560 | 2,750 |
| Fotal suspended solids Redox | mg/L mV | 5 0.1 | <5 -164.1 | <5 -132.1 | <5 -41.5 | <5 25.8 | <5 -157.1 | <5 -94.1 | <5 -305.5 | 22 -168.9 | 36 -241.8 | 41 -204.7 | <5 -221.4 | 6 -243.1 | <5 -209.0 |
| Nater type ^a | - | | Na-Cl | Na-Cl | Na-Cl | Na-Cl | Na-Cl | Na-Cl | Na-Cl | Na-Cl | Na-Cl- HCO ₃ | Na-Cl | Na-Cl- HCO ₃ | Na-HCO ₃ - Cl | Na-HCC Cl |
| _aboratory analytes | | | | | | | | | | | - | | | | |
| Hydroxide alkalinity as CaCO ₃ | mg/L mg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 298 | <1 229 | <1 252 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 184 | <1 309 | <1 129 |
| Bicarbonate alkalinity as CaCO ₃ | mg/L | 1 | 972 972 | 792 792 | 895 895 | 239 537 | 189 418 | 224 476 | 683 683 | 434 434 | 793 793 | 595 595 | 876 1,060 | 1,070 1,380 | 1,510 1,640 |
| Fotal alkalinity as CaCO ₃ Sulfate as SO ₄ ²⁻ | mg/L mg/L | 1 | <10 | <1 | <10 | <1 | <1 | <1 | <1 | 135 | 5 | <1 | <1 | <1 | 19 |
| Chloride Calcium | mg/L mg/L | 1 | 3,120 270 | 3,170 382 | 3,280 326 | 1,760 6 | 2,290 14 | 2,310 10 | 2,710 174 | 2,470 182 | 1,370 210 | 1,760 246 | 934 16 | 734 24 | 774 31 |
| Magnesium | mg/L | 1 | 81 | 83 | 88 | 14 | 11 | 11 | 211 | 186 | 87 | 105 | 61 | 57 | 61 |
| Sodium Potassium | mg/L mg/L | 1 | 1,740 37 | 1,770 30 | 1,890 30 | 1,610 22 | 1,570 16 | 1,600 16 | 1,360 24 | 1,260 19 | 1,040 29 | 1,040 24 | 957 32 | 1,020 36 | 1,050 39 |
| Silica (reactive) Fluoride | mg/L mg/L | 0.05 | 11.6 0.2 | 9.66 0.2 | 10.4 0.1 | 7.30 0.4 | 6.24 0.4 | 6.68 0.6 | 10.6 0.3 | 7.38 0.2 | 21.2 0.3 | 16.0 0.3 | 9.77 0.2 | 11.0 <0.1 | 11.8 <0.1 |
| Bromide | mg/L | 0.01 | 5.88 | 8.89 | 6.79 | 4.71 | 6.05 | 4.68 | 6.06 | 8.30 | 3.98 | 5.75 | 2.35 | 2.25 | 1.98 |
| Γotal cyanide Dissolved metals | mg/L | 0.0004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 |
| Aluminium | mg/L | 0.01 | <0.01 <0.001 | <0.01 <0.001 | <0.01 <0.001 | <0.01 <0.001 | <0.01 0.001 | 0.01 | <0.01 <0.001 | <0.01 <0.001 | <0.01 <0.001 | <0.01 <0.001 | 0.01 <0.001 | 0.01 | <0.01 |
| Antimony Arsenic | mg/L mg/L | 0.001 | <0.001 | <0.001 | <0.001 | 0.001 | 0.001 | 0.001 | <0.001 | 0.010 | 0.014 | 0.010 | 0.020 | 0.020 | 0.032 |
| Barium Beryllium | mg/L mg/L | 0.001 | 40.0 <0.001 | 33.2 <0.001 | 33.0 <0.001 | 4.35 <0.001 | 5.07 <0.001 | 4.64 <0.001 | 2.64 <0.001 | 0.362 <0.001 | 9.7800 <0.001 | 8.78 <0.001 | 4.49 <0.001 | 6.28 <0.001 | 7.63 <0.00 |
| Cadmium | mg/L | 0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.000 |
| Chromium Cobalt | mg/L mg/L | 0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.00 |
| Copper | mg/L | 0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | 0.002 <0.001 | 0.001 <0.001 | 0.003 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | 0.002 |
| Lead Manganese | mg/L mg/L | 0.001 | 0.026 | 0.024 | 0.024 | <0.001 | 0.001 | <0.001 | 0.300 | 0.390 | 0.666 | 0.384 | 0.002 | 0.003 | 0.005 |
| Mercury Molybdenum | mg/L mg/L | 0.0001 | <0.0001 <0.001 | <0.0001 <0.001 | <0.0001 <0.001 | <0.0001 0.004 | <0.0001 0.004 | <0.0001 0.005 | <0.0001 <0.001 | <0.0001 <0.001 | <0.0001 <0.001 | <0.0001 <0.001 | <0.0001 0.005 | <0.0001 0.004 | <0.000 |
| Nickel | mg/L | 0.001 | <0.001 | 0.003 | 0.001 | <0.001 | <0.001 | <0.001 | 0.002 | <0.001 | 0.002 | 0.003 | <0.001 | <0.001 | < 0.00 |
| Selenium Strontium | mg/L mg/L | 0.01 0.001 | <0.01 8.18 | <0.01 9.97 | <0.01 8.47 | <0.01 1.64 | <0.01 1.83 | <0.01 1.58 | <0.01 6.33 | <0.01 4.08 | <0.01 5.14 | <0.01 3.97 | <0.01 2.06 | <0.01 1.54 | <0.01 2.23 |
| Jranium /anadium | mg/L | 0.001 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.001 <0.01 | <0.00 |
| /anadium Zinc | mg/L mg/L | 0.005 | 0.017 | 0.033 | 0.031 | 0.053 | 0.039 | 0.051 | <0.005 | <0.005 | 0.020 | 0.006 | 0.006 | <0.005 | 0.008 |
| Boron | mg/L mg/L | 0.05 | <0.05 4.81 | <0.05 4.73 | <0.05 4.42 | 0.15 <0.05 | 0.15 <0.05 | 0.17 <0.05 | <0.05 0.19 | <0.05 4.28 | <0.05 2.97 | <0.05 4.87 | <0.05 <0.05 | <0.05 0.10 | 0.05 |
| Bromine | mg/L | 0.1 | 7.9 | 7.8 | 7.0 | 5.8 | 5.7 | 5.4 | 7.2 | 6.4 | 4.5 | 5.1 | 2.7 | 2.2 | 2.0 |
| Nutrients Ammonia as N | mg/L | 0.01 | 4.30 | 4.55 | 4.43 | 3.73 | 3.36 | 3.33 | 1.71 | 1.54 | 1.88 | 2.20 | 2.37 | 2.76 | 2.87 |
| Nitrite as N Nitrate as N | mg/L mg/L | 0.01 | 0.04 <0.01 | <0.01 <0.01 | <0.01 <0.01 | 0.04 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 |
| Fotal phosphorus | mg/L | 0.01 | <0.01 | 0.02 | 0.04 | <0.01 | <0.01 | 0.01 | 0.08 | 0.04 | 0.12 | 0.07 | 0.04 | <0.01 | 0.04 |
| Reactive phosphorus Fotal organic carbon | mg/L mg/L | 0.01 | <0.01 <1 | 0.03 <1 | 0.03 5 | <0.01 28 | <0.01 38 | 0.01 40 | 0.06 3 | <0.01 <1 | 0.06 9 | 0.05 <1 | <0.01 10 | 0.03 <1 | 0.05 10 |
| Dissolved gases Methane | mg/L | 0.01 | 33.10 | 22.80 | 39.10 | 40.90 | 36.80 | 38.10 | 29.90 | 33.90 | 21.70 | 18.40 | 18.80 | 18.20 | 41.70 |
| Ethane Ethene | mg/L | 0.01 | 0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | 0.01 <0.01 | 0.01 <0.01 | 1.48 <0.01 | 1.65 <0.01 | 0.93 <0.01 | 0.80 <0.01 | 0.12 <0.01 | 0.06 <0.01 | 0.11 <0.01 |
| Propane | mg/L mg/L | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | 0.42 | 0.38 | 0.19 | 0.20 | 0.03 | 0.02 | 0.02 |
| Propene Butane | mg/L mg/L | 0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 0.07 | <0.01 0.06 | <0.01 0.04 | <0.01 0.04 | <0.01 <0.01 | <0.01 <0.01 | <0.01 |
| Butene | mg/L | 0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Phenolic compounds Phenol | μg/L | 1 | <1 | <1 | <1 | <1 | <1 | <1 | 2 | <1 | 1.8 | <1 | 1.5 | <1 | <1 |
| 2-Chlorophenol 2-Methylphenol | μg/L μg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| 3-&4-Methylphenol | µg/L | 2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 <1 | <2 <1 | <2 | <2 |
| 2-Nitrophenol 2.4-Dimethylphenol | μg/L μg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 | <1 | <1 <1 | <1 <1 |
| 2.4-Dichlorophenol 2.6-Dichlorophenol | μg/L μg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| 4-Chloro-3-Methylphenol | µg/L | 1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| 2.4.6-Trichlorophenol 2.4.5-Trichlorophenol | μg/L μg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Pentachlorophenol Polycyclic aromatic hydrocarbons | μg/L | 2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 |
| Acenaphthylene | µg/L | 1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| Acenaphthene Fluorene | μg/L μg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Phenanthrene Anthracene | μg/L μg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Fluoranthene | µg/L | 1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| ⁵ yrene Benz(a)anthracene | μg/L μg/L | 1 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Chrysene Benzo(b)fluoranthene | μg/L μg/L | 1 | <1 <1 | <1 - | <1 - | <1 <1 | <1 | <1 | <1 <1 | <1 - | <1 <1 | <1 | <1 <1 | <1 | <1 |
| 3enzo(k)fluoranthene | µg/L | 1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| Benzo(a)pyrene ndeno(1.2.3.cd)pyrene | μg/L μg/L | 0.5 1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 | <0.5 <1 |
| Dibenz(a.h)anthracene | µg/L | 1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Benzo(g.h.i)perylene Sum of PAHs | µg/L ug/L | 1 0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 |
| Fotal petroleum hydrocarbons C ₆ -C ₉ fraction | µg/L | 20 | <20 | <20 | <20 | 70 | 90 | 70 | 210 | 110 | 210 | 80 | 310 | 210 | 190 |
| C ₁₀ -C ₁₄ fraction | µg/L | 50 | <50 <100 | <50 <100 | <50 <100 | 90 <100 | <50 <100 | <50 <100 | <50 <100 | <50 <100 | <50 <100 | <50 <100 | <50 <100 | <50 <100 | <50 <100 |
| C ₁₅ -C ₂₈ fraction C ₂₉ -C ₃₆ fraction | μg/L μg/L | 100 50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 |
| C ₁₀ -C ₃₆ fraction (sum) Fotal recoverable hydrocarbons | µg/L | 50 | <50 | <50 | <50 | 90 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 | <50 |
| C ₆ -C ₁₀ fraction | µg/L | 20 | <20 | <20 | <20 | 70 | 90 | 70 | 290 | 110 | 220 | 80 | 310 | 210 | 190 |
| C ₆ -C ₁₀ fraction minus BTEX -C ₁₀ -C ₁₆ fraction | μg/L μg/L | 20 100 | <20 <100 | <20 <100 | <20 <100 | 20 <100 | 40 <100 | 30 <100 | 290 <100 | 90 <100 | 210 <100 | 80 <100 | 170 <100 | 100 <100 | 70 <100 |
| C ₁₆ -C ₃₄ fraction | µg/L | 100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 |
| >C ₃₄ -C ₄₀ fraction >C ₁₀ -C ₄₀ fraction (sum) | μg/L μg/L | 100 100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 |
| Aromatic hydrocarbons Benzene | μg/L | 1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| Foluene | µg/L | 2 | 8 | <2 | <2 | 48 | 46 | 39 | <2 | 19 | 10 | <2 | 139 | 106 | 124 |
| Ethylbenzene n&p-Xylenes | μg/L μg/L | 2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 |
| | μg/L | 2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 <2 |
| • | 110/ | 2 | e? | e) | <i> </i> | e9 | e) | - | ~ ~ | ~ ~ / | | 21 | 2) | د٢ | |
| Xylenes Total xlyenes Sum of BTEX | μg/L μg/L | 2 | <2 8 <1 | <2 <1 <1 | <2 <1 <1 | <2 48 <1 | <2 46 <1 | <2 39 <1 | <2 <1 <1 | <2 19 <1 | <2 10 <1 | <2 <1 <1 | <2 139 <1 | <2 106 <1 | 124 |

SUMMARY TABLE B.2: Water quality results Menangle Park site and Nepean River (2014-2015)

