

CERTIFICATE OF ANALYSIS

Work Order	ES1623644	Page	: 1 of 8
Client	: ORIGIN ENERGY RESOURCES LTD	Laboratory	: Environmental Division Sydney
Contact		Contact	
Address	ORIGIN ENERGY RESOURCES LIMITED GPO BOX 148	Address	: 277-289 Woodpark Road Smithfield NSW Australia 2164
	BRISBANE QLD, AUSTRALIA 4001		
Telephone		Telephone	
Project	BEETALOO GROUNDWATER MONITORING	Date Samples Received	: 21-Oct-2016 13:45
Order number	: 16231417	Date Analysis Commenced	: 21-Oct-2016
C-O-C number	: 19102016_SU_BEETALOO FB	Issue Date	: 02-Nov-2016 16:43
Sampler	:		Iac-MRA NATA
Site	: BEETALOO		
Quote number	:		Accreditation No. 825
No. of samples received	: 3		Accredited for compliance with
No. of samples analysed	: 3		ISO/IEC 17025 - Testing

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

This Certificate of Analysis contains the following information:

- General Comments
- Analytical Results
- Surrogate Control Limits

Additional information pertinent to this report will be found in the following separate attachments: Quality Control Report, QA/QC Compliance Assessment to assist with Quality Review and Sample Receipt Notification.

Signatories

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

Signatories	Position	Accreditation Category
	Inorganic Chemist	Sydney Inorganics, Smithfield, NSW
	Inorganic Chemist	Sydney Inorganics, Smithfield, NSW
	Senior Spectroscopist	Sydney Inorganics, Smithfield, NSW
		Sydney Inorganics, Smithfield, NSW
	Organic Coordinator	Sydney Organics, Smithfield, NSW
	Lab technician	Sydney Inorganics, Smithfield, NSW
i	Metals Teamleader	Radionuclides, Fyshwick, ACT
	Inorganics Coordinator	Sydney Inorganics, Smithfield, NSW



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
 - ø = ALS is not NATA accredited for these tests.
 - ~ = Indicates an estimated value.
- EG035: Poor matrix spike recovery was obtained for Mercury on sample ES1623644# 2 due to high matrix interference. Confirmed by re-analysis.
- EG020/ED093: LOR's have been raised for sample ID ES1623644 #001, 2 & 3 due to matrix interference. (High Total Dissolved Solids)
- ED041G: LOR raised for (Sulfate analysis) on sample no:1, due to sample matrix.
- Gross Alpha and Beta Activity analyses are performed by ALS Fyshwick (NATA Accreditation number 992).
- EP050: The MBAS reported is calculated as LAS, mol wt ____342___
- EA016: Calculated TDS is determined from Electrical conductivity using a conversion factor of 0.65.
- EA250 LSC : LOR for Gross Alpha and Beta raised due to high solid content
- Benzo(a)pyrene Toxicity Equivalent Quotient (TEQ) is the sum total of the concentra ion 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), 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.



Sub-Matrix: WATER (Matrix: WATER)		Clie	ent sample ID	BET-PW001_Fe_11.5%	BET-PW001_Fe_12.15 %	BET-PW001_Fe_12.5%	
	Ci	lient sampli	ng date / time	15-Oct-2016 19:00	17-Oct-2016 09:00	19-Oct-2016 08:00	
Compound	CAS Number	LOR	Unit	ES1623644-001	ES1623644-002	ES1623644-003	
				Result	Result	Result	
EA005P: pH by PC Titrator							
pH Value		0.01	pH Unit	6.47	6.50	6.45	
EA010P: Conductivity by PC Titrator							
Electrical Conductivity @ 25°C		1	µS/cm	49000	50500	51100	
EA015: Total Dissolved Solids dried at 1	80 ± 5 °C						
Total Dissolved Solids @180°C		10	mg/L	38800	37400	39000	
EA016: Calculated TDS (from Electrical 0							
Total Dissolved Solids (Calc.)		1	mg/L	31800	32800	33200	
EA065: Total Hardness as CaCO3							
Total Hardness as CaCO3		1	mg/L	4090	4110	4500	
		1	ing/L	4050	4110	4500	
EA250: Gross Alpha and Beta Activity		0.05	D-4	0.00	0.01		
Gross alpha		0.05	Bq/L	8.82	6.31	8.38	
Gross beta activity - 40K		0.1	Bq/L	15.4	7.55	8.31	
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	<1	<1	
Bicarbonate Alkalinity as CaCO3	71-52-3	1	mg/L	506	474	472	
Total Alkalinity as CaCO3		1	mg/L	506	474	472	
ED041G: Sulfate (Turbidimetric) as SO4 2	2- by DA						
Sulfate as SO4 - Turbidimetric	14808-79-8	1	mg/L	<10	26	38	
ED045G: Chloride by Discrete Analyser							
Chloride	16887-00-6	1	mg/L	18200	18300	18800	
ED093F: Dissolved Major Cations							
Calcium	7440-70-2	1	mg/L	1220	1230	1360	
Magnesium	7439-95-4	1	mg/L	253	252	269	
Sodium	7440-23-5	1	mg/L	12100	11700	12300	
Potassium	7440-09-7	1	mg/L	72	70	74	
EG020F: Dissolved Metals by ICP-MS							
Arsenic	7440-38-2	0.001	mg/L	<0.010	<0.010	<0.010	
Boron	7440-42-8	0.05	mg/L	45.9	44.5	43.5	
Barium	7440-39-3	0.001	mg/L	51.9	53.9	59.1	
Beryllium	7440-41-7		mg/L	<0.010	<0.010	<0.010	
Cadmium	7440-43-9		mg/L	<0.0010	<0.0010	<0.0010	
Cobalt	7440-48-4	0.001	mg/L	<0.010	<0.010	<0.010	
	1440-40-4						



