

5B  
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

Instrument ID: NT8

Project: GAS WORKS PARK-PLAY AREA

DFTPP Injection Date: 01/13/15

DFTPP Injection Time: 1833

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	18.6
68	Less than 2.0% of mass 69	0.0 ( 0.0)1
69	Mass 69 relative abundance	48.6
70	Less than 2.0% of mass 69	0.3 ( 0.7)1
127	10.0 - 80.0% of mass 198	48.3
197	Less than 2.0% of mass 198	0.0
198	Base Peak, 100% relative abundance	100.0
199	5.0 to 9.0% of mass 198	7.7
275	10.0 - 60.0% of mass 198	31.4
365	Greater than 1.0% of mass 198	3.71
441	0.0 - 24.0% of mass 442	9.0 ( 16.2)2
442	50.0 - 200.0% of mass 198	55.7
443	15.0 - 24.0% of mass 442	12.0 ( 21.6)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	ICV0113	ICV0113	01131519	01/13/15	1846
02	ZP06LCSS1	ZP06LCSS1	01131524	01/13/15	2053
03	ZP06LCSDS1	ZP06LCSDS1	01131525	01/13/15	2118
04	PAI-9-12.5-13.0	ZP16A	01131530	01/13/15	2325
05	PAI-9-12.5-13-DU	ZP16B	01131531	01/13/15	2350
06	PAI-10-9.5-10.0	ZP16C	01131532	01/14/15	0016
07	PAI-10-19.5-20.0	ZP16D	01131533	01/14/15	0041
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5B  
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

Instrument ID: NT8

Project: GAS WORKS PARK-PLAY AREA

DFTPP Injection Date: 01/14/15

DFTPP Injection Time: 1044

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	17.9
68	Less than 2.0% of mass 69	0.0 ( 0.0)1
69	Mass 69 relative abundance	51.9
70	Less than 2.0% of mass 69	0.4 ( 0.7)1
127	10.0 - 80.0% of mass 198	48.8
197	Less than 2.0% of mass 198	0.3
198	Base Peak, 100% relative abundance	100.0
199	5.0 to 9.0% of mass 198	7.6
275	10.0 - 60.0% of mass 198	31.3
365	Greater than 1.0% of mass 198	3.42
441	0.0 - 24.0% of mass 442	9.3 ( 15.9)2
442	50.0 - 200.0% of mass 198	58.5
443	15.0 - 24.0% of mass 442	11.7 ( 20.1)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	ICV0114	ICV0114	01141502	01/14/15	1057
02	ZP06MBS1	ZP06MBS1	01141503	01/14/15	1123
03	PAI-10-24.5-25.0	ZP16E	01141506	01/14/15	1238
04	PAI-9-12.5-13.0	ZP16A	01141507	01/14/15	1304
05	PAI-9-12.5-13-DU	ZP16B	01141508	01/14/15	1329
06	PAI-10-9.5-10.0	ZP16C	01141509	01/14/15	1354
07	PAI-10-19.5-20.0	ZP16D	01141510	01/14/15	1420
08	PAI-10-24.5-25.0	ZP16E	01141511	01/14/15	1445
09	PAI-9-12.5-13.0	ZP16A	01141512	01/14/15	1510
10	PAI-10-9.5-10.0	ZP16C	01141513	01/14/15	1535
11	PAI-10-19.5-20.0	ZP16D	01141514	01/14/15	1601
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## SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT8

Cont. Calib. Date: 01/13/15

Init. Calib. Date: 01/05/15

Cont. Calib. Time: 1846

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
=====	=====	=====	=====	=====	=====
Naphthalene	1.150	1.113	0.700	AVRG	-3.2
2-Methylnaphthalene	0.673	0.702	0.400	AVRG	4.3
Acenaphthylene	1.724	1.772	0.900	AVRG	2.8
Acenaphthene	1.171	1.177	0.900	AVRG	0.5
Dibenzofuran	1.645	1.576	0.800	AVRG	-4.2
Fluorene	1.333	1.339	0.900	AVRG	0.4
Phenanthrene	1.088	1.104	0.700	AVRG	1.5
Anthracene	1.078	1.115	0.700	AVRG	3.4
Fluoranthene	1.316	1.281	0.600	AVRG	-2.6
Pyrene	1.231	1.262	0.600	AVRG	2.5
Benzo(a)anthracene	1.218	1.209	0.800	AVRG	-0.7
Chrysene	1.182	1.148	0.700	AVRG	-2.9
Benzo(b)fluoranthene	1.133	1.166	0.700	AVRG	2.9
Benzo(k)fluoranthene	1.178	1.150	0.700	AVRG	-2.4
Benzo(j)fluoranthene	1.160	1.090	0.010	AVRG	-6.0
Benzo(a)pyrene	1.116	1.099	0.700	AVRG	-1.5
Indeno(1,2,3-cd)pyrene	1.241	1.348	0.500	AVRG	8.6
Dibenzo(a,h)anthracene	1.024	1.134	0.400	AVRG	10.7
Benzo(g,h,i)perylene	1.073	1.166	0.500	AVRG	8.7
1-methylnaphthalene	0.622	0.663	0.010	AVRG	6.6
Perylene	1.122	1.096	0.010	AVRG	-2.3
=====	=====	=====	=====	=====	=====
2-Methylnaphthalene-d10	0.707	0.711	0.010	AVRG	0.6
Dibenzo(a,h)anthracene-d14	0.820	0.959	0.010	AVRG	17.0
Fluoranthene-d10	1.185	1.179	0.010	AVRG	-0.5

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF



## SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT8

Cont. Calib. Date: 01/14/15

Init. Calib. Date: 01/05/15

Cont. Calib. Time: 1057

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
=====	=====	=====	=====	=====	=====
Naphthalene	1.150	1.077	0.700	AVRG	-6.3
2-Methylnaphthalene	0.673	0.697	0.400	AVRG	3.6
Acenaphthylene	1.724	1.820	0.900	AVRG	5.6
Acenaphthene	1.171	1.195	0.900	AVRG	2.0
Dibenzofuran	1.645	1.642	0.800	AVRG	-0.2
Fluorene	1.333	1.370	0.900	AVRG	2.8
Phenanthrene	1.088	1.086	0.700	AVRG	-0.2
Anthracene	1.078	1.099	0.700	AVRG	1.9
Fluoranthene	1.316	1.297	0.600	AVRG	-1.4
Pyrene	1.231	1.231	0.600	AVRG	0.0
Benzo(a)anthracene	1.218	1.210	0.800	AVRG	-0.6
Chrysene	1.182	1.141	0.700	AVRG	-3.5
Benzo(b)fluoranthene	1.133	1.174	0.700	AVRG	3.6
Benzo(k)fluoranthene	1.178	1.196	0.700	AVRG	1.5
Benzo(j)fluoranthene	1.160	1.131	0.010	AVRG	-2.5
Benzo(a)pyrene	1.116	1.106	0.700	AVRG	-0.9
Indeno(1,2,3-cd)pyrene	1.241	1.246	0.500	AVRG	0.4
Dibenzo(a,h)anthracene	1.024	1.045	0.400	AVRG	2.0
Benzo(g,h,i)perylene	1.073	1.077	0.500	AVRG	0.4
1-methylnaphthalene	0.622	0.675	0.010	AVRG	8.5
Perylene	1.122	1.108	0.010	AVRG	-1.2
=====	=====	=====	=====	=====	=====
2-Methylnaphthalene-d10	0.707	0.697	0.010	AVRG	-1.4
Dibenzo(a,h)anthracene-d14	0.820	0.893	0.010	AVRG	8.9
Fluoranthene-d10	1.185	1.194	0.010	AVRG	0.8

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF

## SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT8

Cont. Calib. Date: 01/15/15

Init. Calib. Date: 01/05/15

Cont. Calib. Time: 0956

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
Naphthalene	1.150	1.107	0.700	AVRG	-3.7
2-Methylnaphthalene	0.673	0.686	0.400	AVRG	1.9
Acenaphthylene	1.724	1.794	0.900	AVRG	4.1
Acenaphthene	1.171	1.195	0.900	AVRG	2.0
Dibenzofuran	1.645	1.657	0.800	AVRG	0.7
Fluorene	1.333	1.358	0.900	AVRG	1.9
Phenanthrene	1.088	1.092	0.700	AVRG	0.4
Anthracene	1.078	1.077	0.700	AVRG	-0.1
Fluoranthene	1.316	1.318	0.600	AVRG	0.2
Pyrene	1.231	1.257	0.600	AVRG	2.1
Benzo(a)anthracene	1.218	1.211	0.800	AVRG	-0.6
Chrysene	1.182	1.158	0.700	AVRG	-2.0
Benzo(b)fluoranthene	1.133	1.153	0.700	AVRG	1.8
Benzo(k)fluoranthene	1.178	1.200	0.700	AVRG	1.9
Benzo(j)fluoranthene	1.160	1.191	0.010	AVRG	2.7
Benzo(a)pyrene	1.116	1.134	0.700	AVRG	1.6
Indeno(1,2,3-cd)pyrene	1.241	1.262	0.500	AVRG	1.7
Dibenzo(a,h)anthracene	1.024	1.076	0.400	AVRG	5.1
Benzo(g,h,i)perylene	1.073	1.074	0.500	AVRG	0.1
1-methylnaphthalene	0.622	0.658	0.010	AVRG	5.8
Perylene	1.122	1.151	0.010	AVRG	2.6
2-Methylnaphthalene-d10	0.707	0.705	0.010	AVRG	-0.3
Dibenzo(a,h)anthracene-d14	0.820	0.880	0.010	AVRG	7.3
Fluoranthene-d10	1.185	1.193	0.010	AVRG	0.7