| | | ANZECC 2000 guidelines ^b MPMB01 | | | | | Menangle Park MPMB02 MPMB03 | | | | | | Nepear River | | | |
|--|--------------|---|--------------------------|-----------------|-----------------|-----------------|--------------------------------|----------------------------------|-------------------------------|----------------------------|----------------------------|----------------------------|----------------------------|-----------------------------|----------------------------|----------------------------|
| Hydrogeological unit | | | | | Alluvium | | Hawkest | bury Sandston | e (upper) | Hawkesb | ury Sandstone | e (middle) | Hawkes | 1 | | |
| Sample date | | Typical | | 25/08/2014 | 12/01/2015 | 8/04/2015 | 25/08/2014 | 12/01/2015 | 8/04/2015 | 26/08/2014 | 12/01/2015 | 8/04/2015 | 26/08/2014 | 12/01/2015 | 8/04/2015 | 9/09/2014 |
| General parameters pH (field) | Units | LOR | 6.5 - 8.0 ^c | 5.17 | 5.16 | 5.41 | 6.54 | 6.50 | 7.31 | 7.08 | 6.54 | 7.99 | 9.87 | 9.37 | 9.86 | 7.78 |
| DH (lab) | pH units | 0.01 | 405 0.000 ⁶ | 5.58 908 | 5.65 882 | 5.54 896 | 6.92 925 | 6.86 935 | 6.86 953 | 7.47 | 7.54 1,045 | 7.44 1,059 | 9.70 1,020 | 9.23 1,000 | 9.73 956 | 7.37 129 |
| Electrical conductivity (field) Electrical conductivity (lab) | µS/cm | 1 | 125 - 2,200 ^c | 945 | 930 | 912 | 958 | 967 | 963 | 1,090 | 1,110 | 1,090 | 1,050 | 1,030 | 984 | 114 |
| Femperature | °C | 0.01 | - 80 - 110%* | 17.74 9.10 | 18.77 20.20 | 19.60 21.00 | 18.02 15.10 | 21.84 34.90 | 20.18 9.40 | 17.35 13.50 | 19.20 3.30 | 18.67 4.70 | 16.51 1.00 | 21.78 2.60 | 19.91 1.70 | 16.20 102.00 |
| Dissolved oxygen Fotal dissolved solids (field) | % sat | 0.01 | - | 509 | 574 | 584 | 601 | 608 | 620 | 666 | 678 | 689 | 663 | 650 | 621 | 84 |
| Fotal dissolved solids (lab) Fotal suspended solids | mg/L | 10 5 | - | 507 190 | 514 82 | 526 74 | 440 14 | 455 13 | 420 5 | 536 <5 | 564 7 | 528 6 | 540 24 | 586 10 | 540 <5 | 47 <5 |
| Redox | mg/L mV | 0.1 | - | 2.0 | 85.9 | 88.9 | -151.2 | -84.9 | -101.9 | -121.1 | -110.0 | -138.1 | -31.1 | -111.9 | -2.1 | 153.5 |
| Nater type ^a | - | | - | Na-Mg-Cl | Na-Mg-Cl | Na-Mg-Cl | Na-Mg- HCO ₃ -Cl | Na-Mg-Ca- CI-HCO ₃ | Na-Mg-Cl- HCO ₃ | Na-Ca- HCO ₃ | Na-Ca- HCO ₃ | Na-Ca- HCO ₃ | Na-CI- HCO ₃ | Na-HCO ₃ - Cl | Na-Cl- HCO ₃ | Na-Cl- HCO ₃ |
| _aboratory analytes | | | | | | | | | - | - | | | | | | |
| Hydroxide alkalinity as CaCO ₃ Carbonate alkalinity as CaCO ₃ | mg/L mg/L | 1 | • | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 156 | <1 88 | <1 209 | <1 <1 |
| Bicarbonate alkalinity as CaCO3 | mg/L | 1 | - | 16 | 17 | 14 | 230 | 213 | 238 | 493 | 468 | 522 | 133 | 222 | 110 | 16 |
| Fotal alkalinity as $CaCO_3$ Sulfate as $SO_4^{2^-}$ | mg/L mg/L | 1 | - | 16 3 | 17 3 | 14 2 | 230 5 | 213 4 | 238 4 | 493 <1 | 468 <1 | 522 <1 | 289 <1 | 310 <1 | 319 <1 | 16 4 |
| Chloride | mg/L | 1 | - | 262 | 245 | 258 | 159 | 161 | 170 | 73 | 70 | 77 | 152 | 120 | 126 | 23 |
| Calcium Magnesium | mg/L mg/L | 1 | - | 12 22 | 13 22 | 13 23 | 29 28 | 37 32 | 34 31 | 74 23 | 88 22 | 91 23 | 3 4 | 15 6 | 4 5 | 1 |
| Sodium | mg/L | 1 | • | 115 | 106 | 110 | 110 | 107 | 105 | 124 | 104 | 110 | 246 | 180 | 189 | 14 |
| Potassium Silica (reactive) | mg/L mg/L | 1 0.05 | - | 2 19.4 | 1 18.3 | 1 19.2 | 4 13.7 | 4 12.3 | 3 12.7 | 20 8.63 | 12 8.50 | 14 8.82 | 26 3.59 | 13 4.25 | 15 3.48 | 1 3.06 |
| Fluoride | mg/L | 0.1 | - | <0.1 | <0.1 | <0.1 | 0.2 | 0.2 | 0.2 | 0.2 | 0.2 | 0.2 | 0.6 | 0.5 | 0.6 | <0.1 |
| Bromide Fotal cyanide | mg/L mg/L | 0.01 | - 0.01 | 0.597 <0.004 | 0.764 <0.004 | 0.601 <0.004 | 0.354 <0.004 | 0.596 <0.004 | 0.409 <0.004 | 0.145 <0.004 | 0.868 <0.004 | 0.213 <0.004 | 0.299 <0.004 | 0.365 <0.004 | 0.275 <0.004 | <0.004 |
| Dissolved metals | _ | | | | | | | | | | | | | | | |
| Aluminium Antimony | mg/L mg/L | 0.01 | 0.06 | 0.10 <0.001 | 0.07 <0.001 | 0.01 <0.001 | <0.01 <0.001 | <0.01 <0.001 | <0.01 <0.001 | 0.06 <0.001 | <0.01 <0.001 | <0.01 <0.001 | 0.02 <0.001 | <0.01 <0.001 | <0.01 <0.001 | 0.04 <0.00 |
| Arsenic | mg/L | 0.001 | 0.013 (As V) | <0.001 | <0.001 | <0.001 | 0.003 | 0.003 | 0.002 | 0.031 | 0.025 | 0.018 | 0.002 | 0.004 | 0.002 | < 0.00 |
| Barium Beryllium | mg/L mg/L | 0.001 | - ID | 0.667 <0.001 | 0.691 <0.001 | 0.664 <0.001 | 0.515 <0.001 | 0.629 <0.001 | 0.584 <0.001 | 3.54 <0.001 | 3.48 <0.001 | 3.27 <0.001 | 0.795 <0.001 | 1.23 <0.001 | 0.753 <0.001 | 0.029 <0.001 |
| Cadmium | mg/L | 0.0001 | 0.00 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.000 |
| Chromium Cobalt | mg/L mg/L | 0.001 | 0.001 ID | <0.001 0.042 | <0.001 0.045 | <0.001 0.041 | <0.001 0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.001 0.004 | <0.001 0.004 | <0.001 0.004 | <0.001 <0.001 | <0.001 <0.001 | <0.001 <0.001 | <0.00 <0.00 |
| Copper | mg/L | 0.001 | 0.001 | 0.002 | 0.016 | 0.003 | <0.001 | <0.001 | <0.001 | 0.003 | <0.001 | <0.001 | 0.001 | 0.001 | 0.001 | <0.00 |
| lead Manganese | mg/L mg/L | 0.001 | 0.003 | <0.001 0.472 | <0.001 0.517 | 0.002 0.488 | <0.001 0.149 | <0.001 0.160 | <0.001 0.144 | <0.001 0.047 | <0.001 0.047 | <0.001 0.047 | <0.001 0.003 | <0.001 0.009 | <0.001 <0.001 | <0.00 |
| Mercury | mg/L | 0.0001 | 0.00 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 | 0.00 | <0.000 |
| Nolybdenum Nickel | mg/L mg/L | 0.001 | ID 0.011 | <0.001 0.016 | <0.001 0.019 | <0.001 0.016 | <0.001 0.002 | <0.001 <0.001 | <0.001 <0.001 | <0.001 0.005 | <0.001 0.004 | <0.001 0.004 | 0.006 <0.001 | 0.004 <0.001 | 0.005 <0.001 | 0.001 |
| Selenium | mg/L | 0.01 | 0.011 (total) | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Strontium Jranium | mg/L mg/L | 0.001 | - ID | 0.175 <0.001 | 0.155 <0.001 | 0.143 <0.001 | 0.359 <0.001 | 0.355 <0.001 | 0.352 <0.001 | 1.03 <0.001 | 0.957 <0.001 | 0.887 <0.001 | 0.245 <0.001 | 0.30 <0.001 | 0.198 <0.001 | 0.022 <0.00 |
| /anadium | mg/L | 0.01 | ID | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Zinc Boron | mg/L mg/L | 0.005 | 0.008 | 0.057 <0.05 | 0.062 <0.05 | 0.066 <0.05 | 0.015 <0.05 | 0.058 <0.05 | 0.018 <0.05 | 0.017 <0.05 | <0.005 <0.05 | 0.012 <0.05 | 0.026 <0.05 | 0.086 <0.05 | 0.047 <0.05 | <0.00 <0.05 |
| ron | mg/L | 0.05 | ID | 0.18 | <0.05 | 0.06 | 3.76 | 4.04 | 4.21 | 0.87 | 0.86 | 1.13 | <0.05 | <0.05 | <0.05 | 0.23 |
| Bromine Nutrients | mg/L | 0.1 | ID | 0.6 | 0.5 | 0.7 | 0.4 | 0.4 | 0.4 | 0.1 | 0.2 | 0.2 | 0.3 | 0.2 | 0.3 | <0.1 |
| Ammonia as N | mg/L | 0.01 | 0.02 ^c | 0.04 | 0.04 | 0.02 | 0.08 | 0.08 | 0.06 | 0.90 | 0.94 | 0.92 | 0.96 | 0.97 | 0.86 | 0.03 |
| Nitrite as N Nitrate as N | mg/L mg/L | 0.01 | - 0.70 | <0.01 0.14 | <0.01 0.15 | <0.01 0.16 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 0.18 |
| Fotal phosphorus | mg/L | 0.01 | 0.05° | 0.14 | 0.15 <0.01 | 0.05 <0.01 | 0.03 <0.01 | 0.11 <0.01 | 0.02 | 0.02 <0.01 | 0.07 <0.01 | <0.01 <0.01 | 0.02 <0.01 | <0.01 <0.01 | <0.01 0.01 | <0.01 |
| Reactive phosphorus Fotal organic carbon | mg/L mg/L | 0.01 | 0.02 ^c | <0.01 3 | <0.01 | <0.01 | 4 | <0.01 | <1 | <0.01 | <0.01 | 4 | 20 | 14 | 22 | <0.0 |
| Dissolved gases Methane | mg/L | 0.01 | - | <0.01 | <0.01 | <0.01 | 0.092 | 0.323 | 0.273 | 28.000 | 37.000 | 44.600 | 9.390 | 41.100 | 52.300 | 0.010 |
| Ethane | mg/L | 0.01 | - | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Ethene Propane | mg/L mg/L | 0.01 | - | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 | <0.01 |
| Propene | mg/L | 0.01 | - | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Butane Butene | mg/L mg/L | 0.01 | • | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 <0.01 | <0.01 |
| Phenolic compounds | | | 200 | | | | | | | | | | | | | |
| Phenol 2-Chlorophenol | μg/L μg/L | 1 | 320 490 | 1.2 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | 8.1 <1 | 3.1 <1 | 2.8 <1 | <1 <1 |
| 2-Methylphenol | µg/L | 1 | - | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| 3-&4-Methylphenol 2-Nitrophenol | μg/L μg/L | 2 | - ID | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 | <2 <1 |
| 2.4-Dimethylphenol | µg/L | 1 | ID 160 | <1 <1 | <1 | <1 <1 | <1 | <1 <1 | <1 | <1 | <1 <1 | <1 | <1 <1 | <1 | <1 <1 | <1 <1 |
| 2.4-Dichlorophenol 2.6-Dichlorophenol | μg/L μg/L | 1 | 160 ID | <1 | <1 <1 | <1 | <1 <1 | <1 | <1 <1 | <1 <1 | <1 | <1 <1 | <1 | <1 <1 | <1 | <1 |
| 1-Chloro-3-Methylphenol | µg/L | 1 | - | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| 2.4.6-Trichlorophenol 2.4.5-Trichlorophenol | μg/L μg/L | 1 | 20 ID | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Pentachlorophenol | μg/L | 2 | ID | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 |
| Polycyclic aromatic hydrocarbons Acenaphthylene | µg/L | 1 | 16 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| Acenaphthene Fluorene | μg/L μg/L | 1 | - | <1 <1 | <1 <1 | 1.1 1.3 | <1 <1 | 4.4 3.7 | 1.9 1.7 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Phenanthrene | μg/L μg/L | 1 | - | <1 | <1 | 1.8 | <1 | 5.6 | 4.7 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| Anthracene Fluoranthene | μg/L μg/L | 1 | ID ID | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Pyrene | μg/L | 1 | ID | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| Benz(a)anthracene Chrysene | μg/L μg/L | 1 | - | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Benzo(b)fluoranthene | µg/L | 1 | - | <1 | - | - | <1 | - | - | <1 | - | - | <1 | - | - | - |
| Benzo(k)fluoranthene Benzo(a)pyrene | μg/L μg/L | 1 0.5 | - | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 | <1 <0.5 |
| ndeno(1.2.3.cd)pyrene | µg/L | 1 | ID | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| Dibenz(a.h)anthracene Benzo(g.h.i)perylene | μg/L μg/L | 1 | - | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 | <1 <1 |
| Sum of PAHs | μg/L ug/L | 0.5 | - | <0.5 | <0.5 | 4.2 | <0.5 | 17 | 8.3 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 |
| Fotal petroleum hydrocarbons C ₆ -C ₉ fraction | µg/L | 20 | ID | <20 | <20 | <20 | <20 | <20 | <20 | 190 | <20 | <20 | 200 | 80 | 110 | <20 |
| C ₁₀ -C ₁₄ fraction | µg/L | 50 | ID | <50 | <50 | <50 | <50 | 450 | 70 | <50 | <50 | <50 | <50 | <50 | <50 | <50 |
| C ₁₅ -C ₂₈ fraction C ₂₉ -C ₃₆ fraction | μg/L μg/L | 100 50 | ID ID | <100 <50 | <100 <50 | <100 <50 | <100 <50 | 120 <50 | <100 <50 | <100 <50 | <100 <50 | <100 <50 | <100 <50 | <100 <50 | <100 <50 | <100 <50 |
| C ₁₀ -C ₃₆ fraction (sum) | μg/L μg/L | 50 | ID | <50 | <50 | <50 | <50 | 570 | 70 | <50 | <50 | <50 | <50 | <50 | <50 | <50 |
| Fotal recoverable hydrocarbons C ₆ -C ₁₀ fraction | µg/L | 20 | - | <20 | <20 | <20 | <20 | <20 | <20 | 200 | <20 | <20 | 210 | 90 | 110 | <20 |
| C6-C10 fraction minus BTEX | µg/L | 20 | - | <20 | <20 | <20 | <20 | <20 | <20 | 110 | <20 | <20 | 110 | <20 | 40 | <20 |
| C_{10} - C_{16} fraction C_{16} - C_{34} fraction | μg/L μg/L | 100 100 | - | <100 <100 | <100 <100 | <100 <100 | <100 <100 | 490 100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 <100 | <100 |
| C ₃₄ -C ₄₀ fraction | μg/L μg/L | 100 | - | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 |
| C ₁₀ -C ₄₀ fraction (sum) Aromatic hydrocarbons | µg/L | 100 | - | <100 | <100 | <100 | <100 | 590 | <100 | <100 | <100 | <100 | <100 | <100 | <100 | <100 |
| Aromatic nydrocarbons Benzene | µg/L | 1 | 950 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 | <1 |
| | μg/L μg/L | 2 | ID ID | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | 94 <2 | <2 <2 | <2 <2 | 97 <2 | 79 <2 | 66 <2 | <2 |
| oluene | | 4 | | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 | <2 |
| thylbenzene n&p-Xylenes | µg/L | 2 | ID | | | | | | | | | | | | | |
| thylbenzene n&p-Xylenes -Xylenes | μg/L μg/L | 2 | ID | <2 | <2 | <2 | <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | <2 <2 | |
| thylbenzene | µg/L | | | | | | | | <2 <2 <1 <1 | <2 <2 94 <1 | <2 <2 <1 <1 | <2 <2 <1 <1 | <2 <2 97 <1 | <2 <2 79 <1 | <2 <2 66 <1 | <2 <2 <1 <1 |

Appendix C

ALS laboratory reports





| CERTIFICATE OF ANALYSIS | | | | | | | | | | |
|-------------------------|---|-------------------------|---|--|--|--|--|--|--|--|
| Work Order | ES1418991 | Page | : 1 of 13 | | | | | | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney | | | | | | | |
| Contact | : MS ANDREA MADDEN | Contact | : Client Services | | | | | | | |
| Address | | Address | : 277-289 Woodpark Road Smithfield NSW Australia 2164 | | | | | | | |
| E-mail | SYDNEY NSW, AUSTRALIA 2001 : amadden@pb.com.au | E-mail | : sydney@alsglobal.com | | | | | | | |
| Telephone | : +61 02 9272 5100 | Telephone | : +61-2-8784 8555 | | | | | | | |
| Facsimile | : +61 02 9272 5101 | Facsimile | : +61-2-8784 8500 | | | | | | | |
| Project | : 2193361A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement | | | | | | | |
| Order number | : | | | | | | | | | |
| C-O-C number | : | Date Samples Received | : 26-AUG-2014 | | | | | | | |
| Sampler | : CR | Issue Date | : 02-SEP-2014 | | | | | | | |
| Site | : | | | | | | | | | |
| | | No. of samples received | : 8 | | | | | | | |
| Quote number | : EN/008/14 | No. of samples analysed | : 8 | | | | | | | |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results
- Surrogate Control Limits

Signatories NATA Accredited Laboratory 825 This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11. Accredited for compliance with NATA ISO/IEC 17025. Signatories Position Accreditation Category Inorganic Chemist Ankit Joshi Sydney Inorganics Senior Spectroscopist Celine Conceicao Sydney Inorganics WORLD RECOGNISED ACCREDITATION Dian Dao Sydney Inorganics Senior Organic Chemist Pabi Subba Sydney Organics

Address 277-289 Woodpark Road Smithfield NSW Australia 2164 PHONE +61-2-8784 8555 Facsimile +61-2-8784 8500 Environmental Division Sydney ABN 84 009 936 029 Part of the ALS Group An ALS Limited Company



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General Comments

The analytical procedures used by the Environmental Division have been developed from established internationally recognized procedures such as those published by the USEPA, APHA, AS and NEPM. In house developed procedures are employed in the absence of documented standards or by client request.

Where moisture determination has been performed, results are reported on a dry weight basis.

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

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

When sampling time information is not provided by the client, sampling dates are shown without a time component. In these instances, the time component has been assumed by the laboratory for processing purposes.

Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

Key: CAS Number = CAS registry number from database maintained by Chemical Abstracts Services. The Chemical Abstracts Service is a division of the American Chemical Society. LOR = Limit of reporting

* = This result is computed from individual analyte detections at or above the level of reporting

- Benzo(a)pyrene Toxicity Equivalent Quotient (TEQ) is the sum total of the concentration of the eight carcinogenic PAHs multiplied by their Toxicity Equivalence Factor (TEF) relative to Benzo(a)pyrene. TEF values are provided in brackets as follows: Benz(a)anthracene (0.1), Chrysene (0.01), Benzo(b+j) & Benzo(k)fluoranthene (0.1), Benzo(a)pyrene (1.0), Indeno(1.2.3.cd)pyrene (0.1), Dibenz(a.h)anthracene (1.0), Benzo(g.h.i)perylene (0.01). Less than LOR results for 'TEQ Zero' are treated as zero.
- EG020: Bromine quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- Ionic Balance out of acceptable limits due to analytes not quantified in this report.

Page: 3 of 13Work Order: ES1418991Client: PARSONS BRINCKERHOFF AUST P/LProject: 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB01 | GLMB02 | GLMB03 | MPMB01 | MPMB02 |
|---------------------------------------|-------------|--------------|----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | C | ient samplii | ng date / time | 25-AUG-2014 11:45 | 25-AUG-2014 13:30 | 25-AUG-2014 12:15 | 25-AUG-2014 16:00 | 25-AUG-2014 15:00 |
| Compound | CAS Number | LOR | Unit | ES1418991-001 | ES1418991-002 | ES1418991-003 | ES1418991-004 | ES1418991-005 |
| EA005P: pH by PC Titrator | | | | | | | | |
| pH Value | | 0.01 | pH Unit | 7.40 | 7.39 | 8.77 | 5.58 | 6.92 |
| EA010P: Conductivity by PC Titrator | | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | µS/cm | 9640 | 6810 | 5200 | 945 | 958 |
| EA015: Total Dissolved Solids | | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 5320 | 3360 | 2460 | 507 | 440 |
| EA025: Suspended Solids | | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | <5 | 36 | <5 | 190 | 14 |
| ED009: Anions | | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 6.06 | 3.98 | 2.35 | 0.597 | 0.354 |
| ED037P: Alkalinity by PC Titrator | | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | <1 | <1 | <1 |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | <1 | <1 | 184 | <1 | <1 |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 683 | 793 | 876 | 16 | 230 |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 683 | 793 | 1060 | 16 | 230 |
| ED041G: Sulfate (Turbidimetric) as SO | 94 2- by DA | | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | <1 | 5 | <1 | 3 | 5 |
| ED045G: Chloride Discrete analyser | | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 2710 | 1370 | 934 | 262 | 159 |
| ED093F: Dissolved Major Cations | | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 174 | 210 | 16 | 12 | 29 |
| Magnesium | 7439-95-4 | 1 | mg/L | 211 | 87 | 61 | 22 | 28 |
| Sodium | 7440-23-5 | 1 | mg/L | 1360 | 1040 | 957 | 115 | 110 |
| Potassium | 7440-09-7 | 1 | mg/L | 24 | 29 | 32 | 2 | 4 |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | <0.01 | <0.01 | 0.01 | 0.10 | <0.01 |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Arsenic | 7440-38-2 | 0.001 | mg/L | <0.001 | 0.014 | 0.020 | <0.001 | 0.003 |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Barium | 7440-39-3 | 0.001 | mg/L | 2.64 | 9.78 | 4.49 | 0.667 | 0.515 |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Copper | 7440-50-8 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | 0.002 | <0.001 |
| Cobalt | 7440-48-4 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | 0.042 | 0.001 |

Page : 4 of 13 Work Order : ES1418991 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB01 | GLMB02 | GLMB03 | MPMB01 | MPMB02 |
|---------------------------------------|--------------------|-------------|----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Ci | ient sampli | ng date / time | 25-AUG-2014 11:45 | 25-AUG-2014 13:30 | 25-AUG-2014 12:15 | 25-AUG-2014 16:00 | 25-AUG-2014 15:00 |
| Compound | CAS Number | LOR | Unit | ES1418991-001 | ES1418991-002 | ES1418991-003 | ES1418991-004 | ES1418991-005 |
| EG020F: Dissolved Metals by ICP-MS - | Continued | | | | | | | |
| Nickel | 7440-02-0 | 0.001 | mg/L | 0.002 | 0.002 | <0.001 | 0.016 | 0.002 |
| Lead | 7439-92-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Zinc | 7440-66-6 | 0.005 | mg/L | <0.005 | 0.020 | 0.006 | 0.057 | 0.015 |
| Manganese | 7439-96-5 | 0.001 | mg/L | 0.300 | 0.666 | 0.002 | 0.472 | 0.149 |
| Molybdenum | 7439-98-7 | 0.001 | mg/L | <0.001 | <0.001 | 0.005 | <0.001 | <0.001 |
| Selenium | 7782-49-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Strontium | 7440-24-6 | 0.001 | mg/L | 6.33 | 5.14 | 2.06 | 0.175 | 0.359 |
| Uranium | 7440-61-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Vanadium | 7440-62-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Boron | 7440-42-8 | 0.05 | mg/L | <0.05 | <0.05 | <0.05 | <0.05 | <0.05 |
| Iron | 7439-89-6 | 0.05 | mg/L | 0.19 | 2.97 | <0.05 | 0.18 | 3.76 |
| Bromine | 7726-95-6 | 0.1 | mg/L | 7.2 | 4.5 | 2.7 | 0.6 | 0.4 |
| G035F: Dissolved Mercury by FIMS | | | | | | | | |
| Mercury | 7439-97-6 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 |
| G052G: Silica by Discrete Analyser | | | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 10.6 | 21.2 | 9.77 | 19.4 | 13.7 |
| K026SF: Total CN by Segmented Flo | w Analyser | | | | | | | |
| Total Cyanide | 57-12-5 | 0.004 | mg/L | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 |
| K040P: Fluoride by PC Titrator | | | | | | | | |
| Fluoride | 16984-48-8 | 0.1 | mg/L | 0.3 | 0.3 | 0.2 | <0.1 | 0.2 |
| K055G: Ammonia as N by Discrete A | nalyser | | | | | | | |
| Ammonia as N | 7664-41-7 | 0.01 | mg/L | 1.71 | 1.88 | 2.37 | 0.04 | 0.08 |
| K057G: Nitrite as N by Discrete Anal | vser | | | | | | | |
| Nitrite as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| EK058G: Nitrate as N by Discrete Ana | lvser | | | | | | | |
| Nitrate as N | 14797-55-8 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | 0.14 | <0.01 |
| K059G: Nitrite plus Nitrate as N (NO) | () by Discrete Ana | lvser | | | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | 0.14 | <0.01 |
| K067G: Total Phosphorus as P by Dis | screte Analyser | | | | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | 0.08 | 0.12 | 0.04 | 0.14 | 0.03 |
| K071G: Reactive Phosphorus as P by | / discrete analyse | | | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | 0.06 | 0.06 | <0.01 | <0.01 | <0.01 |
| N055: Ionic Balance | | | 3 | | | | | |