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	Cl	ient samplii	ng date / time	15-Oct-2016 19:00	17-Oct-2016 09:00	19-Oct-2016 08:00		
Compound	CAS Number	LOR	Unit	ES1623644-001	ES1623644-002	ES1623644-003		
				Result	Result	Result	—	
EG020F: Dissolved Metals by ICP-M	S - Continued							
Chromium	7440-47-3	0.001	mg/L	0.042	0.030	0.033		
Copper	7440-50-8	0.001	mg/L	<0.010	<0.010	<0.010		
Manganese	7439-96-5	0.001	mg/L	2.38	2.24	2.43		
Nickel	7440-02-0	0.001	mg/L	<0.010	<0.010	<0.010		
Lead	7439-92-1	0.001	mg/L	<0.010	<0.010	<0.010		
Selenium	7782-49-2	0.01	mg/L	<0.10	<0.10	<0.10		
Vanadium	7440-62-2	0.01	mg/L	<0.10	<0.10	<0.10		
Zinc	7440-66-6	0.005	mg/L	<0.050	<0.050	<0.050		
EG035F: Dissolved Mercury by FIMS								
Mercury	7439-97-6	0.0001	mg/L	<0.0001	<0.0001	<0.0001		
EK040P: Fluoride by PC Titrator								
Fluoride	16984-48-8	0.1	mg/L	1.1	1.1	1.1		
EK059G: Nitrite plus Nitrate as N (N	Ox) by Discrete Ana	lvser						
Nitrite + Nitrate as N		0.01	mg/L	0.03	0.02	0.12		
EK061G: Total Kjeldahl Nitrogen By	Discrete Analyser							
Total Kjeldahl Nitrogen as N		0.1	mg/L	45.1	65.6	48.0		
EK062G: Total Nitrogen as N (TKN +	NOx) by Discrete Ar	nalvser	_					
Total Nitrogen as N		0.1	mg/L	45.1	65.6	48.1		
EK067G: Total Phosphorus as P by I	Discrete Analyser							
Total Phosphorus as P		0.01	mg/L	0.22	0.16	0.12		
EN055: Ionic Balance								
Total Anions		0.01	meg/L	524	526	540		
Total Cations		0.01	meq/L	610	593	627		
Ionic Balance		0.01	%	7.62	5.95	7.40		
EP033: C1 - C4 Hydrocarbon Gases								
Methane	74-82-8	1	µg/L	5460	7090	5500		
Ethene	74-85-1	1	µg/L	<1	<1	<1		
Ethane	74-84-0	1	µg/L	581	657	575		
Propene	115-07-1	1	µg/L	<1	<1	<1		
Propane	74-98-6	1	µg/L	16	17	13		
Butene	25167-67-3	1	µg/L	<1	<1	<1		
Dutene								-