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF



## SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT8

Cont. Calib. Date: 01/16/15

Init. Calib. Date: 01/05/15

Cont. Calib. Time: 0856

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
=====	=====	=====	=====	=====	=====
Naphthalene	1.150	1.118	0.700	AVRG	-2.8
2-Methylnaphthalene	0.673	0.693	0.400	AVRG	3.0
Acenaphthylene	1.724	1.748	0.900	AVRG	1.4
Acenaphthene	1.171	1.183	0.900	AVRG	1.0
Dibenzofuran	1.645	1.615	0.800	AVRG	-1.8
Fluorene	1.333	1.344	0.900	AVRG	0.8
Phenanthrene	1.088	1.099	0.700	AVRG	1.0
Anthracene	1.078	1.102	0.700	AVRG	2.2
Fluoranthene	1.316	1.349	0.600	AVRG	2.5
Pyrene	1.231	1.229	0.600	AVRG	-0.2
Benzo(a)anthracene	1.218	1.207	0.800	AVRG	-0.9
Chrysene	1.182	1.145	0.700	AVRG	-3.1
Benzo(b)fluoranthene	1.133	1.152	0.700	AVRG	1.7
Benzo(k)fluoranthene	1.178	1.185	0.700	AVRG	0.6
Benzo(j)fluoranthene	1.160	1.153	0.010	AVRG	-0.6
Benzo(a)pyrene	1.116	1.142	0.700	AVRG	2.3
Indeno(1,2,3-cd)pyrene	1.241	1.275	0.500	AVRG	2.7
Dibenzo(a,h)anthracene	1.024	1.042	0.400	AVRG	1.8
Benzo(g,h,i)perylene	1.073	1.099	0.500	AVRG	2.4
1-methylnaphthalene	0.622	0.650	0.010	AVRG	4.5
Perylene	1.122	1.144	0.010	AVRG	2.0
=====	=====	=====	=====	=====	=====
2-Methylnaphthalene-d10	0.707	0.705	0.010	AVRG	-0.3
Dibenzo(a,h)anthracene-d14	0.820	0.878	0.010	AVRG	7.1
Fluoranthene-d10	1.185	1.201	0.010	AVRG	1.4

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/13/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
=====	=====	=====	=====	=====	=====	=====
CCAL	393798	4.60	262932	6.86	448261	8.87
UPPER LIMIT		5.10		7.36		9.37
LOWER LIMIT		4.10		6.36		8.37
01 ZO53MBW1	419217	4.60	279466	6.86	476274	8.88
02 ZO53LCSW1	436630	4.60	287344	6.86	489369	8.88
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IS1 = Naphthalene-d8  
IS2 = Acenaphthene-d10  
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.



SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/13/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
=====	=====	=====	=====	=====	=====	=====
CCAL	482100	13.53	484352	17.30		
UPPER LIMIT		14.03		17.80		
LOWER LIMIT		13.03		16.80		
01 ZO53MBW1	515142	13.54	531933	17.31		
02 ZO53LCSW1	549867	13.54	523541	17.31		
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IS4 = Chrysene-d12

IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/14/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
=====	=====	=====	=====	=====	=====	=====
CCAL	432716	4.59	285009	6.85	514676	8.85
UPPER LIMIT		5.09		7.35		9.35
LOWER LIMIT		4.09		6.35		8.35
01 ZO53LCSDW1	412151	4.58	272668	6.85	470766	8.85
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IS1 = Naphthalene-d8  
IS2 = Acenaphthene-d10  
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.



8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/14/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
=====	=====	=====	=====	=====	=====	=====
CCAL	561143	13.51	529600	17.27		
UPPER LIMIT		14.01		17.77		
LOWER LIMIT		13.01		16.77		
01 ZO53LCSDW1	528600	13.51	439033	17.27		
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IS4 = Chrysene-d12

IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/15/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
=====	=====	=====	=====	=====	=====	=====
CCAL	363073	4.59	238241	6.84	432531	8.85
UPPER LIMIT		5.09		7.34		9.35
LOWER LIMIT		4.09		6.34		8.35
01 MW-36D-14121	609643	4.62	256822	6.85	452479	8.85
02 MW-36S-14121	476880	4.62	269147	6.85	461437	8.85
03 MW-36D-14121	632661	4.62	263212	6.85	466918	8.86
04 MW-36S-14121	485099	4.62	266619	6.85	467656	8.85
05 MW-36D-14121	354198	4.59	230542	6.85	405398	8.85
06 MW-36S-14121	361516	4.59	234780	6.85	409403	8.85
07 MW-36D-14121	381401	4.59	245833	6.85	424418	8.85
08 MW-36S-14121	342594	4.59	231149	6.84	397525	8.85
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23						
24						
25						

IS1 = Naphthalene-d8  
IS2 = Acenaphthene-d10  
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.



8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/15/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
=====	=====	=====	=====	=====	=====	=====
CCAL	471355	13.51	427266	17.27		
UPPER LIMIT		14.01		17.77		
LOWER LIMIT		13.01		16.77		
01 MW-36D-14121	500346	13.50	432368	17.26		
02 MW-36S-14121	506793	13.50	453340	17.26		
03 MW-36D-14121	505814	13.50	435683	17.26		
04 MW-36S-14121	509705	13.50	455118	17.26		
05 MW-36D-14121	467282	13.50	392651	17.26		
06 MW-36S-14121	471825	13.51	400330	17.26		
07 MW-36D-14121	486664	13.50	430000	17.26		
08 MW-36S-14121	454831	13.51	388532	17.27		
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IS4 = Chrysene-d12

IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/16/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
CCAL	353898	4.58	236153	6.84	418975	8.85
UPPER LIMIT		5.08		7.34		9.35
LOWER LIMIT		4.08		6.34		8.35
01 MW-36D-14121	364376	4.59	234794	6.84	416909	8.85
02 MW-36S-14121	358912	4.59	244607	6.84	427198	8.85
03 MW-36D-14121	389742	4.59	257118	6.84	447106	8.85
04 MW-36S-14121	380771	4.59	251175	6.84	432723	8.85
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IS1 = Naphthalene-d8  
IS2 = Acenaphthene-d10  
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
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RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.



8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP15

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/16/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
CCAL	468221	13.51	421063	17.27		
UPPER LIMIT		14.01		17.77		
LOWER LIMIT		13.01		16.77		
01 MW-36D-14121	470542	13.51	417375	17.27		
02 MW-36S-14121	478681	13.51	430521	17.26		
03 MW-36D-14121	503540	13.50	453365	17.26		
04 MW-36S-14121	485546	13.50	423060	17.26		
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 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP16

Project: GAS WORKS PARK-PLAY

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/15/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
CCAL	363073	4.59	238241	6.84	432531	8.85
UPPER LIMIT		5.09		7.34		9.35
LOWER LIMIT		4.09		6.34		8.35
01 RINSATE-1412	350465	4.58	236444	6.85	417746	8.85
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8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP16

Project: GAS WORKS PARK-PLAY

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/15/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
=====	=====	=====	=====	=====	=====	=====
CCAL	471355	13.51	427266	17.27		
UPPER LIMIT		14.01		17.77		
LOWER LIMIT		13.01		16.77		
01 RINSATE-1412	459533	13.51	357415	17.26		
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IS4 = Chrysene-d12

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\* Values outside of QC limits.

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP16

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/13/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
=====	=====	=====	=====	=====	=====	=====
CCAL	393798	4.60	262932	6.86	448261	8.87
UPPER LIMIT		5.10		7.36		9.37
LOWER LIMIT		4.10		6.36		8.37
01 ZP06LCSS1	402831	4.59	270225	6.86	460514	8.87
02 ZP06LCSDS1	388827	4.60	262621	6.86	444839	8.87
03 PAI-9-12.5-1	450772	4.60	282878	6.86	485107	8.88
04 PAI-9-12.5-1	353150	4.60	239110	6.86	413806	8.88
05 PAI-10-9.5-1	462010	4.66	469313	6.88	610062	8.90
06 PAI-10-19.5-	477920	4.61	280009	6.87	467229	8.88
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\* Values outside of QC limits.



8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP16

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/13/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
=====	=====	=====	=====	=====	=====	=====
CCAL	482100	13.53	484352	17.30		
UPPER LIMIT		14.03		17.80		
LOWER LIMIT		13.03		16.80		
01 ZP06LCSS1	501792	13.53	497942	17.30		
02 ZP06LCSDS1	492133	13.54	483539	17.30		
03 PAI-9-12.5-1	527377	13.54	560527	17.31		
04 PAI-9-12.5-1	449873	13.54	458038	17.30		
05 PAI-10-9.5-1	617103	13.56	671044	17.33		
06 PAI-10-19.5-	518884	13.54	526402	17.31		
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IS4 = Chrysene-d12

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AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint

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RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP16

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/14/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
=====	=====	=====	=====	=====	=====	=====
CCAL	432716	4.59	285009	6.85	514676	8.85
UPPER LIMIT		5.09		7.35		9.35
LOWER LIMIT		4.09		6.35		8.35
01 ZP06MBS1	429175	4.59	282514	6.85	484040	8.86
02 PAI-10-24.5-	425344	4.60	288517	6.85	493565	8.86
03 PAI-9-12.5-1	418807	4.59	286732	6.85	486711	8.86
04 PAI-9-12.5-1	409742	4.59	273484	6.85	481890	8.86
05 PAI-10-9.5-1	407927	4.59	271738	6.85	468520	8.85
06 PAI-10-19.5-	417693	4.59	278419	6.85	483336	8.86
07 PAI-10-24.5-	434530	4.59	295410	6.85	522923	8.85
08 PAI-9-12.5-1	409778	4.59	275307	6.85	476785	8.85
09 PAI-10-9.5-1	373128	4.59	253870	6.84	440114	8.85
10 PAI-10-19.5-	367897	4.59	249778	6.85	441906	8.85
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IS1 = Naphthalene-d8  
IS2 = Acenaphthene-d10  
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
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\* Values outside of QC limits.