Page : 5 of 13 Work Order : ES1418991 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | GLMB01 | GLMB02 | GLMB03 | MPMB01 | MPMB02 |
|--------------------------------------|-------------|-------------|----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Cl | ient sampli | ng date / time | 25-AUG-2014 11:45 | 25-AUG-2014 13:30 | 25-AUG-2014 12:15 | 25-AUG-2014 16:00 | 25-AUG-2014 15:00 |
| Compound | CAS Number | LOR | Unit | ES1418991-001 | ES1418991-002 | ES1418991-003 | ES1418991-004 | ES1418991-005 |
| EN055: Ionic Balance - Continued | | | | | | | | |
| Total Anions | | 0.01 | meq/L | 90.1 | 54.6 | 47.5 | 7.77 | 9.18 |
| Total Cations | | 0.01 | meq/L | 85.8 | 63.6 | 48.3 | 7.46 | 8.64 |
| Ionic Balance | | 0.01 | % | 2.44 | 7.62 | 0.73 | 2.04 | 3.08 |
| EP005: Total Organic Carbon (TOC) | | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | 3 | 9 | 10 | 3 | 4 |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 29900 | 21700 | 18800 | <10 | 92 |
| Ethene | 74-85-1 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Ethane | 74-84-0 | 10 | µg/L | 1480 | 931 | 118 | <10 | <10 |
| Propene | 115-07-1 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Propane | 74-98-6 | 10 | µg/L | 419 | 194 | 28 | <10 | <10 |
| Butene | 25167-67-3 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Butane | 106-97-8 | 10 | µg/L | 72 | 44 | <10 | <10 | <10 |
| EP075(SIM)A: Phenolic Compounds | | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | 2.0 | 1.8 | 1.5 | 1.2 | <1.0 |
| 2-Chlorophenol | 95-57-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | <2.0 |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | <2.0 |
| EP075(SIM)B: Polynuclear Aromatic Hy | ydrocarbons | | | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |

Page : 6 of 13 Work Order : ES1418991 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB01 | GLMB02 | GLMB03 | MPMB01 | MPMB02 |
|---|--------------------|-------------|----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Cl | ient sampli | ng date / time | 25-AUG-2014 11:45 | 25-AUG-2014 13:30 | 25-AUG-2014 12:15 | 25-AUG-2014 16:00 | 25-AUG-2014 15:00 |
| Compound | CAS Number | LOR | Unit | ES1418991-001 | ES1418991-002 | ES1418991-003 | ES1418991-004 | ES1418991-005 |
| EP075(SIM)B: Polynuclear Aromatic H | ydrocarbons - Cont | tinued | | | | | | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benz(a)anthracene | 56-55-3 | 1.0 | μg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(b+j)fluoranthene | 205-99-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Sum of polycyclic aromatic hydrocarbons | ; | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 |
| [^] Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 |
| EP080/071: Total Petroleum Hydrocart | bons | | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | 210 | 210 | 310 | <20 | <20 |
| C10 - C14 Fraction | | 50 | µg/L | <50 | <50 | <50 | <50 | <50 |
| C15 - C28 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| C29 - C36 Fraction | | 50 | µg/L | <50 | <50 | <50 | <50 | <50 |
| [^] C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | <50 | <50 | <50 | <50 |
| EP080/071: Total Recoverable Hydroca | arbons - NEPM 201 | 3 | | | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | 290 | 220 | 310 | <20 | <20 |
| [^] C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | 290 | 210 | 170 | <20 | <20 |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| ^ >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| ^ >C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| EP080: BTEXN | | | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | <1 | <1 | <1 |
| Toluene | 108-88-3 | 2 | μg/L | <2 | 10 | 139 | <2 | <2 |
| Ethylbenzene | 100-41-4 | 2 | μg/L | <2 | <2 | <2 | <2 | <2 |
| meta- & para-Xylene | 108-38-3 106-42-3 | 2 | μg/L | <2 | <2 | <2 | <2 | <2 |
| ortho-Xylene | 95-47-6 | 2 | µg/L | <2 | <2 | <2 | <2 | <2 |

Page : 7 of 13 Work Order : ES1418991 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | GLMB01 | GLMB02 | GLMB03 | MPMB01 | MPMB02 |
|-----------------------------------|------------|------------|-----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Cli | ient sampl | ing date / time | 25-AUG-2014 11:45 | 25-AUG-2014 13:30 | 25-AUG-2014 12:15 | 25-AUG-2014 16:00 | 25-AUG-2014 15:00 |
| Compound | CAS Number | LOR | Unit | ES1418991-001 | ES1418991-002 | ES1418991-003 | ES1418991-004 | ES1418991-005 |
| EP080: BTEXN - Continued | | | | | | | | |
| ∑ Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | <2 | <2 | <2 |
| Sum of BTEX | | 1 | µg/L | <1 | 10 | 139 | <1 | <1 |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | <5 | <5 | <5 | <5 |
| EP075(SIM)S: Phenolic Compound Su | urrogates | | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 25.3 | 31.3 | 22.4 | 29.5 | 27.2 |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 48.5 | 61.2 | 46.3 | 60.6 | 56.3 |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 58.8 | 81.2 | 54.7 | 58.8 | 56.4 |
| EP075(SIM)T: PAH Surrogates | | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 58.2 | 72.5 | 58.8 | 81.6 | 75.2 |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 91.6 | 95.2 | 67.8 | 91.4 | 82.9 |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 76.4 | 96.6 | 66.5 | 90.9 | 80.3 |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 111 | 105 | 102 | 106 | 108 |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 98.8 | 97.9 | 95.6 | 95.6 | 94.9 |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 102 | 97.3 | 96.1 | 97.2 | 95.7 |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB03 | MPMB04 | QA | |
|--|-------------|---------------|----------------|-------------------|-------------------|-------------------|------|
| | C | lient samplii | ng date / time | 26-AUG-2014 11:30 | 26-AUG-2014 12:40 | 25-AUG-2014 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1418991-006 | ES1418991-007 | ES1418991-008 | |
| EA005P: pH by PC Titrator | | | | | | | |
| pH Value | | 0.01 | pH Unit | 7.47 | 9.70 | 7.45 | |
| EA010P: Conductivity by PC Titrator | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | μS/cm | 1090 | 1050 | 9650 | |
| EA015: Total Dissolved Solids | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 536 | 540 | 4540 | |
| EA025: Suspended Solids | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | <5 | 24 | <5 | |
| ED009: Anions | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 0.145 | 0.299 | 7.11 | |
| ED037P: Alkalinity by PC Titrator | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | <1 | |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | <1 | 156 | <1 | |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 493 | 133 | 679 | |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 493 | 289 | 679 | |
| ED041G: Sulfate (Turbidimetric) as SO4 | 4 2- by DA | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | <1 | <1 | <1 | |
| ED045G: Chloride Discrete analyser | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 73 | 152 | 2730 | |
| ED093F: Dissolved Major Cations | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 74 | 3 | 185 | |
| Magnesium | 7439-95-4 | 1 | mg/L | 23 | 4 | 220 | |
| Sodium | 7440-23-5 | 1 | mg/L | 124 | 246 | 1460 | |
| Potassium | 7440-09-7 | 1 | mg/L | 20 | 26 | 27 | |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | 0.06 | 0.02 | 0.01 | |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Arsenic | 7440-38-2 | 0.001 | mg/L | 0.031 | 0.002 | <0.001 | |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Barium | 7440-39-3 | 0.001 | mg/L | 3.54 | 0.795 | 2.67 | |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Copper | 7440-50-8 | 0.001 | mg/L | 0.003 | 0.001 | <0.001 | |
| Cobalt | 7440-48-4 | 0.001 | mg/L | 0.004 | <0.001 | <0.001 | |

Page : 9 of 13 Work Order : ES1418991 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | nt sample ID | MPMB03 | MPMB04 | QA | |
|--|---------|-------------|----------------|-------------------|-------------------|-------------------|------|
| | Clie | ent samplir | ng date / time | 26-AUG-2014 11:30 | 26-AUG-2014 12:40 | 25-AUG-2014 15:00 | |
| Compound CAS I | Number | LOR | Unit | ES1418991-006 | ES1418991-007 | ES1418991-008 | |
| EG020F: Dissolved Metals by ICP-MS - Continued | | | | | | | |
| Nickel 74 | 40-02-0 | 0.001 | mg/L | 0.005 | <0.001 | 0.004 | |
| Lead 74 | 39-92-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Zinc 74 | 40-66-6 | 0.005 | mg/L | 0.017 | 0.026 | 0.009 | |
| Manganese 74 | 39-96-5 | 0.001 | mg/L | 0.047 | 0.003 | 0.298 | |
| Molybdenum 74 | 39-98-7 | 0.001 | mg/L | <0.001 | 0.006 | <0.001 | |
| Selenium 77 | 82-49-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| Strontium 74 | 40-24-6 | 0.001 | mg/L | 1.03 | 0.245 | 6.43 | |
| Uranium 74 | 40-61-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Vanadium 74 | 40-62-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| | 40-42-8 | 0.05 | mg/L | <0.05 | <0.05 | <0.05 | |
| Iron 74 | 39-89-6 | 0.05 | mg/L | 0.87 | <0.05 | 0.20 | |
| | 26-95-6 | 0.1 | mg/L | 0.1 | 0.3 | 7.4 | |
| EG035F: Dissolved Mercury by FIMS | | | | | | | |
| | 39-97-6 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | |
| EG052G: Silica by Discrete Analyser | | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 8.63 | 3.59 | 10.3 | |
| EK026SF: Total CN by Segmented Flow Analyse | r | | | | | | |
| | 57-12-5 | 0.004 | mg/L | <0.004 | <0.004 | <0.004 | |
| EK040P: Fluoride by PC Titrator | | | | | | | |
| | 84-48-8 | 0.1 | mg/L | 0.2 | 0.6 | 0.3 | |
| EK055G: Ammonia as N by Discrete Analyser | | | | | | | |
| | 64-41-7 | 0.01 | mg/L | 0.90 | 0.96 | 1.66 | |
| EK057G: Nitrite as N by Discrete Analyser | | | | | | | |
| Nitrite as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| EK058G: Nitrate as N by Discrete Analyser | | | - | | | | |
| | 97-55-8 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| EK059G: Nitrite plus Nitrate as N (NOx) by Discr | | vser | | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| EK067G: Total Phosphorus as P by Discrete Ana | vser | | - | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | 0.02 | 0.02 | 0.13 | |
| EK071G: Reactive Phosphorus as P by discrete a | nalvsor | | - | | | | |
| | 65-44-2 | 0.01 | mg/L | <0.01 | <0.01 | 0.06 | |
| EN055: Ionic Balance | | | , | | | | |
| | | | | | | | |

Page: 10 of 13Work Order: ES1418991Client: PARSONS BRINCKERHOFF AUST P/LProject: 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB03 | MPMB04 | QA | |
|---------------------------------------|------------|-------------|----------------|-------------------|-------------------|-------------------|------|
| | Cl | ient sampli | ng date / time | 26-AUG-2014 11:30 | 26-AUG-2014 12:40 | 25-AUG-2014 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1418991-006 | ES1418991-007 | ES1418991-008 | |
| EN055: Ionic Balance - Continued | | | | | | | |
| Total Anions | | 0.01 | meq/L | 11.9 | 10.1 | 90.6 | |
| Total Cations | | 0.01 | meq/L | 11.5 | 11.8 | 91.5 | |
| Ionic Balance | | 0.01 | % | 1.80 | 8.10 | 0.51 | |
| EP005: Total Organic Carbon (TOC) | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | <1 | 20 | 4 | |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 28000 | 9390 | 28800 | |
| Ethene | 74-85-1 | 10 | µg/L | <10 | <10 | <10 | |
| Ethane | 74-84-0 | 10 | µg/L | <10 | <10 | 1480 | |
| Propene | 115-07-1 | 10 | μg/L | <10 | <10 | <10 | |
| Propane | 74-98-6 | 10 | µg/L | <10 | <10 | 402 | |
| Butene | 25167-67-3 | 10 | μg/L | <10 | <10 | <10 | |
| Butane | 106-97-8 | 10 | µg/L | <10 | <10 | 90 | |
| EP075(SIM)A: Phenolic Compounds | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | <1.0 | 8.1 | <1.0 | |
| 2-Chlorophenol | 95-57-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | |
| EP075(SIM)B: Polynuclear Aromatic Hyd | rocarbons | | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |

Page: 11 of 13Work Order: ES1418991Client: PARSONS BRINCKERHOFF AUST P/LProject: 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB03 | MPMB04 | QA | |
|--|---------------------|------------|----------------|-------------------|-------------------|-------------------|------|
| | Clie | ent sampli | ng date / time | 26-AUG-2014 11:30 | 26-AUG-2014 12:40 | 25-AUG-2014 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1418991-006 | ES1418991-007 | ES1418991-008 | |
| EP075(SIM)B: Polynuclear Aromatic Hy | ydrocarbons - Conti | inued | | | | | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(b+j)fluoranthene | 205-99-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| [^] Sum of polycyclic aromatic hydrocarbons | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | |
| [^] Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | |
| EP080/071: Total Petroleum Hydrocarb | oons | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | 190 | 200 | 240 | |
| C10 - C14 Fraction | | 50 | µg/L | <50 | <50 | <50 | |
| C15 - C28 Fraction | | 100 | µg/L | <100 | <100 | <100 | |
| C29 - C36 Fraction | | 50 | µg/L | <50 | <50 | <50 | |
| [^] C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | <50 | <50 | |
| EP080/071: Total Recoverable Hydroca | rbons - NEPM 201 | 3 | | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | 200 | 210 | 290 | |
| C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | 110 | 110 | 290 | |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | <100 | <100 | |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | <100 | <100 | |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | <100 | |
| ^ >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | <100 | <100 | |
| ^ >C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | <100 | <100 | |
| | | | | | | | |
| EP080: BTEXN Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | <1 | |
| Toluene | 108-88-3 | 2 | μg/L | 94 | 97 | <2 | |
| Ethylbenzene | 100-80-3 | 2 | μg/L | <2 | <2 | <2 | |
| meta- & para-Xylene | 108-38-3 106-42-3 | 2 | μg/L | <2 | <2 | <2 | |
| ortho-Xylene | 95-47-6 | 2 | μg/L | <2 | <2 | <2 | |
| | 90-47-0 | 4 | ₽9′⊏ | 2۲ | ~ <u>~</u> | ~2 | |

Page: 12 of 13Work Order: ES1418991Client: PARSONS BRINCKERHOFF AUST P/LProject: 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | MPMB03 | MPMB04 | QA | |
|-----------------------------------|------------|------------|-----------------|-------------------|-------------------|-------------------|------|
| | Cli | ent sampli | ing date / time | 26-AUG-2014 11:30 | 26-AUG-2014 12:40 | 25-AUG-2014 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1418991-006 | ES1418991-007 | ES1418991-008 | |
| EP080: BTEXN - Continued | | | | | | | |
| ^ Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | <2 | |
| [^] Sum of BTEX | | 1 | µg/L | 94 | 97 | <1 | |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | <5 | <5 | |
| EP075(SIM)S: Phenolic Compound Su | irrogates | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 27.6 | 27.8 | 25.9 | |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 58.6 | 55.0 | 52.4 | |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 51.2 | 64.2 | 61.8 | |
| EP075(SIM)T: PAH Surrogates | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 74.4 | 73.7 | 65.7 | |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 86.6 | 83.5 | 100 | |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 81.5 | 82.1 | 80.6 | |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 105 | 107 | 104 | |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 93.7 | 95.8 | 94.4 | |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 94.9 | 97.3 | 95.3 | |



Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) |
|---|------------|----------|------------|
| Compound | CAS Number | Low | High |
| EP075(SIM)S: Phenolic Compound Surrogates | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 |
| EP075(SIM)T: PAH Surrogates | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 |
| EP080S: TPH(V)/BTEX Surrogates | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 |
| Toluene-D8 | 2037-26-5 | 79 | 131 |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 |



| CERTIFICATE OF ANALYSIS | | | | | | |
|-------------------------|---------------------------------|-------------------------|---|--|--|--|
| Work Order | ES1419407 | Page | : 1 of 8 | | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney | | | |
| Contact | : MS ANDREA MADDEN | Contact | : Client Services | | | |
| Address | : GPO BOX 5394 | Address | : 277-289 Woodpark Road Smithfield NSW Australia 2164 | | | |
| | SYDNEY NSW, AUSTRALIA 2001 | | | | | |
| E-mail | : amadden@pb.com.au | E-mail | : sydney@alsglobal.com | | | |
| Telephone | : +61 02 9272 5100 | Telephone | : +61-2-8784 8555 | | | |
| acsimile | : +61 02 9272 5101 | Facsimile | : +61-2-8784 8500 | | | |
| Project | : 2193361A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement | | | |
| Order number | : | | | | | |
| C-O-C number | : | Date Samples Received | : 29-AUG-2014 | | | |
| Sampler | : CR | Issue Date | : 04-SEP-2014 | | | |
| Site | : | | | | | |
| | | No. of samples received | : 2 | | | |
| Quote number | : SY/743/14 V2 | No. of samples analysed | : 2 | | | |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results
- Surrogate Control Limits

Signatories NATA Accredited Laboratory 825 This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11. Accredited for compliance with NATA ISO/IEC 17025. Signatories Position Accreditation Category Inorganic Chemist Ankit Joshi Sydney Inorganics Senior Spectroscopist Celine Conceicao Sydney Inorganics WORLD RECOGNISED ACCREDITATION Dian Dao Sydney Inorganics Laboratory Manager - Organics Phalak Inthakesone Sydney Organics Metals Coordinator Shobhna Chandra Sydney Inorganics

Address 277-289 Woodpark Road Smithfield NSW Australia 2164 PHONE +61-2-8784 8555 Facsimile +61-2-8784 8500 Environmental Division Svdnev ABN 84 009 936 029 Part of the ALS Group An ALS Limited Company



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General Comments

The analytical procedures used by the Environmental Division have been developed from established internationally recognized procedures such as those published by the USEPA, APHA, AS and NEPM. In house developed procedures are employed in the absence of documented standards or by client request.

Where moisture determination has been performed, results are reported on a dry weight basis.

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

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

When sampling time information is not provided by the client, sampling dates are shown without a time component. In these instances, the time component has been assumed by the laboratory for processing purposes.

Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

Key : CAS Number = CAS registry number from database maintained by Chemical Abstracts Services. The Chemical Abstracts Service is a division of the American Chemical Society. LOR = Limit of reporting ^ = This result is computed from individual analyte detections at or above the level of reporting

- Benzo(a)pyrene Toxicity Equivalent Quotient (TEQ) is the sum total of the concentration of the eight carcinogenic PAHs multiplied by their Toxicity Equivalence Factor (TEF) relative to Benzo(a)pyrene. TEF values are provided in brackets as follows: Benz(a)anthracene (0.1), Chrysene (0.01), Benzo(b+j) & Benzo(k)fluoranthene (0.1), Benzo(a)pyrene (1.0), Indeno(1.2.3.cd)pyrene (0.1), Dibenz(a,h)anthracene (1.0), Benzo(g,h,i)pervlene (0.01), Less than LOR results for 'TEQ Zero' are treated as zero.
- ED041G: LOR raised for Sulfate analysis on sample ID:RMB02 due to matrix interferences.
- EG020: Bromine quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- Ionic Balance out of acceptable limits due to analytes not quantified in this report.
- It has been noted that Nitrite is greater than NOx on sample ID (RMB02,RMB03), however this difference is within the limits of experimental variation.