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Compound	CAS Number	LOR	Unit	ES1623644-001	ES1623644-002	ES1623644-003	
				Result	Result	Result	 -
EP041A: Nonionic Surfactants - Contin	ued						
Nonionic Surfactants as CTAS		5	mg/L	<5	<5	<5	
EP050: Anionic Surfactants as MBAS							
Anionic Surfactants as MBAS		0.1	mg/L	<0.1	0.1	<0.1	
EP075(SIM)A: Phenolic Compounds							
Phenol	108-95-2	1	µg/L	2.3	2.2	2.2	
2-Chlorophenol	95-57-8	1	µg/L	<1.0	<1.0	<1.0	
2-Methylphenol	95-48-7	1	µg/L	<1.0	<1.0	<1.0	
3- & 4-Methylphenol	1319-77-3	2	µg/L	2.6	2.1	2.9	
2-Nitrophenol	88-75-5	1	µg/L	<1.0	<1.0	<1.0	
2.4-Dimethylphenol	105-67-9	1	µg/L	<1.0	<1.0	<1.0	
2.4-Dichlorophenol	120-83-2	1	µg/L	<1.0	<1.0	<1.0	
2.6-Dichlorophenol	87-65-0	1	µg/L	<1.0	<1.0	<1.0	
4-Chloro-3-methylphenol	59-50-7	1	µg/L	<1.0	<1.0	<1.0	
2.4.6-Trichlorophenol	88-06-2	1	µg/L	<1.0	<1.0	<1.0	
2.4.5-Trichlorophenol	95-95-4	1	µg/L	<1.0	<1.0	<1.0	
Pentachlorophenol	87-86-5	2	µg/L	<2.0	<2.0	<2.0	
EP075(SIM)B: Polynuclear Aromatic H	ydrocarbons						
Naphthalene	91-20-3	1	µg/L	<1.0	<1.0	<1.0	
Acenaphthylene	208-96-8	1	µg/L	<1.0	<1.0	<1.0	
Acenaphthene	83-32-9	1	µg/L	<1.0	<1.0	<1.0	
Fluorene	86-73-7	1	µg/L	<1.0	<1.0	<1.0	
Phenanthrene	85-01-8	1	µg/L	<1.0	<1.0	<1.0	
Anthracene	120-12-7	1	µg/L	<1.0	<1.0	<1.0	
Fluoranthene	206-44-0	1	µg/L	<1.0	<1.0	<1.0	
Pyrene	129-00-0	1	µg/L	<1.0	<1.0	<1.0	
Benz(a)anthracene	56-55-3	1	µg/L	<1.0	<1.0	<1.0	
Chrysene	218-01-9	1	µg/L	<1.0	<1.0	<1.0	
Benzo(b+j)fluoranthene	205-99-2 205-82-3	1	µg/L	<1.0	<1.0	<1.0	
Benzo(k)fluoranthene	207-08-9	1	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	µg/L	<1.0	<1.0	<1.0	
Dibenz(a.h)anthracene	53-70-3	1	µg/L	<1.0	<1.0	<1.0	
Benzo(g.h.i)perylene	191-24-2	1	µg/L	<1.0	<1.0	<1.0	
^ Sum of polycyclic aromatic hydrocarbon	s	0.5	µg/L	<0.5	<0.5	<0.5	



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Compound	CAS Number	LOR	Unit	ES1623644-001	ES1623644-002	ES1623644-003	
				Result	Result	Result	
EP075(SIM)B: Polynuclear Aromatic Hy	drocarbons - Cont	tinued					
^ Benzo(a)pyrene TEQ (zero)		0.5	µg/L	<0.5	<0.5	<0.5	
EP080/071: Total Petroleum Hydrocarb	ons						
C6 - C9 Fraction		20	µg/L	60	80	80	
C10 - C14 Fraction		50	µg/L	110	160	240	
C15 - C28 Fraction		100	µg/L	470	<100	100	
C29 - C36 Fraction		50	µg/L	200	<50	<50	
^ C10 - C36 Fraction (sum)		50	µg/L	780	160	340	
EP080/071: Total Recoverable Hydroca	rbons - NEPM 201	3 Fraction	ns				
C6 - C10 Fraction	C6_C10	20	µg/L	60	80	80	
^ C6 - C10 Fraction minus BTEX	C6_C10-BTEX	20	µg/L	50	70	70	
(F1)	-						
>C10 - C16 Fraction		100	µg/L	<100	<100	120	
>C16 - C34 Fraction		100	µg/L	600	110	110	
>C34 - C40 Fraction		100	µg/L	<100	<100	<100	
^ >C10 - C40 Fraction (sum)		100	µg/L	600	110	230	
^ >C10 - C16 Fraction minus Naphthalene		100	µg/L	<100	<100	120	
(F2)							
EP080: BTEXN							
Benzene	71-43-2	1	µg/L	3	4	4	
Toluene	108-88-3	2	µg/L	2	3	3	
Ethylbenzene	100-41-4	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	
^ Total Xylenes	1330-20-7	2	µg/L	2	<2	<2	
^ Sum of BTEX		1	µg/L	7	7	7	
Naphthalene	91-20-3	5	µg/L	<5	<5	<5	
EP075(SIM)S: Phenolic Compound Sur	rogates						
Phenol-d6	13127-88-3	1	%	22.5	21.6	20.5	
2-Chlorophenol-D4	93951-73-6	1	%	53.6	50.3	50.3	
2.4.6-Tribromophenol	118-79-6	1	%	89.7	82.4	81.5	
EP075(SIM)T: PAH Surrogates							
2-Fluorobiphenyl	321-60-8	1	%	63.1	58.3	53.7	
Anthracene-d10	1719-06-8	1	%	81.1	74.5	72.6	
4-Terphenyl-d14	1718-51-0	1	%	105	102	97.9	



Sub-Matrix: WATER (Matrix: WATER)		Clie	ent sample ID	BET-PW001_Fe_11.5%	BET-PW001_Fe_12.15	BET-PW001_Fe_12.5%	
					%		
	Cli	ient sampli	ng date / time	15-Oct-2016 19:00	17-Oct-2016 09:00	19-Oct-2016 08:00	
Compound	CAS Number	LOR	Unit	ES1623644-001	ES1623644-002	ES1623644-003	
				Result	Result	Result	
EP080S: TPH(V)/BTEX Surrogates							
1.2-Dichloroethane-D4	17060-07-0	2	%	125	124	122	
Toluene-D8	2037-26-5	2	%	107	108	101	
4-Bromofluorobenzene	460-00-4	2	%	114	112	109	



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	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	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