8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP16

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/14/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
CCAL	561143	13.51	529600	17.27		
UPPER LIMIT		14.01		17.77		
LOWER LIMIT		13.01		16.77		
01 ZP06MBS1	532275	13.51	460216	17.27		
02 PAI-10-24.5-	557200	13.51	527441	17.26		
03 PAI-9-12.5-1	530451	13.51	507774	17.27		
04 PAI-9-12.5-1	524181	13.51	485015	17.26		
05 PAI-10-9.5-1	533034	13.51	502085	17.27		
06 PAI-10-19.5-	551443	13.51	499876	17.27		
07 PAI-10-24.5-	577863	13.51	526858	17.27		
08 PAI-9-12.5-1	513337	13.51	481033	17.27		
09 PAI-10-9.5-1	510687	13.51	460447	17.27		
10 PAI-10-19.5-	490713	13.51	437587	17.26		
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IS4 = Chrysene-d12

IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
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 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

Metals Analysis  
Report and Summary QC Forms

ARI Job ID: ZP15, ZP16



# Cover Page

INORGANIC ANALYSIS DATA PACKAGE



CLIENT: Geoengineers

PROJECT: Gas Works Park-Paly

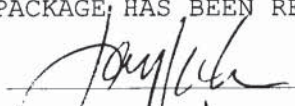
SDG: ZP15

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
MW-36D-141215	ZP15A	14-27518	
MW-36D-141215D	ZP15ADUP	14-27518	
MW-36D-141215S	ZP15ASPK	14-27518	
MW-36S-141215	ZP15B	14-27519	
PBW	ZP15MB1	14-27519	
LCSW	ZP15MB1SPK	14-27519	

Were ICP interelement corrections applied ?                      Yes/No    YES  
Were ICP background corrections applied ?                      Yes/No    YES  
If yes - were raw data generated before  
application of background corrections ?                      Yes/No    NO

Comments: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature:                       Name: Jay Kuhn  
Date: 12/30/14                      Title: Inorganics Director

**INORGANICS ANALYSIS DATA SHEET**

**DISSOLVED METALS**

Page 1 of 1

Sample ID: MW-36D-141215  
SAMPLE

Lab Sample ID: ZP15A

LIMS ID: 14-27518

Matrix: Water

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP15-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: 12/15/14

Date Received: 12/15/14

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/L	Q
6010C	12/18/14	6010C	12/26/14	7440-38-2	Arsenic	0.02	0.2	81.8	
6010C	12/18/14	6010C	12/26/14	7440-70-2	Calcium	0.06	0.2	3.2	
6010C	12/18/14	6010C	12/26/14	7439-89-6	Iron	0.04	0.2	1.6	
6010C	12/18/14	6010C	12/26/14	7439-95-4	Magnesium	0.05	0.2	2.2	
6010C	12/18/14	6010C	12/26/14	7439-96-5	Manganese	0.0014	0.005	0.038	
6010C	12/18/14	6010C	12/26/14	7440-09-7	Potassium	0.3	2	3	
6010C	12/18/14	6010C	12/26/14	7440-23-5	Sodium	0.1	2	979	

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.



**INORGANICS ANALYSIS DATA SHEET  
DISSOLVED METALS**

Sample ID: MW-36S-141215  
SAMPLE

Page 1 of 1

Lab Sample ID: ZP15B


QC Report No: ZP15-Geoengineers

LIMS ID: 14-27519

Project: Gas Works Park-Play Area Investigat

Matrix: Water

0186-846-01 Task 1520

Data Release Authorized: 

Date Sampled: 12/15/14

Reported: 12/30/14

Date Received: 12/15/14

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/L	Q
6010C	12/18/14	6010C	12/26/14	7440-38-2	Arsenic	0.003	0.05	0.07	
6010C	12/18/14	6010C	12/26/14	7440-70-2	Calcium	0.011	0.05	41.6	
6010C	12/18/14	6010C	12/26/14	7439-89-6	Iron	0.008	0.05	14.1	
6010C	12/18/14	6010C	12/26/14	7439-95-4	Magnesium	0.010	0.05	7.06	
6010C	12/18/14	6010C	12/26/14	7439-96-5	Manganese	0.0003	0.001	0.407	
6010C	12/18/14	6010C	12/26/14	7440-09-7	Potassium	0.07	0.5	2.7	
6010C	12/18/14	6010C	12/26/14	7440-23-5	Sodium	0.01	0.5	27.9	

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.



**INORGANICS ANALYSIS DATA SHEET**  
**DISSOLVED METALS**  
 Page 1 of 1

Sample ID: MW-36D-141215  
**MATRIX SPIKE**

Lab Sample ID: ZP15A  
 LIMS ID: 14-27518  
 Matrix: Water  
 Data Release Authorized:  
 Reported: 12/30/14

QC Report No: ZP15-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1520  
 Date Sampled: 12/15/14  
 Date Received: 12/15/14

**MATRIX SPIKE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Arsenic	6010C	81.8	97.8	10.0	160%	H
Calcium	6010C	3.2	51.5	50.0	96.6%	
Iron	6010C	1.6	11.3	10.0	97.0%	
Magnesium	6010C	2.2	50.6	50.0	96.8%	
Manganese	6010C	0.038	2.45	2.50	96.5%	
Potassium	6010C	3	51	50.0	96.0%	
Sodium	6010C	979	1,050	50.0	142%	H

Reported in mg/L

N-Control Limit Not Met  
 H-% Recovery Not Applicable, Sample Concentration Too High  
 NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%



**INORGANICS ANALYSIS DATA SHEET**  
**DISSOLVED METALS**  
Page 1 of 1

Sample ID: MW-36D-141215  
DUPLICATE

Lab Sample ID: ZP15A  
LIMS ID: 14-27518  
Matrix: Water  
Data Release Authorized:  
Reported: 12/30/14

QC Report No: ZP15-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: 12/15/14  
Date Received: 12/15/14



**MATRIX DUPLICATE QUALITY CONTROL REPORT**


Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010C	81.8	66.7	20.3%	+/- 20%	*
Calcium	6010C	3.2	3.3	3.1%	+/- 20%	
Iron	6010C	1.6	1.7	6.1%	+/- 20%	
Magnesium	6010C	2.2	2.2	0.0%	+/- 20%	
Manganese	6010C	0.038	0.040	5.1%	+/- 20%	
Potassium	6010C	3	3	0.0%	+/- 2	L
Sodium	6010C	979	1,010	3.1%	+/- 20%	

Reported in mg/L

\*-Control Limit Not Met  
L-RPD Invalid, Limit = Detection Limit

**INORGANICS ANALYSIS DATA SHEET**  
**DISSOLVED METALS**  
Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: ZP15LCS  
LIMS ID: 14-27519  
Matrix: Water  
Data Release Authorized:   
Reported: 12/30/14

QC Report No: ZP15-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

**BLANK SPIKE QUALITY CONTROL REPORT**


Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Arsenic	6010C	2.14	2.00	107%	
Calcium	6010C	9.81	10.0	98.1%	
Iron	6010C	2.01	2.00	100%	
Magnesium	6010C	10.1	10.0	101%	
Manganese	6010C	0.477	0.500	95.4%	
Potassium	6010C	10.0	10.0	100%	
Sodium	6010C	10.2	10.0	102%	

Reported in mg/L

N-Control limit not met  
Control Limits: 80-120%

**INORGANICS ANALYSIS DATA SHEET**  
**DISSOLVED METALS**  
Page 1 of 1

Sample ID: METHOD BLANK

Lab Sample ID: ZP15MB  
LIMS ID: 14-27519  
Matrix: Water  
Data Release Authorized:   
Reported: 12/30/14

QC Report No: ZP15-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/L	Q
6010C	12/18/14	6010C	12/26/14	7440-38-2	Arsenic	0.003	0.05	0.05	U
6010C	12/18/14	6010C	12/26/14	7440-70-2	Calcium	0.011	0.05	0.05	U
6010C	12/18/14	6010C	12/26/14	7439-89-6	Iron	0.008	0.05	0.05	U
6010C	12/18/14	6010C	12/26/14	7439-95-4	Magnesium	0.010	0.05	0.05	U
6010C	12/18/14	6010C	12/26/14	7439-96-5	Manganese	0.0003	0.001	0.001	U
6010C	12/18/14	6010C	12/26/14	7440-09-7	Potassium	0.07	0.5	0.5	U
6010C	12/18/14	6010C	12/26/14	7440-23-5	Sodium	0.01	0.5	0.5	U

U-Analyte undetected at given DL  
J-Analyte detected between DL and LOQ  
DL-Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.



# Calibration Verification

CLIENT: Geoenigneers

PROJECT: Gas Works Park-Paly

SDG: ZP15

UNITS: ug/L

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	ICP	IF122671	2000.0	2047.58	102.4	2000.0	2054.14	102.7	2052.94	102.6	2056.18	102.8				
Calcium	CA	ICP	IF122671	2000.0	2002.75	100.1	2000.0	1989.73	99.5	1986.28	99.3	1978.80	98.9				
Iron	FE	ICP	IF122671	2000.0	2082.66	104.1	2000.0	2066.14	103.3	2035.73	101.8	2030.86	101.5				
Magnesium	MG	ICP	IF122671	2000.0	1999.58	100.0	2000.0	1997.19	99.9	1998.36	99.9	1984.39	99.2				
Manganese	MN	ICP	IF122671	1000.0	987.78	98.8	1000.0	976.10	97.6	966.38	96.6	968.38	96.8				
Potassium	K	ICP	IF122671	20000.0	20262.32	101.3	20000.0	20110.05	100.6	20059.90	100.3	20100.01	100.5				
Sodium	NA	ICP	IF122671	50000.0	51488.78	103.0	50000.0	50999.11	102.0	51142.82	102.3	50731.04	101.5				

Control Limits: Mercury 80-120; Other Metals 90-110

**CRDL Standard**

CLIENT: Geengineers

PROJECT: Gas Works Park-Paly

SDG: ZP15



UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Arsenic	AS	ICP	IP122671	50.0	50.0	51.18	102.4										
Calcium	CA	ICP	IP122671	50.0	50.0	44.91	89.8										
Iron	FE	ICP	IP122671	50.0	50.0	51.08	102.2										
Magnesium	MG	ICP	IP122671	50.0	50.0	52.11	104.2										
Manganese	MN	ICP	IP122671	1.0	1.0	1.07	107.0										
Potassium	K	ICP	IP122671	500.0	500.0	522.16	104.4										
Sodium	NA	ICP	IP122671	500.0	500.0	461.73	92.3										

Control Limits: no control limits have been established by the EPA at this time.

**Calibration Blanks**

CLIENT: Geoengineers

PROJECT: Gas Works Park-Paly

SDG: ZP15

UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Arsenic	AS	ICP	IPI22671	10.0	50.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U				
Calcium	CA	ICP	IPI22671	5000.0	50.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U				
Iron	FE	ICP	IPI22671	100.0	50.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U				
Magnesium	MG	ICP	IPI22671	5000.0	50.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U				
Manganese	MN	ICP	IPI22671	15.0	1.0	1.0	1.0	U	1.0	U	1.0	U	1.0	U				
Potassium	K	ICP	IPI22671	5000.0	500.0	500.0	500.0	U	500.0	U	500.0	U	500.0	U				
Sodium	NA	ICP	IPI22671	5000.0	500.0	500.0	500.0	U	500.0	U	500.0	U	500.0	U				



# ICP Interference Check Sample



CLIENT: Geoengineers

ICS SOURCE: I.V.