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | RMB02 | RMB03 | | |
|--|-------------|---------------|----------------|---------------|---------------|------|--|
| | Ci | lient samplir | ng date / time | [29-AUG-2014] | [29-AUG-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1419407-001 | ES1419407-002 | | |
| EA005P: pH by PC Titrator | | | | | | | |
| pH Value | | 0.01 | pH Unit | 7.18 | 9.48 | | |
| EA010P: Conductivity by PC Titrator | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | μS/cm | 10600 | 7940 | | |
| EA015: Total Dissolved Solids | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 5440 | 3830 | | |
| EA025: Suspended Solids | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | <5 | | | |
| Suspended Solids (SS) | | 5 | mg/L | | <5 | | |
| ED009: Anions | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 5.88 | 4.71 | | |
| ED037P: Alkalinity by PC Titrator | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | | |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | <1 | 298 | | |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 972 | 239 | | |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 972 | 537 | | |
| ED041G: Sulfate (Turbidimetric) as SO4 | 2- by DA | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | <10 | <1 | | |
| ED045G: Chloride Discrete analyser | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 3120 | 1760 | | |
| ED093F: Dissolved Major Cations | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 270 | 6 | | |
| Magnesium | 7439-95-4 | 1 | mg/L | 81 | 14 | | |
| Sodium | 7440-23-5 | 1 | mg/L | 1740 | 1610 | | |
| Potassium | 7440-09-7 | 1 | mg/L | 37 | 22 | | |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | <0.01 | <0.01 | | |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Arsenic | 7440-38-2 | 0.001 | mg/L | <0.001 | 0.001 | | |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Barium | 7440-39-3 | 0.001 | mg/L | 40.0 | 4.35 | | |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | | |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Cobalt | 7440-48-4 | 0.001 | mg/L | <0.001 | <0.001 | | |

Page : 4 of 8 Work Order : ES1419407 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | RMB02 | RMB03 | | |
|--|------------------------|--------------|----------------|---------------|---------------|------|--|
| | Cl | ient samplii | ng date / time | [29-AUG-2014] | [29-AUG-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1419407-001 | ES1419407-002 | | |
| EG020F: Dissolved Metals by ICP-MS - Cont | tinued | | | | | | |
| Copper | 7440-50-8 | 0.001 | mg/L | <0.001 | 0.002 | | |
| Lead | 7439-92-1 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Manganese | 7439-96-5 | 0.001 | mg/L | 0.026 | <0.001 | | |
| Molybdenum | 7439-98-7 | 0.001 | mg/L | <0.001 | 0.004 | | |
| Nickel | 7440-02-0 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Selenium | 7782-49-2 | 0.01 | mg/L | <0.01 | <0.01 | | |
| Strontium | 7440-24-6 | 0.001 | mg/L | 8.18 | 1.64 | | |
| Uranium | 7440-61-1 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Vanadium | 7440-62-2 | 0.01 | mg/L | <0.01 | <0.01 | | |
| Zinc | 7440-66-6 | 0.005 | mg/L | 0.017 | 0.053 | | |
| Boron | 7440-42-8 | 0.05 | mg/L | <0.05 | 0.15 | | |
| Iron | 7439-89-6 | 0.05 | mg/L | 4.81 | <0.05 | | |
| Bromine | 7726-95-6 | 0.1 | mg/L | 7.9 | 5.8 | | |
| EG035F: Dissolved Mercury by FIMS | | | | | | | |
| Mercury | 7439-97-6 | 0.0001 | mg/L | <0.0001 | <0.0001 | | |
| EG052G: Silica by Discrete Analyser | | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 11.6 | 7.30 | | |
| EK026SF: Total CN by Segmented Flow Ar | nalvser | | | | | | |
| Total Cyanide | 57-12-5 | 0.004 | mg/L | <0.004 | <0.004 | | |
| EK040P: Fluoride by PC Titrator | | | | | | | |
| Fluoride | 16984-48-8 | 0.1 | mg/L | 0.2 | 0.4 | | |
| EK055G: Ammonia as N by Discrete Analys | ser | | | | | | |
| Ammonia as N | 7664-41-7 | 0.01 | mg/L | 4.30 | 3.73 | | |
| EK057G: Nitrite as N by Discrete Analyser | | | | | | | |
| Nitrite as N | | 0.01 | mg/L | 0.04 | 0.04 | | |
| EK058G: Nitrate as N by Discrete Analyse | r | | | | | | |
| Nitrate as N | 14797-55-8 | 0.01 | mg/L | <0.01 | <0.01 | | |
| EK059G: Nitrite plus Nitrate as N (NOx) by | y Discr <u>ete Ana</u> | lyser | | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | 0.03 | 0.02 | | |
| EK067G: Total Phosphorus as P by Discret | te Anal <u>yser</u> | | | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | <0.01 | <0.01 | | |
| EK071G: Reactive Phosphorus as P by disc | crete analyser | | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | <0.01 | <0.01 | | |
| | | | - | | | | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | RMB02 | RMB03 | | |
|---------------------------------------|------------|-------------|----------------|---------------|---------------|------|--|
| | CI | ient sampli | ng date / time | [29-AUG-2014] | [29-AUG-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1419407-001 | ES1419407-002 | | |
| EN055: Ionic Balance | | | | | | | |
| Total Anions | | 0.01 | meq/L | 107 | 60.4 | | |
| Total Cations | | 0.01 | meq/L | 96.8 | 72.0 | | |
| Ionic Balance | | 0.01 | % | 5.23 | 8.78 | | |
| EP005: Total Organic Carbon (TOC) | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | <1 | 28 | | |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 33100 | 40900 | | |
| Ethene | 74-85-1 | 10 | µg/L | <10 | <10 | | |
| Ethane | 74-84-0 | 10 | µg/L | 11 | <10 | | |
| Propene | 115-07-1 | 10 | µg/L | <10 | <10 | | |
| Propane | 74-98-6 | 10 | µg/L | <10 | <10 | | |
| Butene | 25167-67-3 | 10 | μg/L | <10 | <10 | | |
| Butane | 106-97-8 | 10 | µg/L | <10 | <10 | | |
| EP075(SIM)A: Phenolic Compounds | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2-Chlorophenol | 95-57-8 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | | |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | | |
| EP075(SIM)B: Polynuclear Aromatic Hyd | drocarbons | | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Anthracene | 120-12-7 | 1.0 | μg/L | <1.0 | <1.0 | | |

Page : 6 of 8 Work Order : ES1419407 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | RMB02 | RMB03 | | |
|--|-------------------|-------------|----------------|---------------|---------------|------|--|
| | Cli | ient sampli | ng date / time | [29-AUG-2014] | [29-AUG-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1419407-001 | ES1419407-002 | | |
| EP075(SIM)B: Polynuclear Aromatic Hy | drocarbons - Cont | inued | | | | | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(b+j)fluoranthene | 205-99-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | | |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| [^] Sum of polycyclic aromatic hydrocarbons | | 0.5 | µg/L | <0.5 | <0.5 | | |
| A Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | | |
| EP080/071: Total Petroleum Hydrocarb | ons | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | <20 | 70 | | |
| C10 - C14 Fraction | | 50 | µg/L | <50 | 90 | | |
| C15 - C28 Fraction | | 100 | µg/L | <100 | <100 | | |
| C29 - C36 Fraction | | 50 | µg/L | <50 | <50 | | |
| ^ C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | 90 | | |
| EP080/071: Total Recoverable Hydroca | rbons - NEPM 201 | 3 | | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | <20 | 70 | | |
| C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | <20 | 20 | | |
| EP080/071: Total Recoverable Hydroca | rbons - NEPM 201 | 3 Fractio | าร | | | | |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | <100 | | |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | <100 | | |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | | |
| ^ >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | <100 | | |
| ^ >C10 - C16 Fraction minus Naphthalene | | 100 | µg/L | <100 | <100 | | |
| (F2) | | | | | | | |
| EP080: BTEXN | | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | | |
| Toluene | 108-88-3 | 2 | µg/L | 8 | 48 | | |
| Ethylbenzene | 100-41-4 | 2 | µg/L | <2 | <2 | | |
| meta- & para-Xylene | 108-38-3 106-42-3 | 2 | µg/L | <2 | <2 | | |

Page: 7 of 8Work Order: ES1419407Client: PARSONS BRINCKERHOFF AUST P/LProject: 2193361A



| | | | | | | | |
|-----------------------------------|-----------------------------|------------------|------|---------------|---------------|---|------|
| Sub-Matrix: WATER (Matrix: WATER) | | Client sample ID | | RMB02 | RMB03 | | |
| | Client sampling date / time | | | [29-AUG-2014] | [29-AUG-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1419407-001 | ES1419407-002 | | |
| EP080: BTEXN - Continued | | | | | | | |
| ortho-Xylene | 95-47-6 | 2 | µg/L | <2 | <2 | | |
| Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | | |
| Sum of BTEX | | 1 | µg/L | 8 | 48 | | |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | <5 | | |
| EP075(SIM)S: Phenolic Compound Su | irrogates | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 25.6 | 29.7 | | |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 53.3 | 62.7 | | |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 46.8 | 80.8 | | |
| EP075(SIM)T: PAH Surrogates | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 62.0 | 74.7 | | |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 74.8 | 81.4 | | |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 75.7 | 85.5 | | |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 100 | 96.9 | | |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 105 | 102 | | |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 86.9 | 81.7 | | |
| | | | | | | 1 | |



Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) |
|---|------------|----------|------------|
| Compound | CAS Number | Low | High |
| EP075(SIM)S: Phenolic Compound Surrogates | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 |
| EP075(SIM)T: PAH Surrogates | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 |
| EP080S: TPH(V)/BTEX Surrogates | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 |
| Toluene-D8 | 2037-26-5 | 79 | 131 |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 |



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|----------------|---------------------------------|-------------------------|---|--|--|--|--|--|--|--|--|
| Work Order | ES1420206 | Page | : 1 of 8 | | | | | | | | |
| Amendment | : 1 | | | | | | | | | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney | | | | | | | | |
| Contact | : MS ANDREA MADDEN | Contact | : Client Services | | | | | | | | |
| Address | : GPO BOX 5394 | Address | : 277-289 Woodpark Road Smithfield NSW Australia 2164 | | | | | | | | |
| | SYDNEY NSW, AUSTRALIA 2001 | | | | | | | | | | |
| E-mail | : amadden@pb.com.au | E-mail | : sydney@alsglobal.com | | | | | | | | |
| Telephone | : +61 02 9272 5100 | Telephone | : +61-2-8784 8555 | | | | | | | | |
| Facsimile | : +61 02 9272 5101 | Facsimile | : +61-2-8784 8500 | | | | | | | | |
| Project | : 2193361A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement | | | | | | | | |
| Order number | : | | | | | | | | | | |
| C-O-C number | : | Date Samples Received | : 09-SEP-2014 | | | | | | | | |
| Sampler | : CR | Issue Date | : 03-AUG-2015 | | | | | | | | |
| Site | : | | | | | | | | | | |
| | | No. of samples received | : 1 | | | | | | | | |
| Quote number | : SY/743/14 V2 | No. of samples analysed | : 1 | | | | | | | | |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

Accredited for compliance with

ISO/IEC 17025.

- General Comments
- Analytical Results
- Surrogate Control Limits



This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11.

| Signatories | Position | Accreditation Category |
|-----------------|------------------------|------------------------|
| Ankit Joshi | Inorganic Chemist | Sydney Inorganics |
| Dian Dao | Inorganic Chemist | Sydney Inorganics |
| Pabi Subba | Senior Organic Chemist | Sydney Organics |
| Shobhna Chandra | Metals Coordinator | Sydney Inorganics |

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General Comments

The analytical procedures used by the Environmental Division have been developed from established internationally recognized procedures such as those published by the USEPA, APHA, AS and NEPM. In house developed procedures are employed in the absence of documented standards or by client request.

Where moisture determination has been performed, results are reported on a dry weight basis.

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

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

When sampling time information is not provided by the client, sampling dates are shown without a time component. In these instances, the time component has been assumed by the laboratory for processing purposes.

Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

Key: CAS Number = CAS registry number from database maintained by Chemical Abstracts Services. The Chemical Abstracts Service is a division of the American Chemical Society. LOR = Limit of reporting

* = This result is computed from individual analyte detections at or above the level of reporting

- Benzo(a)pyrene Toxicity Equivalent Quotient (TEQ) is the sum total of the concentration of the eight carcinogenic PAHs multiplied by their Toxicity Equivalence Factor (TEF) relative to Benzo(a)pyrene. TEF values are provided in brackets as follows: Benz(a)anthracene (0.1), Chrysene (0.01), Benzo(b+j) & Benzo(k)fluoranthene (0.1), Benzo(a)pyrene (1.0), Indeno(1.2.3.cd)pyrene (0.1), Dibenz(a.h)anthracene (1.0), Benzo(g.h.i)perylene (0.01). Less than LOR results for 'TEQ Zero' are treated as zero.
- EG020: Bromine quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- This report has been amended as a result of a request to change sample identification numbers (IDs) received by ALS from Carolina Sardella on the 3rd August 2015. All analysis results are as per the previous report.



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | NR | | |
|--|-------------|---------------|----------------|---------------|------|------|
| | Ci | lient samplii | ng date / time | [09-SEP-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1420206-001 | | |
| EA005P: pH by PC Titrator | | | | | | |
| pH Value | | 0.01 | pH Unit | 7.37 | | |
| EA010P: Conductivity by PC Titrator | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | µS/cm | 114 | | |
| EA015: Total Dissolved Solids | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 47 | | |
| EA025: Suspended Solids | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | <5 | | |
| ED009: Anions | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 0.024 | | |
| ED037P: Alkalinity by PC Titrator | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | | |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | <1 | | |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 16 | | |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 16 | | |
| ED041G: Sulfate (Turbidimetric) as SO4 | 2- by DA | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | 4 | | |
| ED045G: Chloride Discrete analyser | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 23 | | |
| ED093F: Dissolved Major Cations | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 1 | | |
| Magnesium | 7439-95-4 | 1 | mg/L | 1 | | |
| Sodium | 7440-23-5 | 1 | mg/L | 14 | | |
| Potassium | 7440-09-7 | 1 | mg/L | 1 | | |
| EG020F: Dissolved Metals by ICP-MS | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | 0.04 | | |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | | |
| Arsenic | 7440-38-2 | 0.001 | mg/L | <0.001 | | |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | | |
| Barium | 7440-39-3 | 0.001 | mg/L | 0.029 | | |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | | |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | | |
| Cobalt | 7440-48-4 | 0.001 | mg/L | <0.001 | | |
| Copper | 7440-50-8 | 0.001 | mg/L | <0.001 | | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | NR | | |
|---------------------------------------|---------------------|--------------|----------------|---------------|------|------|
| | Cl | ient samplir | ng date / time | [09-SEP-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1420206-001 | | |
| EG020F: Dissolved Metals by ICP-MS | - Continued | | | | | |
| Lead | 7439-92-1 | 0.001 | mg/L | <0.001 | | |
| Manganese | 7439-96-5 | 0.001 | mg/L | 0.024 | | |
| Molybdenum | 7439-98-7 | 0.001 | mg/L | 0.001 | | |
| Nickel | 7440-02-0 | 0.001 | mg/L | <0.001 | | |
| Selenium | 7782-49-2 | 0.01 | mg/L | <0.01 | | |
| Strontium | 7440-24-6 | 0.001 | mg/L | 0.022 | | |
| Uranium | 7440-61-1 | 0.001 | mg/L | <0.001 | | |
| Vanadium | 7440-62-2 | 0.01 | mg/L | <0.01 | | |
| Zinc | 7440-66-6 | 0.005 | mg/L | <0.005 | | |
| Boron | 7440-42-8 | 0.05 | mg/L | <0.05 | | |
| Iron | 7439-89-6 | 0.05 | mg/L | 0.23 | | |
| Bromine | 7726-95-6 | 0.1 | mg/L | <0.1 | | |
| EG035F: Dissolved Mercury by FIMS | | | | | | |
| Mercury | 7439-97-6 | 0.0001 | mg/L | <0.0001 | | |
| EG052G: Silica by Discrete Analyser | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 3.06 | | |
| EK026SF: Total CN by Segmented Flo | ow Analyser | | | | | |
| Total Cyanide | 57-12-5 | 0.004 | mg/L | <0.004 | | |
| EK040P: Fluoride by PC Titrator | | | | | | |
| Fluoride | 16984-48-8 | 0.1 | mg/L | <0.1 | | |
| EK055G: Ammonia as N by Discrete A | nalyser | | | | | |
| Ammonia as N | 7664-41-7 | 0.01 | mg/L | 0.03 | | |
| EK057G: Nitrite as N by Discrete Ana | lyser | | | | | |
| Nitrite as N | 14797-65-0 | 0.01 | mg/L | <0.01 | | |
| EK058G: Nitrate as N by Discrete Ana | alyser | | | | | |
| Nitrate as N | 14797-55-8 | 0.01 | mg/L | 0.18 | | |
| EK059G: Nitrite plus Nitrate as N (NO | x) by Discrete Ana | lyser | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | 0.18 | | |
| EK067G: Total Phosphorus as P by Di | iscrete Analyser | | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | <0.01 | | |
| EK071G: Reactive Phosphorus as P b | y discrete analyser | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | <0.01 | | |
| EN055: Ionic Balance | | | | | | |
| | | | | | | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | NR | | |
|--------------------------------------|------------|-------------|----------------|---------------|------|------|
| | Cl | ient sampli | ng date / time | [09-SEP-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1420206-001 | | |
| EN055: Ionic Balance - Continued | | | | | | |
| Total Anions | | 0.01 | meq/L | 1.05 | | |
| Total Cations | | 0.01 | meq/L | 0.83 | | |
| EP005: Total Organic Carbon (TOC) | | | | | | |
| Total Organic Carbon | | 1 | mg/L | 4 | | |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 10 | | |
| Ethene | 74-85-1 | 10 | µg/L | <10 | | |
| Ethane | 74-84-0 | 10 | µg/L | <10 | | |
| Propene | 115-07-1 | 10 | µg/L | <10 | | |
| Propane | 74-98-6 | 10 | µg/L | <10 | | |
| Butene | 25167-67-3 | 10 | µg/L | <10 | | |
| Butane | 106-97-8 | 10 | µg/L | <10 | | |
| EP075(SIM)A: Phenolic Compounds | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | <1.0 | | |
| 2-Chlorophenol | 95-57-8 | 1.0 | μg/L | <1.0 | | |
| 2-Methylphenol | 95-48-7 | 1.0 | μg/L | <1.0 | | |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | μg/L | <2.0 | | |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | | |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | μg/L | <1.0 | | |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | μg/L | <1.0 | | |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | | |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | μg/L | <1.0 | | |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | μg/L | <1.0 | | |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | μg/L | <1.0 | | |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | | |
| EP075(SIM)B: Polynuclear Aromatic Hy | drocarbons | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | | |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | | |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | | |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | | |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | | |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | | |

Page : 6 of 8 Work Order : ES1420206 Amendment 1 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2193361A



| | Cli | ient samplir | ng date / time | [09-SEP-2014] | | |
|--|------------------|--------------|----------------|---------------|------|------|
| Compound | CAS Number | LOR | Unit | ES1420206-001 | | |
| EP075(SIM)B: Polynuclear Aromatic Hydr | ocarbons - Cont | inued | | | | |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | | |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | | |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | | |
| Benzo(b+j)fluoranthene 20 |)5-99-2 205-82-3 | 1.0 | µg/L | <1.0 | | |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | | |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | | |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | | |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | | |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | | |
| [^] Sum of polycyclic aromatic hydrocarbons | | 0.5 | µg/L | <0.5 | | |
| ^ Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | | |
| EP080/071: Total Petroleum Hydrocarbon | IS | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | <20 | | |
| C10 - C14 Fraction | | 50 | µg/L | <50 | | |
| C15 - C28 Fraction | | 100 | µg/L | <100 | | |
| C29 - C36 Fraction | | 50 | µg/L | <50 | | |
| [^] C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | | |
| EP080/071: Total Recoverable Hydrocarb | ons - NEPM 201 | 3 Fractior | ıs | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | <20 | | |
| C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | <20 | | |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | | |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | | |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | | |
| >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | | |
| ^ >C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | | |
| EP080: BTEXN | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | | |
| Toluene | 108-88-3 | 2 | µg/L | <2 | | |
| Ethylbenzene | 100-41-4 | 2 | µg/L | <2 | | |
| meta- & para-Xylene 10 | 08-38-3 106-42-3 | 2 | µg/L | <2 | | |
| ortho-Xylene | 95-47-6 | 2 | µg/L | <2 | | |
| ^ Total Xylenes | 1330-20-7 | 2 | μg/L | <2 | | |



| Sub-Matrix: WATER (Matrix: WATER) | atrix: WATER (Matrix: WATER) Client sample ID | | | | | |
|-----------------------------------|---|------------|----------------|---------------|------|------|
| | Cli | ent sampli | ng date / time | [09-SEP-2014] | | |
| Compound | CAS Number | LOR | Unit | ES1420206-001 | | |
| EP080: BTEXN - Continued | | | | | | |
| Sum of BTEX | | 1 | µg/L | <1 | | |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | | |
| EP075(SIM)S: Phenolic Compound Su | urrogates | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 31.7 | | |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 62.6 | | |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 74.6 | | |
| EP075(SIM)T: PAH Surrogates | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 72.3 | | |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 87.2 | | |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 79.4 | | |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 104 | | |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 95.2 | | |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 90.1 | | |

ALS

Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) |
|---|------------|----------|------------|
| Compound | CAS Number | Low | High |
| EP075(SIM)S: Phenolic Compound Surrogates | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 |
| EP075(SIM)T: PAH Surrogates | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 |
| EP080S: TPH(V)/BTEX Surrogates | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 |
| Toluene-D8 | 2037-26-5 | 79 | 131 |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 |



| | CERTI | FICATE OF ANALYSIS | ; |
|--------------|---------------------------------|-------------------------|---|
| Work Order | ES1500565 | Page | : 1 of 8 |
| Amendment | : 3 | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney |
| Contact | : MS ANDREA MADDEN | Contact | : Loren Schiavon |
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| Telephone | +61 02 9272 5100 | Telephone | : +61 2 8784 8503 |
| Facsimile | : +61 02 9272 5101 | Facsimile | : +61 2 8784 8500 |
| Project | : 2268518A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement |
| Order number | : | | |
| C-O-C number | : | Date Samples Received | : 13-JAN-2015 |
| Sampler | : CS | Issue Date | : 19-MAR-2015 |
| Site | : | | |
| | | No. of samples received | : 4 |
| Quote number | : SY/743/14 V2 | No. of samples analysed | : 4 |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

Accredited for compliance with

ISO/IEC 17025.