PROJECT: Gas Works Park-Paly

RUNID: IP122671

SDG: ZP15

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSAB1	%R	ICSA2	ICSAB2	%R	ICSA3	ICSAB3	%R
Aluminum	200000	200000	201994.8	203508.1	101.8						
Antimony	1000	1000	14.7	1068.8	106.9						
Arsenic	1000	1000	25.4	1079.3	107.9						
Barium	1000	1000	-0.4	1064.7	106.5						
Beryllium	1000	1000	0.1	1027.6	102.8						
Boron			-3.8		-3.7						
Cadmium	1000	1000	0.7	1049.5	105.0						
Calcium	100000	100000	103106.5	103136.0	103.1						
Chromium	1000	1000	-1.4	1051.7	105.2						
Cobalt	1000	1000	2.5	990.9	99.1						
Copper	1000	1000	0.0	1073.4	107.3						
Iron	200000	200000	201409.4	202663.2	101.3						
Lead	1000	1000	-12.5	1014.0	101.4						
Magnesium	100000	100000	104937.3	100979.2	101.0						
Manganese	1000	1000	0.2	990.2	99.0						
Molybdenum			3.7		3.2						
Nickel	1000	1000	0.6	1004.7	100.5						
Potassium			5.1		22.0						
Selenium	1000	1000	37.1	1095.5	109.6						
Silicon			4.6		4.1						
Silver	1000	1000	-1.0	1108.2	110.8						
Sodium			-41.4		-47.9						
Strontium			3.0		3.0						
Thallium	1000	1000	2.1	978.9	97.9						
Tin			-20.1		-21.3						
Titanium			0.8		0.5						
Vanadium	1000	1000	-1.3	1020.8	102.1						
Zinc	1000	1000	2.7	1004.9	100.5						

# IDLs and ICP Linear Ranges



CLIENT: Geoengineers

PROJECT: Gas Works Park-Paly

SDG: ZP15

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
Arsenic	AS	ICP	OPTIMA ICP 2	197.20		10	50.0	4/1/2012	30000.0	6/10/2014
Calcium	CA	ICP	OPTIMA ICP 2	317.93		5000	50.0	4/1/2012	500000.0	6/10/2014
Iron	FE	ICP	OPTIMA ICP 2	259.94		100	50.0	4/1/2012	250000.0	6/10/2014
Magnesium	MG	ICP	OPTIMA ICP 2	279.08		5000	50.0	4/1/2012	500000.0	6/10/2014
Manganese	MN	ICP	OPTIMA ICP 2	257.61		15	1.0	4/1/2012	30000.0	6/10/2014
Potassium	K	ICP	OPTIMA ICP 2	766.49		5000	500.0	4/1/2012	500000.0	6/10/2014
Sodium	NA	ICP	OPTIMA ICP 2	589.00		5000	500.0	4/1/2012	5000000.0	6/10/2014

# ICP Interelement Correction Factors



CLIENT: Geoenigneers

PROJECT: Gas Works Park-Paly

SDG: ZP15

IEC DATE: 10/27/2014

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	AL	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	14.0198100	0.000000	0.000000
Arsenic	188.98	0.000000	0.000000	0.000000	0.000000	0.0764100	0.000000	-1.0860250	1.6016350	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.1911190	0.000000	0.000000	0.1532030
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0098370	0.000000	0.000000
Boron	249.67	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	2.1785010	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	5.7686950	0.000000	0.000000	0.000000	0.000000	0.1134910	0.000000	0.000000	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.000000	0.000000	0.000000	0.000000	0.0105610	0.000000	0.000000	0.000000	0.000000	0.000000
Cobalt	228.62	0.000000	0.000000	0.0853470	0.000000	0.000000	0.000000	0.000000	-0.0421260	0.000000	-0.0413720
Copper	324.75	0.000000	0.000000	0.000000	0.000000	0.0021980	0.000000	-0.1622490	-0.0152760	0.000000	-0.0447290
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.9508650	0.000000	0.000000
Lead	220.35	-0.2096680	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-1.7516130	1.5325290	0.0499030
Magnesium	279.08	0.000000	0.000000	0.000000	0.000000	0.1119060	0.000000	-1.6746270	-1.1710960	0.000000	0.6610200
Manganese	257.61	0.0060150	0.000000	0.000000	0.000000	0.0041750	0.000000	0.0137770	0.000000	0.000000	-0.0046620
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.0156160	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.5290750	0.000000	0.000000	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	-3.5634780	0.000000	-0.5863590	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Thallium	190.80	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.1475120	0.000000	6.1573050	0.4310930	0.000000	-0.1254120
Titanium	334.90	0.000000	0.000000	0.000000	0.000000	0.0797880	0.000000	0.000000	0.000000	0.000000	0.000000
Vanadium	292.40	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-3.7864850	0.000000	0.0538250
Zinc	206.20	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000



# ICP Interlement Correction Factors



CLIENT: Geoenigneers

PROJECT: Gas Works Park-Paly

SDG: ZP15

IEC DATE: 10/27/2014

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PB	SB	TI	TL	V	ZN
Aluminum	308.22	0.000000	0.000000	15.2105350	0.000000	0.000000	0.000000	1.9181250	0.000000	14.9692830	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.6574750	0.000000	-4.0497020	0.000000
Arsenic	188.98	0.000000	0.000000	3.6569730	0.000000	0.000000	0.000000	-26.2017890	0.000000	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.1112510	0.000000	0.000000	0.000000	0.000000	0.2137080	0.000000
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0122940	0.000000	0.2849470	0.000000
Boron	249.67	0.000000	0.000000	-1.1347080	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	0.000000	0.000000	-0.9151470	0.000000	0.000000	0.000000	0.000000	0.0642140	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.1073910	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.3457620	0.000000
Cobalt	228.62	0.000000	0.000000	-0.1331780	0.1621790	0.000000	0.000000	1.7359160	0.000000	0.000000	0.000000
Copper	324.75	0.0055170	0.000000	0.3194440	0.000000	0.000000	0.000000	0.1761040	0.000000	0.000000	0.000000
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	7.5009230	0.000000
Lead	220.35	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Magnesium	279.08	0.000000	0.000000	-5.2575470	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Manganese	257.61	0.000000	0.000000	0.000000	0.000000	-0.1944900	0.000000	0.000000	0.000000	-0.0205160	0.000000
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	-0.5357680	0.000000	0.4509940	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.6217450	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.2780320	0.000000
Thallium	190.80	0.000000	0.000000	-1.4387970	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	81.4212140
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.0427690	-0.5371860	-0.2280960	0.000000	0.000000	0.000000
Titanium	334.90	0.000000	0.000000	0.9643590	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Vanadium	292.40	0.000000	-0.1516390	-0.4437580	0.000000	0.000000	0.000000	0.5348410	0.000000	0.000000	0.000000
Zinc	206.20	0.000000	0.000000	0.2610130	0.000000	-0.0494060	0.000000	0.000000	0.000000	0.000000	0.000000

FORM XI

# Preparation Log



CLIENT: Geoengineers  
PROJECT: Gas Works Park-Paly  
SDG: ZP15

ANALYSIS METHOD: ICP  
ARI PREP CODE: WMN  
PREPDATE: 12/18/2014

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
MW-36D-141215	ZP15A	0.000	50.0	50.0
MW-36D-141215D	ZP15ADUP	0.000	50.0	50.0
MW-36D-141215S	ZP15ASPK	0.000	50.0	50.0
MW-36S-141215	ZP15B	0.000	50.0	50.0
PBW	ZP15MB1	0.000	50.0	50.0
LCSW	ZP15MB1SPK	0.000	50.0	50.0







# Cover Page

INORGANIC ANALYSIS DATA PACKAGE



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP16

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
PAI-9-12.5-13.0	ZP16A	14-27409	
PAI-9-12.5-13.0D	ZP16ADUP	14-27409	
PAI-9-12.5-13.0S	ZP16ASPK	14-27409	
PAI-9-12.5-13-DUP	ZP16B	14-27410	
PBS	ZP16MB1	14-27410	
LCSS	ZP16MB1SPK	14-27410	
PAI-10-9.5-10.0	ZP16C	14-27411	
PAI-10-24.5-25.0	ZP16E	14-27413	
PAI-9-23-23.5	ZP16I	14-27417	
PBW	ZP16MB2	14-27430	
LCSW	ZP16MB2SPK	14-27430	
RINSATE-141211	ZP16V	14-27430	
RINSATE-141211D	ZP16VDUP	14-27430	
RINSATE-141211S	ZP16VSPK	14-27430	

Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES

If yes - were raw data generated before application of background corrections ? Yes/No NO

Comments: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature: 

Name: Jay Kuhn

Date: 12/30/14

Title: Inorganics Director

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1


Sample ID: PAI-9-12.5-13.0

**SAMPLE**

Lab Sample ID: ZP16A

LIMS ID: 14-27409

Matrix: Soil

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: 12/11/14

Date Received: 12/11/14

Percent Total Solids: 83.2%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/26/14	7440-38-2	Arsenic	0.55	6	478

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**  
Page 1 of 1

Sample ID: PAI-9-12.5-13-DUP  
SAMPLE

Lab Sample ID: ZP16B  
LIMS ID: 14-27410  
Matrix: Soil  
Data Release Authorized:  
Reported: 12/30/14



QC Report No: ZP16-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

Percent Total Solids: 82.2%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/26/14	7440-38-2	Arsenic	0.54	6	463

U-Analyte undetected at given DL  
J-Analyte detected between DL and LOQ  
DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1


Sample ID: PAI-10-9.5-10.0

**SAMPLE**

Lab Sample ID: ZP16C

LIMS ID: 14-27411

Matrix: Soil

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: 12/11/14

Date Received: 12/11/14

Percent Total Solids: 61.1%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/29/14	7440-38-2	Arsenic	1.8	19.9	19.3 J

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**


Page 1 of 1

Sample ID: PAI-10-24.5-25.0  
SAMPLE

Lab Sample ID: ZP16E

LIMS ID: 14-27413

Matrix: Soil

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: 12/11/14

Date Received: 12/11/14

Percent Total Solids: 58.4%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/29/14	7440-38-2	Arsenic	1.8	20	610

U-Analyte undetected at given DL  
J-Analyte detected between DL and LOQ  
DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**


Page 1 of 1

Sample ID: PAI-9-23-23.5  
SAMPLE

Lab Sample ID: ZP16I

LIMS ID: 14-27417

Matrix: Soil

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: 12/11/14

Date Received: 12/11/14

Percent Total Solids: 81.4%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/26/14	7440-38-2	Arsenic	0.52	6	60