- General Comments
- Analytical Results
- Surrogate Control Limits



| NATA Accredited Laboratory 825 | Signatories |
|--------------------------------|-------------|
|--------------------------------|-------------|

This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11.

| Signatories | Position | Accreditation Category |
|---------------------------------|--|--|
| Ankit Joshi Celine Conceicao | Inorganic Chemist Senior Spectroscopist | Sydney Inorganics Sydney Inorganics |
| Edwandy Fadjar | Organic Coordinator | Sydney Organics |

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General Comments

The analytical procedures used by the Environmental Division have been developed from established internationally recognized procedures such as those published by the USEPA, APHA, AS and NEPM. In house developed procedures are employed in the absence of documented standards or by client request.

Where moisture determination has been performed, results are reported on a dry weight basis.

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

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

When sampling time information is not provided by the client, sampling dates are shown without a time component. In these instances, the time component has been assumed by the laboratory for processing purposes.

Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

Key: CAS Number = CAS registry number from database maintained by Chemical Abstracts Services. The Chemical Abstracts Service is a division of the American Chemical Society. LOR = Limit of reporting

* = This result is computed from individual analyte detections at or above the level of reporting

- Benzo(a)pyrene Toxicity Equivalent Quotient (TEQ) is the sum total of the concentration of the eight carcinogenic PAHs multiplied by their Toxicity Equivalence Factor (TEF) relative to Benzo(a)pyrene. TEF values are provided in brackets as follows: Benz(a)anthracene (0.1), Chrysene (0.01), Benzo(b+j) & Benzo(k)fluoranthene (0.1), Benzo(a)pyrene (1.0), Indeno(1.2.3.cd)pyrene (0.1), Dibenz(a.h)anthracene (1.0), Benzo(g.h.i)perylene (0.01). Less than LOR results for 'TEQ Zero' are treated as zero.
- EG020: Bromine quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- This report has been amended and re-released to allow the reporting of additional analytical data.



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB01 | MPMB02 | MPMB03 | MPMB04 | |
|---------------------------------------|-------------|---------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | C | lient samplii | ng date / time | 12-JAN-2015 12:30 | 12-JAN-2015 12:30 | 12-JAN-2015 11:30 | 12-JAN-2015 12:15 | |
| Compound | CAS Number | LOR | Unit | ES1500565-001 | ES1500565-002 | ES1500565-003 | ES1500565-004 | |
| EA005P: pH by PC Titrator | | | | | | | | |
| pH Value | | 0.01 | pH Unit | 5.65 | 6.86 | 7.54 | 9.23 | |
| EA010P: Conductivity by PC Titrator | | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | µS/cm | 930 | 967 | 1110 | 1030 | |
| EA015: Total Dissolved Solids | | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 514 | 455 | 564 | 586 | |
| EA025: Suspended Solids | | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | 82 | 13 | 7 | 10 | |
| ED009: Anions | | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 0.764 | 0.596 | 0.868 | 0.365 | |
| ED037P: Alkalinity by PC Titrator | | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | <1 | <1 | |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | <1 | <1 | <1 | 88 | |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 17 | 213 | 468 | 222 | |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 17 | 213 | 468 | 310 | |
| ED041G: Sulfate (Turbidimetric) as SO | 4 2- by DA | | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | 3 | 4 | <1 | <1 | |
| ED045G: Chloride Discrete analyser | | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 245 | 161 | 70 | 120 | |
| ED093F: Dissolved Major Cations | | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 13 | 37 | 88 | 15 | |
| Magnesium | 7439-95-4 | 1 | mg/L | 22 | 32 | 22 | 6 | |
| Sodium | 7440-23-5 | 1 | mg/L | 106 | 107 | 104 | 180 | |
| Potassium | 7440-09-7 | 1 | mg/L | 1 | 4 | 12 | 13 | |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | 0.07 | <0.01 | <0.01 | <0.01 | |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Arsenic | 7440-38-2 | 0.001 | mg/L | <0.001 | 0.003 | 0.025 | 0.004 | |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Barium | 7440-39-3 | 0.001 | mg/L | 0.691 | 0.629 | 3.48 | 1.23 | |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | <0.0001 | |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Cobalt | 7440-48-4 | 0.001 | mg/L | 0.045 | <0.001 | 0.004 | <0.001 | |
| Copper | 7440-50-8 | 0.001 | mg/L | 0.016 | <0.001 | <0.001 | 0.001 | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB01 | MPMB02 | MPMB03 | MPMB04 | |
|--|----------------------|--------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | CI | ient samplir | ng date / time | 12-JAN-2015 12:30 | 12-JAN-2015 12:30 | 12-JAN-2015 11:30 | 12-JAN-2015 12:15 | |
| Compound | CAS Number | LOR | Unit | ES1500565-001 | ES1500565-002 | ES1500565-003 | ES1500565-004 | |
| EG020F: Dissolved Metals by ICP-MS - Co | ontinued | | | | | | | |
| Lead | 7439-92-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Manganese | 7439-96-5 | 0.001 | mg/L | 0.517 | 0.160 | 0.047 | 0.009 | |
| Molybdenum | 7439-98-7 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | 0.004 | |
| Nickel | 7440-02-0 | 0.001 | mg/L | 0.019 | <0.001 | 0.004 | <0.001 | |
| Selenium | 7782-49-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | |
| Strontium | 7440-24-6 | 0.001 | mg/L | 0.155 | 0.355 | 0.957 | 0.300 | |
| Tin | 7440-31-5 | 0.001 | mg/L | <0.001 | 0.002 | <0.001 | <0.001 | |
| Uranium | 7440-61-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Vanadium | 7440-62-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | |
| Zinc | 7440-66-6 | 0.005 | mg/L | 0.062 | 0.058 | <0.005 | 0.086 | |
| Boron | 7440-42-8 | 0.05 | mg/L | <0.05 | <0.05 | <0.05 | <0.05 | |
| Iron | 7439-89-6 | 0.05 | mg/L | <0.05 | 4.04 | 0.86 | <0.05 | |
| Bromine | 7726-95-6 | 0.1 | mg/L | 0.5 | 0.4 | 0.2 | 0.2 | |
| EG035F: Dissolved Mercury by FIMS | | | | | | | | |
| Mercury | 7439-97-6 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | <0.0001 | |
| EG052G: Silica by Discrete Analyser | | | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 18.3 | 12.3 | 8.50 | 4.25 | |
| EK026SF: Total CN by Segmented Flow | Analyser | | | | | | | |
| Total Cyanide | 57-12-5 | 0.004 | mg/L | <0.004 | <0.004 | <0.004 | <0.004 | |
| EK040P: Fluoride by PC Titrator | | | | | | | | |
| Fluoride | 16984-48-8 | 0.1 | mg/L | <0.1 | 0.2 | 0.2 | 0.5 | |
| EK055G: Ammonia as N by Discrete Anal | vser | | | | | | | |
| Ammonia as N | 7664-41-7 | 0.01 | mg/L | 0.04 | 0.08 | 0.94 | 0.97 | |
| EK055G-NH4: Ammonium as N by DA | | | | | | | | |
| Ammonium as N | | 0.01 | mg/L | 0.04 | 0.08 | 0.94 | 0.47 | |
| EK057G: Nitrite as N by Discrete Analyse | er | | | | | | | |
| Nitrite as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | |
| EK058G: Nitrate as N by Discrete Analys | er | | | | | | | |
| Nitrate as N | 14797-55-8 | 0.01 | mg/L | 0.15 | <0.01 | <0.01 | <0.01 | |
| EK059G: Nitrite plus Nitrate as N (NOx) | by Discrete Ana | lyser | | | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | 0.15 | <0.01 | <0.01 | <0.01 | |
| EK067G: Total Phosphorus as P by Discr | ete Analyse <u>r</u> | | | | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | 0.15 | 0.11 | 0.07 | <0.01 | |
| | | | | | 1 | 1 | • | |



| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | MPMB01 | MPMB02 | MPMB03 | MPMB04 | |
|-------------------------------------|-------------------|------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | Clie | ent sampli | ng date / time | 12-JAN-2015 12:30 | 12-JAN-2015 12:30 | 12-JAN-2015 11:30 | 12-JAN-2015 12:15 | |
| Compound | CAS Number | LOR | Unit | ES1500565-001 | ES1500565-002 | ES1500565-003 | ES1500565-004 | |
| EK071G: Reactive Phosphorus as P by | discrete analyser | | | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | |
| EK084: Un-ionized Hydrogen Sulfide | | | | | | | | |
| Unionized Hydrogen Sulfide | | 0.1 | mg/L | <0.1 | <0.1 | <0.1 | <0.1 | |
| EK085M: Sulfide as S2- | | | | | | | | |
| Sulfide as S2- | 18496-25-8 | 0.1 | mg/L | <0.1 | <0.1 | <0.1 | <0.1 | |
| EN055: Ionic Balance | | | | | | | | |
| Total Anions | | 0.01 | meq/L | 7.31 | 8.88 | 11.3 | 9.58 | |
| Total Cations | | 0.01 | meq/L | 7.10 | 9.24 | 11.0 | 9.40 | |
| Ionic Balance | | 0.01 | % | 1.52 | 1.95 | 1.32 | 0.96 | |
| EP005: Total Organic Carbon (TOC) | | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | <1 | <1 | <1 | 14 | |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | <10 | 323 | 37000 | 41100 | |
| Ethene | 74-85-1 | 10 | µg/L | <10 | <10 | <10 | <10 | |
| Ethane | 74-84-0 | 10 | µg/L | <10 | <10 | <10 | <10 | |
| Propene | 115-07-1 | 10 | µg/L | <10 | <10 | <10 | <10 | |
| Propane | 74-98-6 | 10 | µg/L | <10 | <10 | <10 | <10 | |
| Butene | 25167-67-3 | 10 | µg/L | <10 | <10 | <10 | <10 | |
| Butane | 106-97-8 | 10 | µg/L | <10 | <10 | <10 | <10 | |
| EP075(SIM)A: Phenolic Compounds | | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | 3.1 | |
| 2-Chlorophenol | 95-57-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB01 | MPMB02 | MPMB03 | MPMB04 | |
|---|--------------------|--------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | Cl | ient samplii | ng date / time | 12-JAN-2015 12:30 | 12-JAN-2015 12:30 | 12-JAN-2015 11:30 | 12-JAN-2015 12:15 | |
| Compound | CAS Number | LOR | Unit | ES1500565-001 | ES1500565-002 | ES1500565-003 | ES1500565-004 | |
| EP075(SIM)B: Polynuclear Aromatic I | Hydrocarbons - Con | tinued | | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | 3.3 | <1.0 | <1.0 | |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | 4.4 | <1.0 | <1.0 | |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | 3.7 | <1.0 | <1.0 | |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | 5.6 | <1.0 | <1.0 | |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(b+j)fluoranthene | 205-99-2 205-82-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Sum of polycyclic aromatic hydrocarbon | IS | 0.5 | µg/L | <0.5 | 17.0 | <0.5 | <0.5 | |
| ^ Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | |
| EP080/071: Total Petroleum Hydroca | rbons | | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | <20 | <20 | <20 | 80 | |
| C10 - C14 Fraction | | 50 | µg/L | <50 | 450 | <50 | <50 | |
| C15 - C28 Fraction | | 100 | µg/L | <100 | 120 | <100 | <100 | |
| C29 - C36 Fraction | | 50 | µg/L | <50 | <50 | <50 | <50 | |
| [^] C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | 570 | <50 | <50 | |
| EP080/071: Total Recoverable Hydrod | carbons - NEPM 201 | 3 Fractio | าร | | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | <20 | <20 | <20 | 90 | |
| [^] C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | <20 | <20 | <20 | <20 | |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | 490 | <100 | <100 | |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | 100 | <100 | <100 | |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | |
| ^ >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | 590 | <100 | <100 | |
| >C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | 490 | <100 | <100 | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB01 | MPMB02 | MPMB03 | MPMB04 | |
|-----------------------------------|-------------------|------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | Cli | ent sampli | ng date / time | 12-JAN-2015 12:30 | 12-JAN-2015 12:30 | 12-JAN-2015 11:30 | 12-JAN-2015 12:15 | |
| Compound | CAS Number | LOR | Unit | ES1500565-001 | ES1500565-002 | ES1500565-003 | ES1500565-004 | |
| EP080: BTEXN | | | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | <1 | <1 | |
| Toluene | 108-88-3 | 2 | µg/L | <2 | <2 | <2 | 79 | |
| Ethylbenzene | 100-41-4 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| meta- & para-Xylene | 108-38-3 106-42-3 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| ortho-Xylene | 95-47-6 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| Sum of BTEX | | 1 | μg/L | <1 | <1 | <1 | 79 | |
| Naphthalene | 91-20-3 | 5 | μg/L | <5 | <5 | <5 | <5 | |
| EP075(SIM)S: Phenolic Compound S | Surrogates | | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 25.9 | 26.3 | 24.8 | 18.0 | |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 62.9 | 60.0 | 58.6 | 28.9 | |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 67.3 | 70.7 | 59.8 | 23.1 | |
| EP075(SIM)T: PAH Surrogates | | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 71.1 | 69.9 | 72.7 | 59.6 | |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 82.5 | 73.8 | 79.4 | 66.0 | |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 86.2 | 86.0 | 80.7 | 68.7 | |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 112 | 112 | 112 | 111 | |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 126 | 126 | 125 | 124 | |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 109 | 109 | 112 | 114 | |

ALS

Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) |
|---|------------|----------|------------|
| Compound | CAS Number | Low | High |
| EP075(SIM)S: Phenolic Compound Surrogates | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 |
| EP075(SIM)T: PAH Surrogates | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 |
| EP080S: TPH(V)/BTEX Surrogates | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 |
| Toluene-D8 | 2037-26-5 | 79 | 131 |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 |



| CERTIFICATE OF ANALYSIS | | | | | | | | | |
|-------------------------|---------------------------------|-------------------------|---|--|--|--|--|--|--|
| Work Order | ES1500625 | Page | : 1 of 8 | | | | | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney | | | | | | |
| Contact | : MS ANDREA MADDEN | Contact | : Loren Schiavon | | | | | | |
| Address | : GPO BOX 5394 | Address | : 277-289 Woodpark Road Smithfield NSW Australia 2164 | | | | | | |
| | SYDNEY NSW, AUSTRALIA 2001 | | | | | | | | |
| E-mail | : amadden@pb.com.au | E-mail | : loren.schiavon@alsglobal.com | | | | | | |
| Telephone | +61 02 9272 5100 | Telephone | : +61 2 8784 8503 | | | | | | |
| Facsimile | : +61 02 9272 5101 | Facsimile | : +61 2 8784 8500 | | | | | | |
| Project | : 2268518A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement | | | | | | |
| Order number | : | | | | | | | | |
| C-O-C number | : | Date Samples Received | : 14-JAN-2015 | | | | | | |
| Sampler | : CS | Issue Date | : 22-JAN-2015 | | | | | | |
| Site | : | | | | | | | | |
| | | No. of samples received | : 2 | | | | | | |
| Quote number | : SY/743/14 V2 | No. of samples analysed | : 2 | | | | | | |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results
- Surrogate Control Limits

Signatories NATA Accredited Laboratory 825 This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11. Accredited for compliance with NATA ISO/IEC 17025. Signatories Position Accreditation Category Inorganic Chemist Ankit Joshi Sydney Inorganics Senior Spectroscopist Celine Conceicao Sydney Inorganics WORLD RECOGNISED ACCREDITATION Organic Coordinator Edwandy Fadjar Sydney Organics Metals Coordinator Shobhna Chandra Sydney Inorganics

Address 277-289 Woodpark Road Smithfield NSW Australia 2164 PHONE +61-2-8784 8555 Facsimile +61-2-8784 8500 Environmental Division Sydney ABN 84 009 936 029 Part of the ALS Group An ALS Limited Company



RIGHT SOLUTIONS RIGHT PARTNER



General Comments

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Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

- EG020: Bromine quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- EK071G: It has been noted that Reactive P is greater than Total P for sample ID(RMB02), however this difference is within the limits of experimental variation.