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Method Detection Limit


Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: PAI-9-12.5-13.0  
MATRIX SPIKE

Lab Sample ID: ZP16A  
LIMS ID: 14-27409  
Matrix: Soil  
Data Release Authorized:   
Reported: 12/30/14

QC Report No: ZP16-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

**MATRIX SPIKE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Arsenic	6010C	478	771	239	123%	

Reported in mg/kg-dry

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked


Percent Recovery Limits: 75-125%

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: PAI-9-12.5-13.0  
DUPLICATE

Lab Sample ID: ZP16A  
LIMS ID: 14-27409  
Matrix: Soil  
Data Release Authorized:   
Reported: 12/30/14

QC Report No: ZP16-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

**MATRIX DUPLICATE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010C	478	590	21.0%	+/- 20%	*

Reported in mg/kg-dry

\*-Control Limit Not Met  
L-RPD Invalid, Limit = Detection Limit



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: ZP16LCS

LIMS ID: 14-27410

Matrix: Soil

Data Release Authorized:

Reported: 12/30/14



QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: NA

Date Received: NA

**BLANK SPIKE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Arsenic	6010C	213	200	106%	

Reported in mg/kg-dry

N-Control limit not met

NA-Not Applicable, Analyte Not Spiked

Control Limits: 80-120%

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

**Sample ID: METHOD BLANK**

Page 1 of 1

Lab Sample ID: ZP16MB


QC Report No: ZP16-Geoengineers

LIMS ID: 14-27410

Project: Gas Works Park-Play Area Investigat

Matrix: Soil

0186-846-01 Task 1520

Data Release Authorized: 

Date Sampled: NA

Reported: 12/30/14

Date Received: NA

Percent Total Solids: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/26/14	7440-38-2	Arsenic	0.46	5	5 U

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1


Sample ID: RINSATE-141211

SAMPLE

Lab Sample ID: ZP16V

LIMS ID: 14-27430

Matrix: Water

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: 12/11/14

Date Received: 12/11/14

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/L	Q
3010A	12/17/14	6010C	12/26/14	7440-38-2	Arsenic	0.003	0.05	0.05	U

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

**Sample ID: RINSATE-141211  
MATRIX SPIKE**

Lab Sample ID: ZP16V

LIMS ID: 14-27430

Matrix: Water

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1520

Date Sampled: 12/11/14

Date Received: 12/11/14

**MATRIX SPIKE QUALITY CONTROL REPORT**

<b>Analyte</b>	<b>Analysis Method</b>	<b>Sample</b>	<b>Spike</b>	<b>Spike Added</b>	<b>% Recovery</b>	<b>Q</b>
Arsenic	6010C	0.05 U	2.19	2.00	110%	

Reported in mg/L

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: RINSATE-141211  
DUPLICATE

Lab Sample ID: ZP16V  
LIMS ID: 14-27430  
Matrix: Water  
Data Release Authorized:  
Reported: 12/30/14



QC Report No: ZP16-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

**MATRIX DUPLICATE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010C	0.05 U	0.05 U	0.0%	+/- 0.05	L

Reported in mg/L

\*-Control Limit Not Met  
L-RPD Invalid, Limit = Detection Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

**Sample ID: LAB CONTROL**

Lab Sample ID: ZP16LCS  
LIMS ID: 14-27430  
Matrix: Water  
Data Release Authorized  
Reported: 12/30/14



QC Report No: ZP16-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

**BLANK SPIKE QUALITY CONTROL REPORT**

<b>Analyte</b>	<b>Analysis Method</b>	<b>Spike Found</b>	<b>Spike Added</b>	<b>% Recovery</b>	<b>Q</b>
Arsenic	6010C	2.20	2.00	110%	

Reported in mg/L

N-Control limit not met  
Control Limits: 80-120%



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**


Page 1 of 1

**Sample ID: METHOD BLANK**

Lab Sample ID: ZP16MB

LIMS ID: 14-27430

Matrix: Water

Data Release Authorized: 

Reported: 12/30/14

QC Report No: ZP16-Geoengineers

Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1520

Date Sampled: NA

Date Received: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/L	Q
3010A	12/17/14	6010C	12/26/14	7440-38-2	Arsenic	0.003	0.05	0.05	U

U-Analyte undetected at given DL  
J-Analyte detected between DL and LOQ  
DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.

# Calibration Verification



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP16

UNITS: ug/L

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	ICP	IP122671	2000.0	2047.58	102.4	2000.0	2054.14	102.7	2052.94	102.6	2056.18	102.8	2094.84	104.7	2097.45	104.9

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

# Calibration Verification



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

UNITS: ug/L

SDG: ZP16

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	ICP	IP122971	2000.0	2046.13	102.3	2000.0	2031.70	101.6	2021.54	101.1	2098.61	104.9	2122.52	106.1	2127.04	106.4

Control Limits: Mercury 80-120; Other Metals 90-110



# Calibration Verification



CLIENT: Geoenineers

PROJECT: Gas Works Park-Play

SDG: ZP16

UNITS: ug/L

ANALYTE	EL	M	RUN	CCVTV	CCV6	%R	CCV7	%R	CCV8	%R	CCV9	%R	CCV10	%R	CCV11	%R
Arsenic	AS	ICP	IP122971	2000.0	2148.74	107.4										

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

ZP15 : 00200

**CRDL Standard**

CLIENT: Geoenineers

PROJECT: Gas Works Park-Play

SDG: ZP16



UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Arsenic	AS	ICP	IP122671	50.0		51.18	102.4										
Arsenic	AS	ICP	IP122971	50.0		50.05	100.1										

Control Limits: no control limits have been established by the EPA at this time.

# Calibration Blanks

CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP16



UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	CCB1	CCB2	CCB3	CCB4	CCB5	C
Arsenic	AS	ICP	IPI22671	10.0	50.0	50.0	50.0	50.0	50.0	50.0	50.0	U

# Calibration Blanks



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP16

UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Arsenic	AS	ICP	IP122971	10.0	50.0	50.0	v	50.0	v	50.0	v	50.0	v	50.0	v	50.0	v



# Calibration Blanks



CLIENT: Geotechnicians

PROJECT: Gas Works Park-Play

SDG: ZP16

UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	CCB6	CCB7	CCB8	CCB9	CCB10	CCB11	C
Arsenic	AS	ICP	IP122971	10.0	50.0	50.0						U

# ICP Interference Check Sample



CLIENT: Geoenigneers

ICS SOURCE: I.V.

PROJECT: Gas Works Park-Play

RUNID: IP122671

SDG: ZP16

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSAB1	%R	ICSA2	ICSAB2	%R	ICSA3	ICSAB3	%R
Aluminum	200000	200000	201994.8	203508.1	101.8						
Antimony	1000	1000	14.7	1068.8	106.9						
Arsenic	1000	1000	25.4	1079.3	107.9						
Barium	1000	1000	-0.4	1064.7	106.5						
Beryllium	1000	1000	0.1	1027.6	102.8						
Boron			-3.8		-3.7						
Cadmium	1000	1000	0.7	1049.5	105.0						
Calcium	100000	100000	103106.5	103136.0	103.1						
Chromium	1000	1000	-1.4	1051.7	105.2						
Cobalt	1000	1000	2.5	990.9	99.1						
Copper	1000	1000	0.0	1073.4	107.3						
Iron	200000	200000	201409.4	202663.2	101.3						
Lead	1000	1000	-12.5	1014.0	101.4						
Magnesium	100000	100000	104937.3	100979.2	101.0						
Manganese	1000	1000	0.2	990.2	99.0						
Molybdenum			3.7		3.2						
Nickel	1000	1000	0.6	1004.7	100.5						
Potassium			5.1		22.0						
Selenium	1000	1000	37.1	1095.5	109.6						
Silicon			4.6		4.1						
Silver	1000	1000	-1.0	1108.2	110.8						
Sodium			-41.4		-47.9						
Strontium			3.0		3.0						
Thallium	1000	1000	2.1	978.9	97.9						
Tin			-20.1		-21.3						
Titanium			0.8		0.5						
Vanadium	1000	1000	-1.3	1020.8	102.1						
Zinc	1000	1000	2.7	1004.9	100.5						

# ICP Interference Check Sample



CLIENT: Geoenigneers  
PROJECT: Gas Works Park-Play  
SDG: ZP16

ICS SOURCE: I.V.  
RUNID: IP122971  
INSTRUMENT ID: OPTIMA ICP 2  
UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSAB1	%R	ICSA2	ICSAB2	%R	ICSA3	ICSAB3	%R
Aluminum	200000	200000	205572.8	211758.7	105.9						
Antimony		1000	16.6	1069.0	106.9						
Arsenic		1000	18.7	1067.1	106.7						
Barium		1000	-3.2	1025.0	102.5						
Beryllium		1000	0.1	1017.4	101.7						
Boron			-3.6	-4.4							
Cadmium		1000	0.9	1021.4	102.1						
Calcium	100000	100000	103567.3	103318.8	103.3						
Chromium		1000	-0.5	1033.1	103.3						
Cobalt		1000	2.1	970.3	97.0						
Copper		1000	0.1	1059.9	106.0						
Iron	200000	200000	202170.5	208089.7	104.0						
Lead		1000	-11.7	1012.0	101.2						
Magnesium	100000	100000	106670.6	105141.9	105.1						
Manganese		1000	0.0	968.3	96.8						
Molybdenum			3.2	2.9							
Nickel		1000	0.3	1006.1	100.6						
Potassium			-18.4	-26.3							
Selenium		1000	32.4	1083.4	108.3						
Silicon			1.3	3.0							
Silver		1000	-1.1	1115.0	111.5						
Sodium			-68.8	-72.5							
Strontium			3.0	3.0							
Thallium		1000	0.9	971.9	97.2						
Tin			-16.9	-20.3							
Titanium			1.3	1.9							
Vanadium		1000	-2.3	990.0	99.0						
Zinc		1000	3.5	990.4	99.0						

ZP15 : 00200

# IDLs and ICP Linear Ranges



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP16

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
Arsenic	AS	ICP	OPTIMA ICP 2	197.20		10	50.0	4/1/2012	30000.0	6/10/2014



# ICP Interelement Correction Factors



CLIENT: Geoengeers

PROJECT: Gas Works Park-Play

IEC DATE: 10/27/2014

SDG: ZP16

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	AL	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	14.0198100	0.000000	0.000000
Arsenic	188.98	0.000000	0.000000	0.000000	0.000000	0.0764100	0.000000	-1.0860250	1.6016350	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.1911190	0.000000	0.000000	0.1532030
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0098370	0.000000	0.000000
Boron	249.67	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	2.1785010	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	5.7686950	0.000000	0.000000	0.000000	0.000000	0.1134910	0.000000	0.000000	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.000000	0.000000	0.000000	0.000000	0.0105610	0.000000	0.000000	0.000000	0.000000	-0.0413720
Cobalt	228.62	0.000000	0.000000	0.0853470	0.000000	0.000000	0.000000	0.000000	-0.0421260	0.000000	0.000000
Copper	324.75	0.000000	0.000000	0.000000	0.000000	0.0021980	0.000000	-0.1622490	-0.0152760	0.000000	-0.0447290
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.9508650	0.000000	0.000000
Lead	220.35	-0.2096680	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-1.7516130	1.5325290	0.0499030
Magnesium	279.08	0.000000	0.000000	0.000000	0.000000	0.1119060	0.000000	-1.6746270	-1.1710960	0.000000	0.6610200
Manganese	257.61	0.0060150	0.000000	0.000000	0.000000	0.0041750	0.000000	0.0137770	0.000000	0.000000	-0.0046620
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.0156160	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.5290750	0.000000	0.000000	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	-3.5634780	0.000000	-0.5863590	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Thallium	190.80	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	6.1573050	0.4310930	0.000000	-0.1254120
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.1475120	0.000000	0.000000	0.000000	0.000000	0.000000
Titanium	334.90	0.000000	0.000000	0.000000	0.000000	0.0797880	0.000000	0.000000	0.1965610	0.000000	0.000000
Vanadium	292.40	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-3.7864850	0.000000	0.0538250
Zinc	206.20	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000

FORM XI



# ICP Interelement Correction Factors



CLIENT: Geoenigneers

PROJECT: Gas Works Park-Play

SDG: ZP16

IEC DATE: 10/27/2014

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PB	SB	TI	TL	V	ZN
Aluminum	308.22	0.000000	0.000000	15.2105350	0.000000	0.000000	0.000000	1.9181250	0.000000	14.9692830	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.6574750	0.000000	-4.0497020	0.000000
Arsenic	188.98	0.000000	0.000000	3.6569730	0.000000	0.000000	0.000000	-26.2017890	0.000000	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.1112510	0.000000	0.000000	0.000000	0.000000	0.2137080	0.000000
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0122940	0.000000	0.2849470	0.000000
Boron	249.67	0.000000	0.000000	-1.1347080	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	0.000000	0.000000	-0.9151470	0.000000	0.000000	0.000000	0.000000	0.0642140	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.1073910	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.3457620	0.000000
Cobalt	228.62	0.000000	0.000000	-0.1331780	0.1621790	0.000000	0.000000	1.7359160	0.000000	0.000000	0.000000
Copper	324.75	0.0055170	0.000000	0.3194440	0.000000	0.000000	0.000000	0.1761040	0.000000	0.000000	0.000000
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	7.5009230	0.000000
Lead	220.35	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Magnesium	279.08	0.000000	0.000000	-5.2575470	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Manganese	257.61	0.000000	0.000000	0.000000	0.000000	-0.1944900	0.000000	0.000000	0.000000	-0.0205160	0.000000
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	-0.5357680	0.000000	0.4509940	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.6217450	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.2780320	0.000000
Thallium	190.80	0.000000	0.000000	-1.4387970	0.000000	0.000000	0.000000	57.1408130	0.000000	0.000000	81.4212140
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.0427690	0.000000	-0.2280960	0.000000	3.7649150	0.000000
Titanium	334.90	0.000000	0.000000	0.9643590	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Vanadium	292.40	0.000000	-0.1516390	-0.4437580	0.000000	0.000000	0.000000	0.5348410	0.000000	0.000000	0.000000
Zinc	206.20	0.000000	0.000000	0.2610130	0.000000	-0.0494060	0.000000	0.000000	0.000000	0.000000	0.000000

FORM XI

# Preparation Log



CLIENT: Geoengineers

ANALYSIS METHOD: ICP

PROJECT: Gas Works Park-Play

ARI PREP CODE: TWC

SDG: ZP16

PREPDATE: 12/17/2014

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
PBW	ZP16MB2	0.000	50.0	50.0
LCSW	ZP16MB2SPK	0.000	50.0	50.0
RINSATE-141211	ZP16V	0.000	50.0	50.0
RINSATE-141211D	ZP16VDUP	0.000	50.0	50.0
RINSATE-141211S	ZP16VSPK	0.000	50.0	50.0

# Preparation Log



CLIENT: Geoengineers  
PROJECT: Gas Works Park-Play  
SDG: ZP16

ANALYSIS METHOD: ICP  
ARI PREP CODE: SWC  
PREPDATE: 12/23/2014

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
PAI-9-12.5-13.0	ZP16A	1.000	0.0	50.0
PAI-9-12.5-13.0D	ZP16ADUP	1.001	0.0	50.0
PAI-9-12.5-13.0S	ZP16ASPK	1.005	0.0	50.0
PAI-9-12.5-13-DUP	ZP16B	1.041	0.0	50.0
PAI-10-9.5-10.0	ZP16C	1.026	0.0	50.0
PAI-10-24.5-25.0	ZP16E	1.077	0.0	50.0
PAI-9-23-23.5	ZP16I	1.090	0.0	50.0
PBS	ZP16MB1	1.000	0.0	50.0
LCSS	ZP16MB1SPK	1.000	0.0	50.0





**Analysis Run Log**

CLIENT: Geoengineers  
 PROJECT: Gas Works Park-Play  
 SDG: ZP16  
 INSTRUMENT ID: OPTIMA ICP 2  
 RUNID: IP1222671  
 METHOD: ICP  
 START DATE: 12/26/2014  
 END DATE: 12/26/2014

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN					
CCB	CCB3	1.00	12040																																			
PBS	ZP16MB1	2.00	12081																																			
ZZZZZZ	ZP58F	1.00	12121																																			
PAP-9-12.5-13-DUP	ZP16B	2.00	12163																																			
PAP-10-9.5-10.0	ZP16C	2.00	12203																																			
PAP-10-24.5-25.0	ZP16E	2.00	12250																																			
PAP-9-23-23.5	ZP16I	2.00	12292																																			
PAP-9-12.5-13.0D	ZP16ADUP	2.00	12332																																			
PAP-9-12.5-13.0	ZP16A	2.00	12372																																			
PAP-9-12.5-13.0S	ZP16ASP	2.00	12412																																			
LCSS	ZP16MB1SPK	2.00	12452																																			
CCV	CCV4	1.00	12492																																			
CCB	CCB4	1.00	12533																																			
PBS	ZP16MB1	2.00	12574																																			
ZZZZZZ	ZP58MB1	1.00	13020																																			
ZZZZZZ	ZP58E	1.00	13061																																			
ZZZZZZ	ZP58ADUP	1.00	13103																																			
ZZZZZZ	ZP58A	1.00	13144																																			
ZZZZZZ	ZP58ASP	1.00	13260																																			
ZZZZZZ	ZP58B	1.00	13300																																			
ZZZZZZ	ZP58C	1.00	13341																																			
ZZZZZZ	ZP58ME2SPK	1.00	13383																																			
ZZZZZZ	ZP58MB1SPK	1.00	13423																																			
CCV	CCV5	1.00	13463																																			
CCB	CCB5	1.00	13503																																			

# Analysis Run Log

CLIENT: Geoenigneers  
 PROJECT: Gas Works Park-Play  
 SDG: ZP16  
 INSTRUMENT ID: OPTIMA ICP 2  
 RUNID: IP122971  
 METHOD: ICP  
 START DATE: 12/29/2014  
 END DATE: 12/29/2014

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN					
S0			1.00	09561																																		
S2			1.00	10003																																		
S3			1.00	10023																																		
S4			1.00	10050																																		
S5			1.00	10072																																		
ICV			1.00	10225																																		
ICB			1.00	10270																																		
CRI			1.00	10311																																		
ICSA			1.00	10353																																		
ICSAB			1.00	10394																																		
CCV			1.00	10452																																		
CCB			1.00	10493																																		
DI			1.00	10534																																		
ZQ14MB2			1.00	10580																																		
ZQ14EDUP			1.00	11021																																		
ZQ14E			1.00	11063																																		
ZQ14ESPK			1.00	11104																																		
ZP65B			2.00	11144																																		
ZP65H			2.00	11191																																		
ZQ14B			2.00	11235																																		
ZQ14C			2.00	11275																																		
ZQ14MB2SPK			1.00	11315																																		
CCV2			1.00	11355																																		
CCB2			1.00	11395																																		
ZQ14MB1			2.00	11441																																		
ZQ14ADUP			2.00	11482																																		
ZQ14A			2.00	11523																																		
ZQ14ASPK			2.00	11563																																		
ZQ14D			2.00	12003																																		
ZP11A			2.00	12043																																		
ZP11B			2.00	12090																																		
ZP11C			2.00	12130																																		
ZP11D			2.00	12170																																		
ZQ14MB1SPK			2.00	12213																																		
CCV3			1.00	12253																																		







# Analysis Run Log



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP16

INSTRUMENT ID: OPTIMA ICP 2

START DATE: 12/29/2014

RUNID: IP122971

END DATE: 12/29/2014

CLIENT ID

DIL. TIME

%R AG AL AS B BA BE CA CD CO CR CU FE HG K MG MN MO NA NI PB SB SE SI SN TI TL U V ZN

CCB

CCB6

1.00 14582

X

General Chemistry Analysis  
Report and Summary QC Forms

ARI Job ID: ZP15, ZP16

**SAMPLE RESULTS-CONVENTIONALS**  
**ZP15-Geoengineers**



Matrix: Groundwater  
 Data Release Authorized:  
 Reported: 01/23/15

Project: Gas Works Park-Play Area Inv  
 Event: 0186-846-01 Task 1520  
 Date Sampled: 12/15/14  
 Date Received: 12/15/14