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | RMB02 | RMB03 | | |
|--|-----------------------------|--------|-------------------|-------------------|---------------|------|--|
| | Client sampling date / time | | 13-JAN-2015 14:15 | 13-JAN-2015 15:00 | | | |
| Compound | CAS Number | LOR | Unit | ES1500625-001 | ES1500625-002 | | |
| EA005P: pH by PC Titrator | | | | | | | |
| pH Value | | 0.01 | pH Unit | 7.39 | 9.60 | | |
| EA010P: Conductivity by PC Titrator | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | µS/cm | 10500 | 7690 | | |
| EA015: Total Dissolved Solids | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 5980 | 4130 | | |
| EA025: Suspended Solids | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | <5 | <5 | | |
| ED009: Anions | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 8.89 | 6.05 | | |
| ED037P: Alkalinity by PC Titrator | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | | |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | <1 | 229 | | |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 792 | 189 | | |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 792 | 418 | | |
| ED041G: Sulfate (Turbidimetric) as SO4 | 2- by DA | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | <1 | <1 | | |
| ED045G: Chloride Discrete analyser | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 3170 | 2290 | | |
| ED093F: Dissolved Major Cations | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 382 | 14 | | |
| Magnesium | 7439-95-4 | 1 | mg/L | 83 | 11 | | |
| Sodium | 7440-23-5 | 1 | mg/L | 1770 | 1570 | | |
| Potassium | 7440-09-7 | 1 | mg/L | 30 | 16 | | |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | <0.01 | <0.01 | | |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | 0.001 | | |
| Arsenic | 7440-38-2 | 0.001 | mg/L | <0.001 | 0.001 | | |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Barium | 7440-39-3 | 0.001 | mg/L | 33.2 | 5.07 | | |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | | |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Cobalt | 7440-48-4 | 0.001 | mg/L | <0.001 | <0.001 | | |
| Copper | 7440-50-8 | 0.001 | mg/L | <0.001 | 0.001 | | |

Page : 4 of 8 Work Order : ES1500625 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Client sample ID | RMB02 | RMB03 | | | |
|--|-----------------------------|------------------|-------------------|-------------------|--|---|--|
| | Client sampling date / time | | 13-JAN-2015 14:15 | 13-JAN-2015 15:00 | | | |
| Compound CAS N | umber LC | DR Unit | ES1500625-001 | ES1500625-002 | | | |
| EG020F: Dissolved Metals by ICP-MS - Continued | | | | | | | |
| Lead 743 | 9-92-1 0.0 | 01 mg/L | <0.001 | <0.001 | | | |
| Manganese 743 | 9-96-5 0.0 | 01 mg/L | 0.024 | 0.001 | | | |
| Molybdenum 743 | 9-98-7 0.0 | 01 mg/L | <0.001 | 0.004 | | | |
| Nickel 744 | 0-02-0 0.0 | 01 mg/L | 0.003 | <0.001 | | | |
| Selenium 778 | 2-49-2 0.0 |)1 mg/L | <0.01 | <0.01 | | | |
| Strontium 744 | 0-24-6 0.0 | 01 mg/L | 9.97 | 1.83 | | | |
| Tin 744 | 0-31-5 0.0 | 01 mg/L | <0.001 | <0.001 | | | |
| Uranium 744 | 0-61-1 0.0 | 01 mg/L | <0.001 | <0.001 | | | |
| Vanadium 744 |)-62-2 0.0 |)1 mg/L | <0.01 | <0.01 | | | |
| Zinc 744 | 0-66-6 0.0 | 05 mg/L | 0.033 | 0.039 | | | |
| |)-42-8 0.0 |)5 mg/L | <0.05 | 0.15 | | | |
| Iron 743 | 9-89-6 0.0 |)5 mg/L | 4.73 | <0.05 | | | |
| Bromine 772 | 6-95-6 0. | 1 mg/L | 7.8 | 5.7 | | | |
| EG035F: Dissolved Mercury by FIMS | | | | | | | |
| | 9-97-6 0.00 | 001 mg/L | <0.0001 | <0.0001 | | | |
| EG052G: Silica by Discrete Analyser | | | | | | | |
| Reactive Silica | 0.0 |)5 mg/L | 9.66 | 6.24 | | | |
| EK026SF: Total CN by Segmented Flow Analyser | | | | | | | |
| | 7-12-5 0.0 | 04 mg/L | <0.004 | <0.004 | | | |
| EK040P: Fluoride by PC Titrator | | | | | | | |
| | 4-48-8 0. | 1 mg/L | 0.2 | 0.4 | | | |
| EK055G: Ammonia as N by Discrete Analyser | | | | | | | |
| | 4-41-7 0.0 |)1 mg/L | 4.55 | 3.36 | | | |
| EK055G-NH4: Ammonium as N by DA | | | | | | | |
| Ammonium as N | 0.0 |)1 mg/L | 4.55 | 1.23 | | | |
| EK057G: Nitrite as N by Discrete Analyser | | | | | | | |
| Nitrite as N | 0.0 |)1 mg/L | <0.01 | <0.01 | | | |
| EK058G: Nitrate as N by Discrete Analyser | | | | | | | |
| | 7-55-8 0.0 |)1 mg/L | <0.01 | <0.01 | | | |
| EK059G: Nitrite plus Nitrate as N (NOx) by Discr | te Analyser | | | | | | |
| Nitrite + Nitrate as N | 0.0 | | <0.01 | <0.01 | | | |
| EK067G: Total Phosphorus as P by Discrete Anal | | | | | | | |
| Total Phosphorus as P | 0.0 |)1 mg/L | 0.02 | <0.01 | | | |
| | | | | + | | l | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | RMB02 | RMB03 | | |
|---------------------------------------|-------------------|------------|----------------|-------------------|-------------------|------|--|
| | Clie | ent sampli | ng date / time | 13-JAN-2015 14:15 | 13-JAN-2015 15:00 | | |
| Compound | CAS Number | LOR | Unit | ES1500625-001 | ES1500625-002 | | |
| EK071G: Reactive Phosphorus as P by o | discrete analyser | | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | 0.03 | <0.01 | | |
| EK084: Un-ionized Hydrogen Sulfide | | | | | | | |
| Unionized Hydrogen Sulfide | | 0.1 | mg/L | <0.1 | <0.1 | | |
| EK085M: Sulfide as S2- | | | | | | | |
| Sulfide as S2- | 18496-25-8 | 0.1 | mg/L | 0.3 | 0.2 | | |
| EN055: Ionic Balance | | | | | | | |
| Total Anions | | 0.01 | meq/L | 105 | 73.0 | | |
| Total Cations | | 0.01 | meq/L | 104 | 70.3 | | |
| Ionic Balance | | 0.01 | % | 0.77 | 1.87 | | |
| EP005: Total Organic Carbon (TOC) | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | <1 | 38 | | |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 22800 | 36800 | | |
| Ethene | 74-85-1 | 10 | µg/L | <10 | <10 | | |
| Ethane | 74-84-0 | 10 | µg/L | <10 | 12 | | |
| Propene | 115-07-1 | 10 | µg/L | <10 | <10 | | |
| Propane | 74-98-6 | 10 | µg/L | <10 | <10 | | |
| Butene | 25167-67-3 | 10 | µg/L | <10 | <10 | | |
| Butane | 106-97-8 | 10 | µg/L | <10 | <10 | | |
| EP075(SIM)A: Phenolic Compounds | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2-Chlorophenol | 95-57-8 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | | |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | | |

Page : 6 of 8 Work Order : ES1500625 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | RMB02 | RMB03 | | |
|---|-------------------|-------------|----------------|-------------------|-------------------|------|--|
| | Cl | ient sampli | ng date / time | 13-JAN-2015 14:15 | 13-JAN-2015 15:00 | | |
| Compound | CAS Number | LOR | Unit | ES1500625-001 | ES1500625-002 | | |
| EP075(SIM)B: Polynuclear Aromatic I | | tinued | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(b+j)fluoranthene | 205-99-2 205-82-3 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | | |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | <1.0 | | |
| Sum of polycyclic aromatic hydrocarbon | s | 0.5 | µg/L | <0.5 | <0.5 | | |
| [^] Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | | |
| EP080/071: Total Petroleum Hydrocar | rbons | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | <20 | 90 | | |
| 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 | | |
| EP080/071: Total Recoverable Hydrod | arbons - NEPM 201 | 3 Fractio | าร | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | <20 | 90 | | |
| C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | <20 | 40 | | |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | <100 | | |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | <100 | | |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | | |
| ^ >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | <100 | | |
| >C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | <100 | | |

Page : 7 of 8 Work Order : ES1500625 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | Hatrix: WATER (Matrix: WATER) | | ent sample ID | RMB02 | RMB03 | | |
|-----------------------------------|-------------------------------|-------------|----------------|-------------------|-------------------|------|--|
| | Cli | ent samplii | ng date / time | 13-JAN-2015 14:15 | 13-JAN-2015 15:00 | | |
| Compound | CAS Number | LOR | Unit | ES1500625-001 | ES1500625-002 | | |
| EP080: BTEXN | | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | | |
| Toluene | 108-88-3 | 2 | μg/L | <2 | 46 | | |
| Ethylbenzene | 100-41-4 | 2 | µg/L | <2 | <2 | | |
| meta- & para-Xylene | 108-38-3 106-42-3 | 2 | μg/L | <2 | <2 | | |
| ortho-Xylene | 95-47-6 | 2 | µg/L | <2 | <2 | | |
| ^ Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | | |
| Sum of BTEX | | 1 | µg/L | <1 | 46 | | |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | <5 | | |
| EP075(SIM)S: Phenolic Compound | Surrogates | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 20.6 | 19.3 | | |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 45.9 | 22.9 | | |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 45.2 | 27.6 | | |
| EP075(SIM)T: PAH Surrogates | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 57.9 | 70.5 | | |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 57.0 | 71.1 | | |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 64.8 | 80.6 | | |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 88.9 | 87.2 | | |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 106 | 109 | | |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 110 | 110 | | |

ALS

Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) |
|---|------------|----------|------------|
| Compound | CAS Number | Low | High |
| EP075(SIM)S: Phenolic Compound Surrogates | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 |
| EP075(SIM)T: PAH Surrogates | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 |
| EP080S: TPH(V)/BTEX Surrogates | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 |
| Toluene-D8 | 2037-26-5 | 79 | 131 |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 |



| | CERTIFICATE OF ANALYSIS | | | | | | | | | | |
|--------------|---------------------------------|-------------------------|---|--|--|--|--|--|--|--|--|
| Work Order | ES1500816 | Page | : 1 of 8 | | | | | | | | |
| Amendment | : 1 | | | | | | | | | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney | | | | | | | | |
| Contact | : MS ANDREA MADDEN | Contact | : Loren Schiavon | | | | | | | | |
| Address | : GPO BOX 5394 | Address | : 277-289 Woodpark Road Smithfield NSW Australia 2164 | | | | | | | | |
| | SYDNEY NSW, AUSTRALIA 2001 | | | | | | | | | | |
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| Facsimile | : +61 02 9272 5101 | Facsimile | : +61 2 8784 8500 | | | | | | | | |
| Project | : 2268518A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement | | | | | | | | |
| Order number | : | | | | | | | | | | |
| C-O-C number | : | Date Samples Received | : 15-JAN-2015 | | | | | | | | |
| Sampler | : CS | Issue Date | : 19-MAR-2015 | | | | | | | | |
| Site | : | | | | | | | | | | |
| | | No. of samples received | : 4 | | | | | | | | |
| Quote number | : SY/743/14 V2 | No. of samples analysed | : 4 | | | | | | | | |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

Accredited for compliance with

ISO/IEC 17025.

- General Comments
- Analytical Results
- Surrogate Control Limits



| NATA Accredited Laboratory 8 | Signatorie | s |
|------------------------------|----------------|---|
| NATA Accredited Laboratory 8 | S25 Olynatoric | J |

This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11.

| Signatories | Position | Accreditation Category |
|------------------|-----------------------|------------------------|
| Ankit Joshi | Inorganic Chemist | Sydney Inorganics |
| Ashesh Patel | Inorganic Chemist | Sydney Inorganics |
| Celine Conceicao | Senior Spectroscopist | Sydney Inorganics |
| Edwandy Fadjar | Organic Coordinator | Sydney Organics |

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General Comments

The analytical procedures used by the Environmental Division have been developed from established internationally recognized procedures such as those published by the USEPA, APHA, AS and NEPM. In house developed procedures are employed in the absence of documented standards or by client request.

Where moisture determination has been performed, results are reported on a dry weight basis.

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

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

When sampling time information is not provided by the client, sampling dates are shown without a time component. In these instances, the time component has been assumed by the laboratory for processing purposes.

Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

Key: CAS Number = CAS registry number from database maintained by Chemical Abstracts Services. The Chemical Abstracts Service is a division of the American Chemical Society. LOR = Limit of reporting

* = This result is computed from individual analyte detections at or above the level of reporting

- Benzo(a)pyrene Toxicity Equivalent Quotient (TEQ) is the sum total of the concentration of the eight carcinogenic PAHs multiplied by their Toxicity Equivalence Factor (TEF) relative to Benzo(a)pyrene. TEF values are provided in brackets as follows: Benz(a)anthracene (0.1), Chrysene (0.01), Benzo(b+j) & Benzo(k)fluoranthene (0.1), Benzo(a)pyrene (1.0), Indeno(1.2.3.cd)pyrene (0.1), Dibenz(a.h)anthracene (1.0), Benzo(g.h.i)perylene (0.01). Less than LOR results for 'TEQ Zero' are treated as zero.
- EG020: 'Bromine/lodine' quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- EK026SF: Poor spike recovery for Total Cyanide due to matrix interferences (confirmed by re-analysis).
- EK071G: It has been noted that Reactive P is greater than Total P on sample 3 and 4, however this difference is within the limits of experimental variation.
- This report has been amended and re-released to allow the reporting of additional analytical data.



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB01 | GLMB02 | GLMB03 | QA | |
|---------------------------------------|-------------|---------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | C | lient samplii | ng date / time | 14-JAN-2015 12:45 | 14-JAN-2015 11:00 | 14-JAN-2015 11:15 | 14-JAN-2015 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1500816-001 | ES1500816-002 | ES1500816-003 | ES1500816-004 | |
| EA005P: pH by PC Titrator | | | | | | | | |
| pH Value | | 0.01 | pH Unit | 7.48 | 7.66 | 8.85 | 8.76 | |
| EA010P: Conductivity by PC Titrator | | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | μS/cm | 8360 | 7140 | 4930 | 4960 | |
| EA015: Total Dissolved Solids | | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 5110 | 4320 | 2560 | 2740 | |
| EA025: Suspended Solids | | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | 22 | 41 | 6 | 10 | |
| ED009: Anions | | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 8.30 | 5.75 | 2.25 | 2.31 | |
| ED037P: Alkalinity by PC Titrator | | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | <1 | <1 | |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | <1 | <1 | 309 | 252 | |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 434 | 595 | 1070 | 1110 | |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 434 | 595 | 1380 | 1370 | |
| ED041G: Sulfate (Turbidimetric) as SO | 4 2- by DA | | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | 135 | <1 | <1 | <1 | |
| ED045G: Chloride Discrete analyser | | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 2470 | 1760 | 734 | 743 | |
| ED093F: Dissolved Major Cations | | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 182 | 246 | 24 | 23 | |
| Magnesium | 7439-95-4 | 1 | mg/L | 186 | 105 | 57 | 55 | |
| Sodium | 7440-23-5 | 1 | mg/L | 1260 | 1040 | 1020 | 1070 | |
| Potassium | 7440-09-7 | 1 | mg/L | 19 | 24 | 36 | 38 | |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | <0.01 | <0.01 | 0.01 | 0.01 | |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Arsenic | 7440-38-2 | 0.001 | mg/L | 0.010 | 0.010 | 0.020 | 0.022 | |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Barium | 7440-39-3 | 0.001 | mg/L | 0.362 | 8.78 | 6.28 | 6.11 | |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | <0.0001 | |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Cobalt | 7440-48-4 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Copper | 7440-50-8 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB01 | GLMB02 | GLMB03 | QA | |
|--|-----------------------------|--------|-------------------|-------------------|-------------------|-------------------|---------------|--|
| | Client sampling date / time | | 14-JAN-2015 12:45 | 14-JAN-2015 11:00 | 14-JAN-2015 11:15 | 14-JAN-2015 15:00 | | |
| Compound | CAS Number | LOR | Unit | ES1500816-001 | ES1500816-002 | ES1500816-003 | ES1500816-004 | |
| EG020F: Dissolved Metals by ICP-MS - Cont | tinued | | | | | | | |
| Lead | 7439-92-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Manganese | 7439-96-5 | 0.001 | mg/L | 0.390 | 0.384 | 0.003 | 0.002 | |
| Molybdenum | 7439-98-7 | 0.001 | mg/L | <0.001 | <0.001 | 0.004 | 0.004 | |
| Nickel | 7440-02-0 | 0.001 | mg/L | <0.001 | 0.003 | <0.001 | <0.001 | |
| Selenium | 7782-49-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | |
| Strontium | 7440-24-6 | 0.001 | mg/L | 4.08 | 3.97 | 1.54 | 1.54 | |
| Tin | 7440-31-5 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Uranium | 7440-61-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | |
| Vanadium | 7440-62-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | |
| Zinc | 7440-66-6 | 0.005 | mg/L | <0.005 | 0.006 | <0.005 | 0.010 | |
| Boron | 7440-42-8 | 0.05 | mg/L | <0.05 | <0.05 | <0.05 | <0.05 | |
| Iron | 7439-89-6 | 0.05 | mg/L | 4.28 | 4.87 | 0.10 | 0.09 | |
| Bromine | 7726-95-6 | 0.1 | mg/L | 6.4 | 5.1 | 2.2 | 2.1 | |
| EG035F: Dissolved Mercury by FIMS | | | | | | | | |
| Mercury | 7439-97-6 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | <0.0001 | |
| EG052G: Silica by Discrete Analyser | | | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 7.38 | 16.0 | 11.0 | 11.2 | |
| EK026SF: Total CN by Segmented Flow An | nalvser | | | | | | | |
| Total Cyanide | 57-12-5 | 0.004 | mg/L | <0.004 | <0.004 | <0.004 | <0.004 | |
| EK040P: Fluoride by PC Titrator | | | | | | | | |
| Fluoride | 16984-48-8 | 0.1 | mg/L | 0.2 | 0.3 | <0.1 | <0.1 | |
| EK055G: Ammonia as N by Discrete Analys | ser | | | | | | | |
| Ammonia as N | 7664-41-7 | 0.01 | mg/L | 1.54 | 2.20 | 2.76 | 2.79 | |
| EK055G-NH4: Ammonium as N by DA | | | | | | | | |
| Ammonium as N | | 0.01 | mg/L | 1.51 | 2.19 | 2.00 | 2.02 | |
| EK057G: Nitrite as N by Discrete Analyser | | | | | | | | |
| Nitrite as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | |
| EK058G: Nitrate as N by Discrete Analyser | | | | | | | | |
| Nitrate as N | 14797-55-8 | 0.01 | mg/L | <0.01 | 0.01 | <0.01 | <0.01 | |
| EK059G: Nitrite plus Nitrate as N (NOx) by | Discrete Ana | lyser | | | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | <0.01 | 0.01 | <0.01 | <0.01 | |
| EK067G: Total Phosphorus as P by Discret | e Anal <u>yser</u> | | | | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | 0.04 | 0.07 | <0.01 | <0.01 | |
| | | | | | 1 | | 1 | |



| Sub-Matrix: WATER (Matrix: WATER) | | Client sample ID | | GLMB01 | GLMB02 | GLMB03 | QA | |
|-------------------------------------|-----------------------------|------------------|-------|-------------------|-------------------|-------------------|-------------------|--|
| | Client sampling date / time | | | 14-JAN-2015 12:45 | 14-JAN-2015 11:00 | 14-JAN-2015 11:15 | 14-JAN-2015 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1500816-001 | ES1500816-002 | ES1500816-003 | ES1500816-004 | |
| EK071G: Reactive Phosphorus as P by | discrete analyser | | | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | <0.01 | 0.05 | 0.03 | 0.02 | |
| EK084: Un-ionized Hydrogen Sulfide | | | | | | | | |
| Unionized Hydrogen Sulfide | | 0.1 | mg/L | <0.1 | <0.1 | <0.1 | <0.1 | |
| EK085F: Dissolved Sulfide as S2- | | | | | | | | |
| Dissolved Sulfide as S2- | 18496-25-8 | 0.1 | mg/L | <0.1 | <0.1 | <0.1 | <0.1 | |
| EN055: Ionic Balance | | | | | | | | |
| Total Anions | | 0.01 | meq/L | 81.2 | 61.5 | 48.3 | 48.3 | |
| Total Cations | | 0.01 | meq/L | 79.7 | 66.8 | 51.2 | 53.2 | |
| Ionic Balance | | 0.01 | % | 0.93 | 4.07 | 2.87 | 4.74 | |
| EP005: Total Organic Carbon (TOC) | | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | <1 | <1 | <1 | <1 | |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 33900 | 18400 | 18200 | 21900 | |
| Ethene | 74-85-1 | 10 | μg/L | <10 | <10 | <10 | <10 | |
| Ethane | 74-84-0 | 10 | μg/L | 1650 | 798 | 60 | 77 | |
| Propene | 115-07-1 | 10 | μg/L | <10 | <10 | <10 | <10 | |
| Propane | 74-98-6 | 10 | μg/L | 379 | 204 | 15 | 20 | |
| Butene | 25167-67-3 | 10 | μg/L | <10 | <10 | <10 | <10 | |
| Butane | 106-97-8 | 10 | µg/L | 61 | 38 | <10 | <10 | |
| EP075(SIM)A: Phenolic Compounds | | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2-Chlorophenol | 95-57-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB01 | GLMB02 | GLMB03 | QA | |
|---|---------------------|-------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | Cl | ient sampli | ng date / time | 14-JAN-2015 12:45 | 14-JAN-2015 11:00 | 14-JAN-2015 11:15 | 14-JAN-2015 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1500816-001 | ES1500816-002 | ES1500816-003 | ES1500816-004 | |
| EP075(SIM)B: Polynuclear Aromatic I | Hydrocarbons - Cont | inued | | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Acenaphthylene | 208-96-8 | 1.0 | μg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Fluorene | 86-73-7 | 1.0 | μg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(b+j)fluoranthene | 205-99-2 205-82-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | μg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | μg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | μg/L | <1.0 | <1.0 | <1.0 | <1.0 | |
| Sum of polycyclic aromatic hydrocarbon | s | 0.5 | μg/L | <0.5 | <0.5 | <0.5 | <0.5 | |
| ^ Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | |
| EP080/071: Total Petroleum Hydroca | rbons | | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | 110 | 80 | 210 | 190 | |
| C10 - C14 Fraction | | 50 | µg/L | <50 | <50 | <50 | <50 | |
| C15 - C28 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | |
| C29 - C36 Fraction | | 50 | µg/L | <50 | <50 | <50 | <50 | |
| [^] C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | <50 | <50 | <50 | |
| EP080/071: Total Recoverable Hydrod | carbons - NEPM 201 | 3 Fractio | ns | | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | 110 | 80 | 210 | 200 | |
| [^] C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | 90 | 80 | 100 | 100 | |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | <100 | <100 | <100 | |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | |
| >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | <100 | <100 | <100 | |
| >C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | <100 | <100 | <100 | |