**Client ID: MW-36D-141215**  
**ARI ID: 14-27518 ZP15A**

Analyte	Date Batch	Method	Units	RL	Sample
Alkalinity	12/17/14 121714#1	SM 2320	mg/L CaCO3	1.0	1,000
Carbonate	12/17/14	SM 2320	mg/L CaCO3	1.0	112
Bicarbonate	12/17/14	SM 2320	mg/L CaCO3	1.0	893
Hydroxide	12/17/14	SM 2320	mg/L CaCO3	1.0	< 1.0 U
Total Dissolved Solids	12/18/14 121814#1	SM2540C	mg/L	20.0	2,940
Ferrous Iron	12/15/14 121514#1	SM3500 FeD	mg/L	0.040	0.220
N-Nitrate	12/17/14 121714#1	EPA 300.0	mg-N/L	0.5	< 0.5 U
Chloride	12/16/14 121614#1	SM4500-CLE	mg/L	100	497
Sulfate	12/17/14 121714#1	EPA 300.0	mg/L	50.0	604
Sulfide	12/17/14 121714#1	SM4500-S2D	mg/L	10.0	141
Total Organic Carbon	12/19/14 121914#1	SM5310B	mg/L	1.50	7.09
Dissolved Organic Carbon	12/19/14 121914#1	SM5310B	mg/L	1.50	6.85

RL Analytical reporting limit  
 U Undetected at reported detection limit

**SAMPLE RESULTS-CONVENTIONALS**  
**ZP15-Geoengineers**



Matrix: Groundwater  
 Data Release Authorized:  
 Reported: 01/23/15

Project: Gas Works Park-Play Area Inv  
 Event: 0186-846-01 Task 1520  
 Date Sampled: 12/15/14  
 Date Received: 12/15/14

Client ID: MW-36S-141215  
 ARI ID: 14-27519 ZP15B

Analyte	Date Batch	Method	Units	RL	Sample
Alkalinity	12/17/14 121714#1	SM 2320	mg/L CaCO3	1.0	166
Carbonate	12/17/14	SM 2320	mg/L CaCO3	1.0	< 1.0 U
Bicarbonate	12/17/14	SM 2320	mg/L CaCO3	1.0	166
Hydroxide	12/17/14	SM 2320	mg/L CaCO3	1.0	< 1.0 U
Total Dissolved Solids	12/18/14 121814#1	SM2540C	mg/L	5.0	288
Ferrous Iron	12/15/14 121514#1	SM3500 FeD	mg/L	0.800	13.1
N-Nitrate	12/17/14 121714#1	EPA 300.0	mg-N/L	0.1	< 0.1 U
Chloride	12/16/14 121614#1	SM4500-CLE	mg/L	1.0	5.4
Sulfate	12/17/14 121714#1	EPA 300.0	mg/L	1.0	29.2
Sulfide	12/17/14 121714#1	SM4500-S2D	mg/L	0.050	0.512
Total Organic Carbon	12/19/14 121914#1	SM5310B	mg/L	1.50	10.5
Dissolved Organic Carbon	12/19/14 121914#1	SM5310B	mg/L	1.50	8.32

RL Analytical reporting limit  
 U Undetected at reported detection limit



MS/MSD RESULTS-CONVENTIONALS  
ZP15-Geoengineers



Matrix: Groundwater  
Data Release Authorized:  
Reported: 01/23/15

A handwritten signature in black ink, appearing to be 'JL' or similar, written over the 'Data Release Authorized' line.

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: 12/15/14  
Date Received: 12/15/14

Analyte	Method	Date	Units	Sample	Spike	Spike Added	Recovery
ARI ID: ZP15A Client ID: MW-36D-141215							
Ferrous Iron	SM3500 FeD	12/15/14	mg/L	0.220	0.539	0.400	79.8%
Total Organic Carbon	SM5310B	12/19/14	mg/L	7.09	26.5	20.0	97.0%
Dissolved Organic Carbon	SM5310B	12/19/14	mg/L	6.85	27.2	20.0	101.8%

REPLICATE RESULTS-CONVENTIONALS  
ZP15-Geoengineers



Matrix: Groundwater  
Data Release Authorized:  
Reported: 01/23/15

A handwritten signature in black ink, appearing to be 'M. J.', written over the 'Data Release Authorized' text.

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: 12/15/14  
Date Received: 12/15/14

Analyte	Method	Date	Units	Sample	Replicate(s)	RPD/RSD
<b>ARI ID: ZP15A Client ID: MW-36D-141215</b>						
Ferrous Iron	SM3500 FeD	12/15/14	mg/L	0.220	0.200	9.5%
Total Organic Carbon	SM5310B	12/19/14	mg/L	7.09	6.63	6.7%
Dissolved Organic Carbo	SM5310B	12/19/14	mg/L	6.85	6.79	0.9%

LAB CONTROL RESULTS-CONVENTIONALS  
ZP15-Geoengineers



Matrix: Groundwater  
Data Release Authorized  
Reported: 01/23/15


A handwritten signature in black ink, appearing to be 'M' followed by a checkmark-like flourish.

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

Analyte/Method	QC ID	Date	Units	LCS	Spike Added	Recovery
Total Dissolved Solids SM2540C	ICVL	12/18/14	mg/L	488	500	97.6%
Ferrous Iron SM3500 FeD	ICVL	12/15/14	mg/L	0.493	0.500	98.6%
Sulfide SM4500-S2D	ICVL	12/17/14	mg/L	0.521	0.501	104.0%

METHOD BLANK RESULTS-CONVENTIONALS  
ZP15-Geoengineers



Matrix: Groundwater  
Data Release Authorized:   
Reported: 01/23/15

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

Analyte	Method	Date	Units	Blank	ID
Total Dissolved Solids	SM2540C	12/18/14	mg/L	< 5.0 U	
Ferrous Iron	SM3500 FeD	12/15/14	mg/L	< 0.040 U	
N-Nitrate	EPA 300.0	12/17/14	mg-N/L	< 0.1 U	
Chloride	SM4500-CLE	12/16/14	mg/L	< 1.0 U	FB
Sulfate	EPA 300.0	12/17/14	mg/L	< 0.1 U	
Sulfide	SM4500-S2D	12/17/14	mg/L	< 0.050 U	
Total Organic Carbon	SM5310B	12/19/14	mg/L	< 1.50 U	
Dissolved Organic Carbon	SM5310B	12/19/14 12/19/14	mg/L	< 1.50 U 1.70	FB

FB Filtration Blank



STANDARD REFERENCE RESULTS-CONVENTIONALS  
ZP15-Geoengineers




Matrix: Groundwater  
Data Release Authorized  
Reported: 01/23/15

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

Analyte/SRM ID	Method	Date	Units	SRM	True Value	Recovery
Alkalinity ERA #P114506	SM 2320	12/17/14	mg/L CaCO3	58.5	61.7	94.8%
N-Nitrate ERA #320614	EPA 300.0	12/17/14	mg-N/L	2.8	3.0	93.3%
Chloride ERA #290313	SM4500-CLE	12/16/14	mg/L	4.9	5.0	98.0%
Sulfate ERA 131013	EPA 300.0	12/17/14	mg/L	3.0	3.0	100.0%
Total Organic Carbon ERA #0408-13-02	SM5310B	12/19/14	mg/L	21.1	20.0	105.5%
Dissolved Organic Carbon ERA #0408-13-02	SM5310B	12/19/14	mg/L	21.1	20.0	105.5%

SAMPLE RESULTS-CONVENTIONALS  
ZP16-Geoengineers



Matrix: Soil  
Data Release Authorized:   
Reported: 01/05/15

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

Client ID: PAI-10-31.5-32.0  
ARI ID: 14-27414 ZP16F

Analyte	Date	Method	Units	RL	Sample
Preserved Total Solids	12/16/14 121614#1	SM2540G	Percent	0.01	90.01
Sulfide	12/22/14 122214#1	SM4500-S2D	mg/kg	1.10	10.8

RL Analytical reporting limit  
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS  
ZP16-Geoengineers



Matrix: Soil  
Data Release Authorized:  
Reported: 01/05/15

A handwritten signature in black ink, appearing to be 'JW' or similar, written over the 'Data Release Authorized:' text.

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

Client ID: PAI-11-12.0-12.5  
ARI ID: 14-27415 ZP16G

Analyte	Date	Method	Units	RL	Sample
Preserved Total Solids	12/16/14 121614#1	SM2540G	Percent	0.01	68.07
Sulfide	12/22/14 122214#1	SM4500-S2D	mg/kg	291	752

RL Analytical reporting limit  
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS  
ZP16-Geoengineers



Matrix: Soil  
Data Release Authorized:  
Reported: 01/05/15

A handwritten signature in black ink, appearing to be a stylized 'S' or similar character.

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

Client ID: PAI-11-22-22.5  
ARI ID: 14-27416 ZP16H

Analyte	Date	Method	Units	RL	Sample
Preserved Total Solids	12/16/14 121614#1	SM2540G	Percent	0.01	83.03
Sulfide	12/22/14 122214#1	SM4500-S2D	mg/kg	234	851

RL Analytical reporting limit  
U Undetected at reported detection limit



REPLICATE RESULTS-CONVENTIONALS  
ZP16-Geoengineers



Matrix: Soil  
Data Release Authorized:  
Reported: 01/05/15


A handwritten signature in black ink, appearing to be 'M. J.', written over the 'Data Release Authorized' line.