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB01 | GLMB02 | GLMB03 | QA | |
|-----------------------------------|-------------------|------------|----------------|-------------------|-------------------|-------------------|-------------------|--|
| | Cli | ent sampli | ng date / time | 14-JAN-2015 12:45 | 14-JAN-2015 11:00 | 14-JAN-2015 11:15 | 14-JAN-2015 15:00 | |
| Compound | CAS Number | LOR | Unit | ES1500816-001 | ES1500816-002 | ES1500816-003 | ES1500816-004 | |
| EP080: BTEXN | | | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | <1 | <1 | |
| Toluene | 108-88-3 | 2 | µg/L | 19 | <2 | 106 | 99 | |
| Ethylbenzene | 100-41-4 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| meta- & para-Xylene | 108-38-3 106-42-3 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| ortho-Xylene | 95-47-6 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | <2 | <2 | |
| Sum of BTEX | | 1 | µg/L | 19 | <1 | 106 | 99 | |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | <5 | <5 | <5 | |
| EP075(SIM)S: Phenolic Compound | Surrogates | | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 23.6 | 24.0 | 23.1 | 22.3 | |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 53.2 | 56.9 | 46.5 | 40.9 | |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 72.2 | 76.4 | 50.1 | 42.9 | |
| EP075(SIM)T: PAH Surrogates | | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 82.9 | 70.8 | 79.7 | 74.2 | |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 69.7 | 77.1 | 72.5 | 72.8 | |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 82.4 | 84.6 | 81.3 | 84.5 | |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 90.9 | 95.5 | 99.9 | 96.1 | |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 98.3 | 103 | 101 | 99.4 | |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 88.1 | 87.8 | 88.0 | 85.4 | |



Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) |
|---|------------|----------|------------|
| Compound | CAS Number | Low | High |
| EP075(SIM)S: Phenolic Compound Surrogates | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 |
| EP075(SIM)T: PAH Surrogates | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 |
| EP080S: TPH(V)/BTEX Surrogates | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 |
| Toluene-D8 | 2037-26-5 | 79 | 131 |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 |



| | CERTIFICATE OF ANALYSIS | | | | | | | | | | |
|--------------|---------------------------------|-------------------------|---|--|--|--|--|--|--|--|--|
| Work Order | ES1508003 | Page | : 1 of 8 | | | | | | | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney | | | | | | | | |
| Contact | : MS CAROLINA SARDELLA | Contact | : Client Services | | | | | | | | |
| Address | : GPO BOX 5394 | Address | : 277-289 Woodpark Road Smithfield NSW Australia 2164 | | | | | | | | |
| | SYDNEY NSW, AUSTRALIA 2001 | | | | | | | | | | |
| E-mail | : csardella@pb.com.au | E-mail | : sydney@alsglobal.com | | | | | | | | |
| Telephone | : +61 02 92725100 | Telephone | : +61-2-8784 8555 | | | | | | | | |
| Facsimile | : +61 02 92725101 | Facsimile | : +61-2-8784 8500 | | | | | | | | |
| Project | : 2268518A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement | | | | | | | | |
| Order number | : | | | | | | | | | | |
| C-O-C number | : | Date Samples Received | : 08-APR-2015 | | | | | | | | |
| Sampler | : KM & AM | Issue Date | : 15-APR-2015 | | | | | | | | |
| Site | : | | | | | | | | | | |
| | | No. of samples received | : 3 | | | | | | | | |
| Quote number | : SY/239/15 | No. of samples analysed | : 3 | | | | | | | | |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results
- Surrogate Control Limits

Signatories NATA Accredited Laboratory 825 This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11. Accredited for compliance with NATA ISO/IEC 17025. Signatories Position Accreditation Category Inorganic Chemist Ankit Joshi Sydney Inorganics Senior Spectroscopist Celine Conceicao Sydney Inorganics WORLD RECOGNISED ACCREDITATION Organic Coordinator Edwandy Fadjar Sydney Organics Senior Organic Chemist Pabi Subba Sydney Organics

Address 277-289 Woodpark Road Smithfield NSW Australia 2164 PHONE +61-2-8784 8555 Facsimile +61-2-8784 8500 Environmental Division Svdnev ABN 84 009 936 029 Part of the ALS Group An ALS Limited Company



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General Comments

The analytical procedures used by the Environmental Division have been developed from established internationally recognized procedures such as those published by the USEPA, APHA, AS and NEPM. In house developed procedures are employed in the absence of documented standards or by client request.

Where moisture determination has been performed, results are reported on a dry weight basis.

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

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

When sampling time information is not provided by the client, sampling dates are shown without a time component. In these instances, the time component has been assumed by the laboratory for processing purposes.

Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

Key: CAS Number = CAS registry number from database maintained by Chemical Abstracts Services. The Chemical Abstracts Service is a division of the American Chemical Society. LOR = Limit of reporting

* = This result is computed from individual analyte detections at or above the level of reporting

- Benzo(a)pyrene Toxicity Equivalent Quotient (TEQ) is the sum total of the concentration of the eight carcinogenic PAHs multiplied by their Toxicity Equivalence Factor (TEF) relative to Benzo(a)pyrene. TEF values are provided in brackets as follows: Benz(a)anthracene (0.1), Chrysene (0.01), Benzo(b+j) & Benzo(k)fluoranthene (0.1), Benzo(a)pyrene (1.0), Indeno(1.2.3.cd)pyrene (0.1), Dibenz(a.h)anthracene (1.0), Benzo(g.h.i)perylene (0.01). Less than LOR results for 'TEQ Zero' are treated as zero.
- ED041G: LOR raised for Sulfate analysis on sample ID: RMB02, due to matrix interferences.
- EG020: Bromine quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- EK026SF:Spike failed for Total Cyanide analysis due to matrix interference(confirmed by re analysis)
- EK067G: It has been noted that Reactive P is greater than Total P for sample ID(GLMB03), however this difference is within the limits of experimental variation.



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB03 | RMB02 | RMB03 | |
|--|-------------|---------------|----------------|-------------------|-------------------|-------------------|------|
| | C | lient samplir | ng date / time | 07-APR-2015 11:00 | 07-APR-2015 15:00 | 07-APR-2015 16:30 | |
| Compound | CAS Number | LOR | Unit | ES1508003-001 | ES1508003-002 | ES1508003-003 | |
| EA005P: pH by PC Titrator | | | | | | | |
| pH Value | | 0.01 | pH Unit | 8.60 | 7.19 | 9.63 | |
| EA010P: Conductivity by PC Titrator | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | μS/cm | 5010 | 10700 | 7730 | |
| EA015: Total Dissolved Solids | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 2750 | 6170 | 4140 | |
| EA025: Suspended Solids | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | <5 | <5 | <5 | |
| ED009: Anions | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 1.98 | 6.79 | 4.68 | |
| ED037P: Alkalinity by PC Titrator | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | <1 | |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | 129 | <1 | 252 | |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 1510 | 895 | 224 | |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 1640 | 895 | 476 | |
| ED041G: Sulfate (Turbidimetric) as SO4 | 2- by DA | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | 19 | <10 | <1 | |
| ED045G: Chloride Discrete analyser | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 774 | 3280 | 2310 | |
| ED093F: Dissolved Major Cations | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 31 | 326 | 10 | |
| Magnesium | 7439-95-4 | 1 | mg/L | 61 | 88 | 11 | |
| Sodium | 7440-23-5 | 1 | mg/L | 1050 | 1890 | 1600 | |
| Potassium | 7440-09-7 | 1 | mg/L | 39 | 30 | 16 | |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | <0.01 | <0.01 | 0.01 | |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | <0.001 | 0.002 | |
| Arsenic | 7440-38-2 | 0.001 | mg/L | 0.032 | <0.001 | 0.001 | |
| Boron | 7440-42-8 | 0.05 | mg/L | 0.05 | <0.05 | 0.17 | |
| Barium | 7440-39-3 | 0.001 | mg/L | 7.63 | 33.0 | 4.64 | |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | |
| Cobalt | 7440-48-4 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |

Page : 4 of 8 Work Order : ES1508003 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | GLMB03 | RMB02 | RMB03 | |
|--|--------------|-------------|----------------|-------------------|-------------------|-------------------|------|
| | Cl | ient sampli | ng date / time | 07-APR-2015 11:00 | 07-APR-2015 15:00 | 07-APR-2015 16:30 | |
| Compound | CAS Number | LOR | Unit | ES1508003-001 | ES1508003-002 | ES1508003-003 | |
| EG020F: Dissolved Metals by ICP-MS - Contin | nued | | | | | | |
| Copper | 7440-50-8 | 0.001 | mg/L | 0.002 | <0.001 | 0.003 | |
| Manganese | 7439-96-5 | 0.001 | mg/L | 0.005 | 0.024 | <0.001 | |
| Nickel | 7440-02-0 | 0.001 | mg/L | <0.001 | 0.001 | <0.001 | |
| Lead | 7439-92-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Selenium | 7782-49-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| Vanadium | 7440-62-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| Zinc | 7440-66-6 | 0.005 | mg/L | 0.008 | 0.031 | 0.051 | |
| Molybdenum | 7439-98-7 | 0.001 | mg/L | 0.003 | <0.001 | 0.005 | |
| Strontium | 7440-24-6 | 0.001 | mg/L | 2.23 | 8.47 | 1.58 | |
| Uranium | 7440-61-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | |
| Iron | 7439-89-6 | 0.05 | mg/L | 0.12 | 4.42 | <0.05 | |
| Bromine | 7726-95-6 | 0.1 | mg/L | 2.0 | 7.0 | 5.4 | |
| EG035F: Dissolved Mercury by FIMS | | | | | | | |
| Mercury | 7439-97-6 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | |
| EG052G: Silica by Discrete Analyser | | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 11.8 | 10.4 | 6.68 | |
| EK026SF: Total CN by Segmented Flow Ana | lvser | | | | | | |
| Total Cyanide | 57-12-5 | 0.004 | mg/L | <0.004 | <0.004 | <0.004 | |
| EK040P: Fluoride by PC Titrator | | | | | | | |
| Fluoride | 16984-48-8 | 0.1 | mg/L | <0.1 | 0.1 | 0.6 | |
| EK055G: Ammonia as N by Discrete Analyse | r | | | | | | |
| Ammonia as N | 7664-41-7 | 0.01 | mg/L | 2.87 | 4.43 | 3.33 | |
| EK057G: Nitrite as N by Discrete Analyser | | | | | | | |
| Nitrite as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| EK058G: Nitrate as N by Discrete Analyser | | | | | | | |
| Nitrate as N | 14797-55-8 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| EK059G: Nitrite plus Nitrate as N (NOx) by D | Discrete Ana | lvser | | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | |
| EK067G: Total Phosphorus as P by Discrete | Analyser | | | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | 0.04 | 0.04 | 0.01 | |
| EK071G: Reactive Phosphorus as P by discr | ete analyser | | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | 0.05 | 0.03 | 0.01 | |
| EN055: Ionic Balance | | | - | | | | |
| | | | | | | | |

Page : 5 of 8 Work Order : ES1508003 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB03 | RMB02 | RMB03 | |
|---------------------------------------|------------|------------|----------------|-------------------|-------------------|-------------------|------|
| | Clie | ent sampli | ng date / time | 07-APR-2015 11:00 | 07-APR-2015 15:00 | 07-APR-2015 16:30 | |
| Compound | CAS Number | LOR | Unit | ES1508003-001 | ES1508003-002 | ES1508003-003 | |
| EN055: Ionic Balance - Continued | | | | | | | |
| Total Anions | | 0.01 | meq/L | 55.0 | 110 | 74.7 | |
| Total Cations | | 0.01 | meq/L | 53.2 | 106 | 71.4 | |
| Ionic Balance | | 0.01 | % | 1.67 | 1.82 | 2.26 | |
| EP005: Total Organic Carbon (TOC) | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | 10 | 5 | 40 | |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 41700 | 39100 | 38100 | |
| Ethene | 74-85-1 | 10 | µg/L | <10 | <10 | <10 | |
| Ethane | 74-84-0 | 10 | µg/L | 108 | <10 | 11 | |
| Propene | 115-07-1 | 10 | µg/L | <10 | <10 | <10 | |
| Propane | 74-98-6 | 10 | µg/L | 19 | <10 | <10 | |
| Butene | 25167-67-3 | 10 | µg/L | <10 | <10 | <10 | |
| Butane | 106-97-8 | 10 | µg/L | <10 | <10 | <10 | |
| EP075(SIM)A: Phenolic Compounds | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2-Chlorophenol | 95-57-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | |
| EP075(SIM)B: Polynuclear Aromatic Hyd | rocarbons | | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Acenaphthene | 83-32-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |

Page : 6 of 8 Work Order : ES1508003 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | GLMB03 | RMB02 | RMB03 | |
|--|-------------------|------------|----------------|-------------------|-------------------|-------------------|------|
| | Cli | ent sampli | ng date / time | 07-APR-2015 11:00 | 07-APR-2015 15:00 | 07-APR-2015 16:30 | |
| Compound | CAS Number | LOR | Unit | ES1508003-001 | ES1508003-002 | ES1508003-003 | |
| EP075(SIM)B: Polynuclear Aromatic Hyd | Irocarbons - Cont | inued | | | | | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(b+j)fluoranthene 2 | 205-99-2 205-82-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | |
| Sum of polycyclic aromatic hydrocarbons | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | |
| [^] Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | |
| EP080/071: Total Petroleum Hydrocarbo | ns | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | 190 | <20 | 70 | |
| C10 - C14 Fraction | | 50 | µg/L | <50 | <50 | <50 | |
| C15 - C28 Fraction | | 100 | µg/L | <100 | <100 | <100 | |
| C29 - C36 Fraction | | 50 | µg/L | <50 | <50 | <50 | |
| [^] C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | <50 | <50 | |
| EP080/071: Total Recoverable Hydrocarl | bons - NEPM 201 | 3 Fractio | ns | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | 190 | <20 | 70 | |
| C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | 70 | <20 | 30 | |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | <100 | <100 | |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | <100 | <100 | |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | <100 | |
| ^ >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | <100 | <100 | |
| C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | <100 | <100 | |
| EP080: BTEXN | | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | <1 | |
| Toluene | 108-88-3 | 2 | μg/L | 124 | <2 | 39 | |
| Ethylbenzene | 100-41-4 | 2 | μg/L | <2 | <2 | <2 | |
| meta- & para-Xylene | 08-38-3 106-42-3 | 2 | μg/L | <2 | <2 | <2 | |
| ortho-Xylene | 95-47-6 | 2 | μg/L | <2 | <2 | <2 | |

Page : 7 of 8 Work Order : ES1508003 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | GLMB03 | RMB02 | RMB03 | |
|-----------------------------------|------------|-------------|-----------------|-------------------|-------------------|-------------------|------|
| | Cl | ient sampli | ing date / time | 07-APR-2015 11:00 | 07-APR-2015 15:00 | 07-APR-2015 16:30 | |
| Compound | CAS Number | LOR | Unit | ES1508003-001 | ES1508003-002 | ES1508003-003 | |
| EP080: BTEXN - Continued | | | | | | | |
| ^ Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | <2 | |
| [^] Sum of BTEX | | 1 | µg/L | 124 | <1 | 39 | |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | <5 | <5 | |
| EP075(SIM)S: Phenolic Compound S | Surrogates | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 18.5 | 22.3 | 18.9 | |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 41.0 | 49.9 | 18.0 | |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 40.5 | 59.0 | 22.7 | |
| EP075(SIM)T: PAH Surrogates | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 52.1 | 57.3 | 70.6 | |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 76.4 | 87.0 | 80.7 | |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 66.7 | 71.6 | 78.7 | |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 113 | 113 | 124 | |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 118 | 120 | 122 | |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 98.9 | 99.4 | 105 | |

(ALS)

Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) |
|---|------------|----------|------------|
| Compound | CAS Number | Low | High |
| EP075(SIM)S: Phenolic Compound Surrogates | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 |
| EP075(SIM)T: PAH Surrogates | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 |
| EP080S: TPH(V)/BTEX Surrogates | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 |
| Toluene-D8 | 2037-26-5 | 79 | 131 |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 |



| | CERTI | CERTIFICATE OF ANALYSIS | | | | | | | | | | |
|--------------|---------------------------------|-------------------------|---|--|--|--|--|--|--|--|--|--|
| Work Order | ES1508120 | Page | : 1 of 8 | | | | | | | | | |
| Client | : PARSONS BRINCKERHOFF AUST P/L | Laboratory | : Environmental Division Sydney | | | | | | | | | |
| Contact | : MS CAROLINA SARDELLA | Contact | : Client Services | | | | | | | | | |
| Address | : GPO BOX 5394 | Address | : 277-289 Woodpark Road Smithfield NSW Australia 2164 | | | | | | | | | |
| | SYDNEY NSW, AUSTRALIA 2001 | | | | | | | | | | | |
| E-mail | : csardella@pb.com.au | E-mail | : sydney@alsglobal.com | | | | | | | | | |
| Telephone | : +61 02 92725100 | Telephone | +61-2-8784 8555 | | | | | | | | | |
| Facsimile | : +61 02 92725101 | Facsimile | : +61-2-8784 8500 | | | | | | | | | |
| Project | : 2268518A | QC Level | : NEPM 2013 Schedule B(3) and ALS QCS3 requirement | | | | | | | | | |
| Order number | : | | | | | | | | | | | |
| C-O-C number | : | Date Samples Received | : 09-APR-2015 | | | | | | | | | |
| Sampler | : KM.AM | Issue Date | : 16-APR-2015 | | | | | | | | | |
| Site | : | | | | | | | | | | | |
| | | No. of samples received | : 5 | | | | | | | | | |
| Quote number | : SY/239/15 | No. of samples analysed | : 5 | | | | | | | | | |

This report supersedes any previous report(s) with this reference. Results apply to the sample(s) as submitted. All pages of this report have been checked and approved for release.

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results
- Surrogate Control Limits

Signatories NATA Accredited Laboratory 825 This document has been electronically signed by the authorized signatories indicated below. Electronic signing has been carried out in compliance with procedures specified in 21 CFR Part 11. Accredited for compliance with NATA ISO/IEC 17025. Signatories Position Accreditation Category Inorganic Chemist Ankit Joshi Sydney Inorganics Senior Spectroscopist Celine Conceicao Sydney Inorganics WORLD RECOGNISED ACCREDITATION Pabi Subba Senior Organic Chemist Sydney Organics Metals Coordinator Shobhna Chandra Sydney Inorganics

Address 277-289 Woodpark Road Smithfield NSW Australia 2164 PHONE +61-2-8784 8555 Facsimile +61-2-8784 8500 Environmental Division Svdnev ABN 84 009 936 029 Part of the ALS Group An ALS Limited Company



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General Comments

The analytical procedures used by the Environmental Division have been developed from established internationally recognized procedures such as those published by the USEPA, APHA, AS and NEPM. In house developed procedures are employed in the absence of documented standards or by client request.

Where moisture determination has been performed, results are reported on a dry weight basis.

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

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

When sampling time information is not provided by the client, sampling dates are shown without a time component. In these instances, the time component has been assumed by the laboratory for processing purposes.

Where a result is required to meet compliance limits the associated uncertainty must be considered. Refer to the ALS Contact for details.

Key : CAS Number = CAS registry number from database maintained by Chemical Abstracts Services. The Chemical Abstracts Service is a division of the American Chemical Society. LOR = Limit of reporting ^ = This result is computed from individual analyte detections at or above the level of reporting

- EG020: Bromine quantification may be unreliable due to its low solubility in acid, leading to variable volatility during measurement by ICPMS.
- EK026SF:Spike failed for Total Cyanide analysis due to matrix interference(confirmed by re analysis
- EK071G: It has been noted that Reactive P is greater than Total P for sample ID(MPMB02), however this difference is within the limits of experimental variation.



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB04 | MPMB03 | MPMB02 | MPMB01 | QA1 |
|--|-------------|--------------|----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | C | lient sampli | ng date / time | 08-APR-2015 10:00 | 08-APR-2015 09:00 | 08-APR-2015 10:15 | 08-APR-2015 11:15 | 08-APR-2015 15:00 |
| Compound | CAS Number | LOR | Unit | ES1508120-001 | ES1508120-002 | ES1508120-003 | ES1508120-004 | ES1508120-005 |
| EA005P: pH by PC Titrator | | | | | | | | |
| pH Value | | 0.01 | pH Unit | 9.73 | 7.44 | 6.86 | 5.54 | 6.90 |
| EA010P: Conductivity by PC Titrator | | | | | | | | |
| Electrical Conductivity @ 25°C | | 1 | μS/cm | 984 | 1090 | 963 | 912 | 950 |
| EA015: Total Dissolved Solids | | | | | | | | |
| Total Dissolved Solids @180°C | | 10 | mg/L | 540 | 528 | 420 | 526 | 508 |
| EA025: Suspended Solids | | | | | | | | |
| Suspended Solids (SS) | | 5 | mg/L | <5 | 6 | 5 | 74 | <5 |
| ED009: Anions | | | | | | | | |
| Bromide | 24959-67-9 | 0.010 | mg/L | 0.275 | 0.213 | 0.409 | 0.601 | 0.378 |
| ED037P: Alkalinity by PC Titrator | | | | | | | | |
| Hydroxide Alkalinity as CaCO3 | DMO-210-001 | 1 | mg/L | <1 | <1 | <1 | <1 | <1 |
| Carbonate Alkalinity as CaCO3 | 3812-32-6 | 1 | mg/L | 209 | <1 | <1 | <1 | <1 |
| Bicarbonate Alkalinity as CaCO3 | 71-52-3 | 1 | mg/L | 110 | 522 | 238 | 14 | 232 |
| Total Alkalinity as CaCO3 | | 1 | mg/L | 319 | 522 | 238 | 14 | 232 |
| ED041G: Sulfate (Turbidimetric) as SO4 | 4 2- bv DA | | | | | | | |
| Sulfate as SO4 - Turbidimetric | 14808-79-8 | 1 | mg/L | <1 | <1 | 4 | 2 | 4 |
| ED045G: Chloride Discrete analyser | | | | | | | | |
| Chloride | 16887-00-6 | 1 | mg/L | 126 | 77 | 170 | 258 | 171 |
| ED093F: Dissolved Major Cations | | | | | | | | |
| Calcium | 7440-70-2 | 1 | mg/L | 4 | 91 | 34 | 13 | 32 |
| Magnesium | 7439-95-4 | 1 | mg/L | 5 | 23 | 31 | 23 | 30 |
| Sodium | 7440-23-5 | 1 | mg/L | 189 | 110 | 105 | 110 | 105 |
| Potassium | 7440-09-7 | 1 | mg/L | 15 | 14 | 3 | 1 | 3 |
| EG020F: Dissolved Metals by ICP-MS | | | | | | | | |
| Aluminium | 7429-90-5 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | 0.01 | <0.01 |
| Antimony | 7440-36-0 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Arsenic | 7440-38-2 | 0.001 | mg/L | 0.002 | 0.018 | 0.002 | <0.001 | 0.002 |
| Boron | 7440-42-8 | 0.05 | mg/L | <0.05 | <0.05 | <0.05 | <0.05 | <0.05 |
| Barium | 7440-39-3 | 0.001 | mg/L | 0.753 | 3.27 | 0.584 | 0.664 | 0.571 |
| Beryllium | 7440-41-7 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Cadmium | 7440-43-9 | 0.0001 | mg/L | <0.0001 | <0.0001 | <0.0001 | <0.0001 | <0.0001 |
| Cobalt | 7440-48-4 | 0.001 | mg/L | <0.001 | 0.004 | <0.001 | 0.041 | <0.001 |
| Chromium | 7440-47-3 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |

Page : 4 of 8 Work Order : ES1508120 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | MPMB04 | МРМВ03 | MPMB02 | MPMB01 | QA1 |
|--|---------------------|-------------|-----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Cl | ient sampli | ing date / time | 08-APR-2015 10:00 | 08-APR-2015 09:00 | 08-APR-2015 10:15 | 08-APR-2015 11:15 | 08-APR-2015 15:00 |
| Compound | CAS Number | LOR | Unit | ES1508120-001 | ES1508120-002 | ES1508120-003 | ES1508120-004 | ES1508120-005 |
| EG020F: Dissolved Metals by ICP-MS | Continued | | | | | | | |
| Copper | 7440-50-8 | 0.001 | mg/L | 0.001 | <0.001 | <0.001 | 0.003 | <0.001 |
| Manganese | 7439-96-5 | 0.001 | mg/L | <0.001 | 0.047 | 0.144 | 0.488 | 0.139 |
| Nickel | 7440-02-0 | 0.001 | mg/L | <0.001 | 0.004 | <0.001 | 0.016 | <0.001 |
| Lead | 7439-92-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | 0.002 | <0.001 |
| Selenium | 7782-49-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Vanadium | 7440-62-2 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| Zinc | 7440-66-6 | 0.005 | mg/L | 0.047 | 0.012 | 0.018 | 0.066 | 0.015 |
| Molybdenum | 7439-98-7 | 0.001 | mg/L | 0.005 | <0.001 | <0.001 | <0.001 | <0.001 |
| Strontium | 7440-24-6 | 0.001 | mg/L | 0.198 | 0.887 | 0.352 | 0.143 | 0.350 |
| Uranium | 7440-61-1 | 0.001 | mg/L | <0.001 | <0.001 | <0.001 | <0.001 | <0.001 |
| Iron | 7439-89-6 | 0.05 | mg/L | <0.05 | 1.13 | 4.21 | 0.06 | 4.14 |
| Bromine | 7726-95-6 | 0.1 | mg/L | 0.3 | 0.2 | 0.4 | 0.7 | 0.4 |
| EG035F: Dissolved Mercury by FIMS | | | | | | | | |
| Mercury | 7439-97-6 | 0.0001 | mg/L | 0.0003 | <0.0001 | <0.0001 | <0.0001 | <0.0001 |
| G052G: Silica by Discrete Analyser | | | | | | | | |
| Reactive Silica | | 0.05 | mg/L | 3.48 | 8.82 | 12.7 | 19.2 | 12.9 |
| EK026SF: Total CN by Segmented Flo | w Analyser | | | | | | | |
| Total Cyanide | 57-12-5 | 0.004 | mg/L | <0.004 | <0.004 | <0.004 | <0.004 | <0.004 |
| EK040P: Fluoride by PC Titrator | | | | | | | | |
| Fluoride | 16984-48-8 | 0.1 | mg/L | 0.6 | 0.2 | 0.2 | <0.1 | 0.3 |
| K055G: Ammonia as N by Discrete A | nalyser | | | | | | | |
| Ammonia as N | 7664-41-7 | 0.01 | mg/L | 0.86 | 0.92 | 0.06 | 0.02 | 0.09 |
| EK057G: Nitrite as N by Discrete Anal | yser | | | | | | | |
| Nitrite as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | <0.01 | <0.01 |
| EK058G: Nitrate as N by Discrete Ana | lyser | | | | | | | |
| Nitrate as N | 14797-55-8 | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | 0.16 | <0.01 |
| EK059G: Nitrite plus Nitrate as N (NO) | x) by Discrete Ana | lvser | | | | | | |
| Nitrite + Nitrate as N | | 0.01 | mg/L | <0.01 | <0.01 | <0.01 | 0.16 | <0.01 |
| EK067G: Total Phosphorus as P by Di | screte Analyser | | | | | | | |
| Total Phosphorus as P | | 0.01 | mg/L | <0.01 | <0.01 | 0.02 | 0.05 | 0.01 |
| EK071G: Reactive Phosphorus as P by | v discrete analyser | | | | | | | |
| Reactive Phosphorus as P | 14265-44-2 | 0.01 | mg/L | 0.01 | <0.01 | 0.03 | <0.01 | <0.01 |
| EN055: Ionic Balance | | | | | | | | |

Page : 5 of 8 Work Order : ES1508120 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB04 | MPMB03 | MPMB02 | MPMB01 | QA1 |
|--------------------------------------|-------------|-------------|----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Cl | ient sampli | ng date / time | 08-APR-2015 10:00 | 08-APR-2015 09:00 | 08-APR-2015 10:15 | 08-APR-2015 11:15 | 08-APR-2015 15:00 |
| Compound | CAS Number | LOR | Unit | ES1508120-001 | ES1508120-002 | ES1508120-003 | ES1508120-004 | ES1508120-005 |
| EN055: Ionic Balance - Continued | | | | | | | | |
| Total Anions | | 0.01 | meq/L | 9.93 | 12.6 | 9.63 | 7.60 | 9.54 |
| Total Cations | | 0.01 | meq/L | 9.22 | 11.6 | 8.89 | 7.35 | 8.71 |
| Ionic Balance | | 0.01 | % | 3.76 | 4.25 | 4.02 | 1.66 | 4.58 |
| EP005: Total Organic Carbon (TOC) | | | | | | | | |
| Total Organic Carbon | | 1 | mg/L | 22 | 4 | <1 | <1 | <1 |
| EP033: C1 - C4 Hydrocarbon Gases | | | | | | | | |
| Methane | 74-82-8 | 10 | µg/L | 52300 | 44600 | 273 | <10 | 266 |
| Ethene | 74-85-1 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Ethane | 74-84-0 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Propene | 115-07-1 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Propane | 74-98-6 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Butene | 25167-67-3 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| Butane | 106-97-8 | 10 | µg/L | <10 | <10 | <10 | <10 | <10 |
| EP075(SIM)A: Phenolic Compounds | | | | | | | | |
| Phenol | 108-95-2 | 1.0 | µg/L | 2.8 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2-Chlorophenol | 95-57-8 | 1.0 | μg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2-Methylphenol | 95-48-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 3- & 4-Methylphenol | 1319-77-3 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | <2.0 |
| 2-Nitrophenol | 88-75-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4-Dimethylphenol | 105-67-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4-Dichlorophenol | 120-83-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.6-Dichlorophenol | 87-65-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 4-Chloro-3-methylphenol | 59-50-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4.6-Trichlorophenol | 88-06-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| 2.4.5-Trichlorophenol | 95-95-4 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Pentachlorophenol | 87-86-5 | 2.0 | µg/L | <2.0 | <2.0 | <2.0 | <2.0 | <2.0 |
| EP075(SIM)B: Polynuclear Aromatic Hy | /drocarbons | | | | | | | |
| Naphthalene | 91-20-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | 1.1 |
| Acenaphthylene | 208-96-8 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Acenaphthene | 83-32-9 | 1.0 | μg/L | <1.0 | <1.0 | 1.9 | 1.1 | 2.1 |
| Fluorene | 86-73-7 | 1.0 | µg/L | <1.0 | <1.0 | 1.7 | 1.3 | 2.3 |
| Phenanthrene | 85-01-8 | 1.0 | µg/L | <1.0 | <1.0 | 4.7 | 1.8 | 3.5 |
| Anthracene | 120-12-7 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |

Page : 6 of 8 Work Order : ES1508120 Client : PARSONS BRINCKERHOFF AUST P/L Project : 2268518A



| Sub-Matrix: WATER (Matrix: WATER) | | Clie | ent sample ID | MPMB04 | MPMB03 | МРМВ02 | MPMB01 | QA1 |
|---|----------------------|------------|----------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| | Cli | ent sampli | ng date / time | 08-APR-2015 10:00 | 08-APR-2015 09:00 | 08-APR-2015 10:15 | 08-APR-2015 11:15 | 08-APR-2015 15:00 |
| Compound | CAS Number | LOR | Unit | ES1508120-001 | ES1508120-002 | ES1508120-003 | ES1508120-004 | ES1508120-005 |
| EP075(SIM)B: Polynuclear Aromatic | Hydrocarbons - Conti | inued | | | | | | |
| Fluoranthene | 206-44-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Pyrene | 129-00-0 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benz(a)anthracene | 56-55-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Chrysene | 218-01-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(b+j)fluoranthene | 205-99-2 205-82-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(k)fluoranthene | 207-08-9 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(a)pyrene | 50-32-8 | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 |
| Indeno(1.2.3.cd)pyrene | 193-39-5 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Dibenz(a.h)anthracene | 53-70-3 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Benzo(g.h.i)perylene | 191-24-2 | 1.0 | µg/L | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Sum of polycyclic aromatic hydrocarbon | IS | 0.5 | µg/L | <0.5 | <0.5 | 8.3 | 4.2 | 9.0 |
| Benzo(a)pyrene TEQ (zero) | | 0.5 | µg/L | <0.5 | <0.5 | <0.5 | <0.5 | <0.5 |
| EP080/071: Total Petroleum Hydroca | rbons | | | | | | | |
| C6 - C9 Fraction | | 20 | µg/L | 110 | <20 | <20 | <20 | <20 |
| C10 - C14 Fraction | | 50 | µg/L | <50 | <50 | 70 | <50 | 80 |
| C15 - C28 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| C29 - C36 Fraction | | 50 | µg/L | <50 | <50 | <50 | <50 | <50 |
| C10 - C36 Fraction (sum) | | 50 | µg/L | <50 | <50 | 70 | <50 | 80 |
| EP080/071: Total Recoverable Hydro | carbons - NEPM 201 | 3 Fractio | ns | | | | | |
| C6 - C10 Fraction | C6_C10 | 20 | µg/L | 110 | <20 | <20 | <20 | <20 |
| C6 - C10 Fraction minus BTEX (F1) | C6_C10-BTEX | 20 | µg/L | 40 | <20 | <20 | <20 | <20 |
| >C10 - C16 Fraction | >C10_C16 | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| >C16 - C34 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| >C34 - C40 Fraction | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| >C10 - C40 Fraction (sum) | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| >C10 - C16 Fraction minus Naphthalene (F2) | | 100 | µg/L | <100 | <100 | <100 | <100 | <100 |
| EP080: BTEXN | | | | | | | | |
| Benzene | 71-43-2 | 1 | µg/L | <1 | <1 | <1 | <1 | <1 |
| Toluene | 108-88-3 | 2 | µg/L | 66 | <2 | <2 | <2 | <2 |
| Ethylbenzene | 100-41-4 | 2 | µg/L | <2 | <2 | <2 | <2 | <2 |
| meta- & para-Xylene | 108-38-3 106-42-3 | 2 | µg/L | <2 | <2 | <2 | <2 | <2 |
| ortho-Xylene | 95-47-6 | 2 | µg/L | <2 | <2 | <2 | <2 | <2 |

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| Sub-Matrix: WATER (Matrix: WATER) | | Cli | ent sample ID | MPMB04 | MPMB03 | MPMB02 | MPMB01 | QA1 |
|-----------------------------------|------------|-----------------------------|---------------|---------------|-------------------|-------------------|-------------------|-------------------|
| | Cli | Client sampling date / time | | | 08-APR-2015 09:00 | 08-APR-2015 10:15 | 08-APR-2015 11:15 | 08-APR-2015 15:00 |
| Compound | CAS Number | LOR | Unit | ES1508120-001 | ES1508120-002 | ES1508120-003 | ES1508120-004 | ES1508120-005 |
| EP080: BTEXN - Continued | | | | | | | | |
| Total Xylenes | 1330-20-7 | 2 | µg/L | <2 | <2 | <2 | <2 | <2 |
| Sum of BTEX | | 1 | µg/L | 66 | <1 | <1 | <1 | <1 |
| Naphthalene | 91-20-3 | 5 | µg/L | <5 | <5 | <5 | <5 | <5 |
| EP075(SIM)S: Phenolic Compound Su | ırrogates | | | | | | | |
| Phenol-d6 | 13127-88-3 | 0.1 | % | 13.6 | 21.8 | 24.3 | 18.5 | 27.2 |
| 2-Chlorophenol-D4 | 93951-73-6 | 0.1 | % | 26.5 | 43.6 | 49.9 | 47.5 | 62.7 |
| 2.4.6-Tribromophenol | 118-79-6 | 0.1 | % | 29.5 | 38.4 | 48.0 | 57.8 | 84.0 |
| EP075(SIM)T: PAH Surrogates | | | | | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 0.1 | % | 45.3 | 48.9 | 55.9 | 47.7 | 66.0 |
| Anthracene-d10 | 1719-06-8 | 0.1 | % | 90.9 | 92.1 | 90.9 | 43.3 | 67.2 |
| 4-Terphenyl-d14 | 1718-51-0 | 0.1 | % | 60.5 | 61.9 | 72.0 | 42.0 | 66.9 |
| EP080S: TPH(V)/BTEX Surrogates | | | | | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 0.1 | % | 79.0 | 81.1 | 80.3 | 83.0 | 76.0 |
| Toluene-D8 | 2037-26-5 | 0.1 | % | 96.2 | 93.4 | 95.0 | 94.6 | 89.0 |
| 4-Bromofluorobenzene | 460-00-4 | 0.1 | % | 93.1 | 91.1 | 90.9 | 90.6 | 84.9 |

ALS

Surrogate Control Limits

| Sub-Matrix: WATER | | Recovery | Limits (%) | |
|---|------------|----------|------------|--|
| Compound | CAS Number | Low | High | |
| EP075(SIM)S: Phenolic Compound Surrogates | | | | |
| Phenol-d6 | 13127-88-3 | 10.0 | 44 | |
| 2-Chlorophenol-D4 | 93951-73-6 | 14 | 94 | |
| 2.4.6-Tribromophenol | 118-79-6 | 17 | 125 | |
| EP075(SIM)T: PAH Surrogates | | | | |
| 2-Fluorobiphenyl | 321-60-8 | 20 | 104 | |
| Anthracene-d10 | 1719-06-8 | 27.4 | 113 | |
| 4-Terphenyl-d14 | 1718-51-0 | 32 | 112 | |
| EP080S: TPH(V)/BTEX Surrogates | | | | |
| 1.2-Dichloroethane-D4 | 17060-07-0 | 71 | 137 | |
| Toluene-D8 | 2037-26-5 | 79 | 131 | |
| 4-Bromofluorobenzene | 460-00-4 | 70 | 128 | |