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: 12/11/14  
Date Received: 12/11/14

Analyte	Date	Units	Sample	Replicate(s)	RPD/RSD
ARI ID: ZP16F Client ID: PAI-10-31.5-32.0					
Preserved Total Solids	12/16/14	Percent	90.01	88.13	2.1%

LAB CONTROL RESULTS-CONVENTIONALS  
ZP16-Geoengineers




Matrix: Soil  
Data Release Authorized:   
Reported: 01/05/15

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

Analyte/Method	QC ID	Date	Units	LCS	Spike Added	Recovery
Sulfide SM4500-S2D	PREP	12/22/14	mg/kg	7.13	6.95	102.6%

METHOD BLANK RESULTS-CONVENTIONALS  
ZP16-Geoengineers



Matrix: Soil  
Data Release Authorized:   
Reported: 01/05/15

Project: Gas Works Park-Play Area Inv  
Event: 0186-846-01 Task 1520  
Date Sampled: NA  
Date Received: NA

Analyte	Date	Units	Blank	QC ID
Preserved Total Solids	12/16/14	Percent	< 0.01 U	ICB
Sulfide	12/22/14	mg/kg	< 1.00 U	PREP

Total Solids

ARI Job ID: ZP15, ZP16



Volatiles Total Solids-voats  
Data By: Pat Basilio  
Created: 12/29/14

Worklist: 6260  
Analyst: PAB  
Comments:

Oven ID: \_\_\_\_\_

Balance ID: \_\_\_\_\_

Samples In:            Date: \_\_\_\_\_ Time: \_\_\_\_\_ Temp: \_\_\_\_\_ Analyst: \_\_\_\_\_

Samples Out:          Date: \_\_\_\_\_ Time: \_\_\_\_\_ Temp: \_\_\_\_\_ Analyst: \_\_\_\_\_

ARI ID	Tare Wt (g)	Wet Wt (g)	Dry Wt (g)	% Solids
1. ZP16A 14-27409	_____	_____	_____	\$ 89.30
2. ZP16B 14-27410	_____	_____	_____	\$ 89.60
3. ZP16C 14-27411	_____	_____	_____	\$ 57.30
4. ZP16D 14-27412	_____	_____	_____	\$ 78.60
5. ZP16E 14-27413	_____	_____	_____	\$ 69.60

Worklist ID: 6260      Page: 1  
\* - VOA TS Copied From BETX TS  
% - VOA TS Copied From Metals TS  
\$ - VOA TS Copied From Extraction TS  
\$D - VOA TS Copied From Extraction TS/Decanted

**ZP15:00232**

Extractions Total Solids-exttts  
Data By: Susan D. Dunning  
Created: 12/18/14

Worklist: 2600  
Analyst: SDRD  
Comments:

Oven ID: \_\_\_\_\_

Balance ID: \_\_\_\_\_

Samples In:            Date: \_\_\_\_\_ Time: \_\_\_\_\_ Temp: \_\_\_\_\_ Analyst: \_\_\_\_\_

Samples Out:           Date: \_\_\_\_\_ Time: \_\_\_\_\_ Temp: \_\_\_\_\_ Analyst: \_\_\_\_\_

	ARI ID CLIENT ID	Tare Wt (g)	Wet Wt (g)	Dry Wt (g)	% TS	Dcnt	5g	10g	12.5g
1.	ZP16A 14-27409 PAI-9-12.5-13.0	1.19	10.99	9.94	89.3	No	5.60	11.20	14.00
2.	ZP16B 14-27410 PAI-9-12.5-13-DUP	1.19	10.80	9.80	89.6	No	5.58	11.16	13.95
3.	ZP16C 14-27411 PAI-10-9.5-10.0	1.19	10.49	6.52	57.3	No	8.73	17.45	21.82
4.	ZP16D 14-27412 PAI-10-19.5-20.0	1.19	9.27	7.54	78.6	No	6.36	12.72	15.90
5.	ZP16E 14-27413 PAI-10-24.5-25.0	1.18	10.68	7.79	69.6	No	7.18	14.37	17.96

Extractions Total Solids-extts  
Data By: Susan D. Dunning  
Created: 12/18/14

Worklist: 2600  
Analyst: SDRD  
Comments:

Oven ID: 015

Balance ID: B334705934

Samples In: Date: 12/18/14 Time: 14:50 Temp: 105 Analyst: TH

Samples Out: Date: 12/19/14 Time: 0800 Temp: 103 Analyst: SC

ARI ID CLIENT ID	Tare Wt (g)	Wet Wt (g)	Dry Wt (g)	% TS	Dcnt	5g	10g	12.5g
1. ZP16A 14-27409 PAI-9-12.5-13.0	<u>1.19</u>	<u>10.99</u>	<u>9.94</u>		No			
2. ZP16B 14-27410 PAI-9-12.5-13-DUP	<u>1.19</u>	<u>10.80</u>	<u>9.80</u>		No			
3. ZP16C 14-27411 PAI-10-9.5-10.0	<u>1.19</u>	<u>10.49</u>	<u>6.52</u>		No			
4. ZP16D 14-27412 PAI-10-19.5-20.0	<u>1.19</u>	<u>9.27</u>	<u>7.54</u>		No			
5. ZP16E 14-27413 PAI-10-24.5-25.0	<u>1.18</u>	<u>10.68</u>	<u>7.79</u>		No			

Solids Data Entry Report  
Date: 12/23/14

Checked by: (S) Date: 12/23/14  
Data Analyst: DM

Solids Determination performed on 12/22/14 by DM

JOB	SAMPLE	CLIENTID	TAREWEIGHT	SAMPDISH	DRYWEIGHT	SOLIDS
ZP16	A	PAI-9-12.5-13.0	1.019	10.177	8.635	83.16
ZP16	B	PAI-9-12.5-13-DUP	1.010	10.041	8.435	82.22
ZP16	C	PAI-10-9.5-10.0	1.002	10.597	6.866	61.12
ZP16	E	PAI-10-24.5-25.0	0.979	10.129	6.318	58.35
ZP16	I	PAI-9-23-23.5	1.012	10.142	8.446	81.42





# Total Solids Bench Sheet

Laboratory Section Metals

Oven Identification: 07 Balance ID: 068755

Samples in Oven: Date: 12-22-19 Time: 1340 Temp: 105°C Analyst: DM

Removed from Oven: Date: 12-23-14 Time: 0635 Temp: 104°C Analyst: DM

ARI Sample ID	Tare Weight (g)	Tare + Sample Wet (g)	Tare + Sample Dry (g)	Date & Time Last Weight	Final Weighting >12 hrs <sup>1</sup>
ZP14 A	0.974	10.557	9.361	-	✓
" B	1.029	10.390	9.864	-	✓
" C	1.043	10.690	10.025	-	✓
" D	1.049	10.727	10.038	-	✓
ZP06 A	1.011	5.918	1.884	-	✓
ZP35 A	1.042	10.426	6.580	-	✓
" C	1.013	10.786	9.721	-	✓
ZP16 A	1.019	10.177	8.635	-	✓
" B	1.010	10.041	8.436	-	✓
" C	1.002	10.597	6.866	-	✓
" E	0.979	10.129	6.318	-	✓
" I	1.012	10.142	8.446	-	✓
ZP06 A	0.997	10.545	7.063	-	✓
ZP11 A	1.049	10.848	7.924	-	✓
" B	0.989	10.831	7.615	-	✓
" C	1.019	10.500	8.671	-	✓
" D	1.022	10.407	6.775	-	✓
" E	0.981	10.717	8.184	-	✓
" F	1.068	10.351	9.475	-	✓
" G	1.052	10.209	9.191	-	✓
" H	1.012	10.218	8.563	-	✓
" I	1.058	10.241	9.220	-	✓
" J	1.022	10.533	9.534	-	✓

1) Place a check mark in this column if samples have dried > 12 but < 24 hours. When samples have been at 104°C < 12 hours, constant weight must be verified as described in SOP 10023S. Use a 2<sup>nd</sup> bench sheet for additional weightings.

Subcontracted Results  
Sulfur Analyzed by Horizon Labs

ARI Job ID: ZP15, ZP16



General Offices: P.O. Box 995 Price, Utah 84501 435-637-4343  
Laboratory: 545 East 100 North Price, Utah 84501

Submitted to:

December 31, 2014

**Analytical Resources Inc.**

Mark Harris

4611 S. 134th Place

Suite 100

Tukwila, WA 98168-3240

Date Sampled: 12/11-12/2014

Date Received: 12/18/2014

Sample Identification:

See Below

Sampled By: ARI

Identification By: ARI

Analysis Report #: See Below

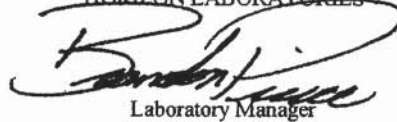
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### CERTIFICATE OF ANALYSIS

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Analysis #	ARI ID	Sulfur (Acid Insoluble)
90188	14-27452 ZO83M	0.37
90189	14-27453 ZO83N	0.63
90190	14-27454 ZO83O	1.31
90191	14-27455 ZO83P	0.28
90192	14-27456 ZO83Q	0.03
90193	14-27418 ZP16J	0.04
90194	14-27419 ZP16K	0.56
90195	14-27420 ZP16L	0.12

Respectfully Submitted,  
HORIZON LABORATORIES



Laboratory Manager

ZP15 : 00238

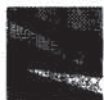
Name	Analysis Date		
Blank	12/29/2014 14:06		
	Sample Mass	Operator	Sulfur Concentration
	1.0000 g	Francesca	0.00%
	1.0000 g	Francesca	0.00%
Name	Analysis Date		
3.25 std	12/29/2014 14:21		
	Sample Mass	Operator	Sulfur Concentration
	0.2036 g	Francesca	3.12%
	0.2049 g	Francesca	3.24%
	0.2598 g	Francesca	3.26%
Name	Analysis Date		
90188	12/29/2014 14:25		
	Sample Mass	Operator	Sulfur Concentration
	0.1170 g	Francesca	0.37%
Name	Analysis Date		
90189	12/29/2014 14:27		
	Sample Mass	Operator	Sulfur Concentration
	0.1197 g	Francesca	0.63%
Name	Analysis Date		
90190	12/29/2014 14:29		
	Sample Mass	Operator	Sulfur Concentration
	0.1228 g	Francesca	1.31%
Name	Analysis Date		
90191	12/29/2014 14:32		
	Sample Mass	Operator	Sulfur Concentration
	0.1084 g	Francesca	0.27%
Name	Analysis Date		
90192	12/29/2014 14:34		
	Sample Mass	Operator	Sulfur Concentration
	0.1442 g	Francesca	0.03%
Name	Analysis Date		
90193	12/29/2014 14:40		
	Sample Mass	Operator	Sulfur Concentration
	0.1820 g	Francesca	0.04%
Name	Analysis Date		
90194	12/29/2014 14:42		
	Sample Mass	Operator	Sulfur Concentration
	0.1086 g	Francesca	0.56%
Name	Analysis Date		
90195	12/29/2014 14:44		
	Sample Mass	Operator	Sulfur Concentration
	0.1644 g	Francesca	0.12%
Name	Analysis Date		
90191	12/29/2014 14:46		
	Sample Mass	Operator	Sulfur Concentration
	0.1097 g	Francesca	0.29%
Name	Analysis Date		
STD BL-4a	12/29/2014 14:48		
	Sample Mass	Operator	Sulfur Concentration
	0.1592 g	Francesca	0.21%



Sample I.D.	Pan Weight	Pan & Sample Wt.	Pan & Dry Sample Wt.	Air Dry Loss (Moisture) %
90188	224.7	289.9	279.8	15.49
90189	227.3	290.9	273.2	27.83
90190	222.2	291.9	279.5	17.79
90191	225	335.7	312.4	21.04
90192	220.8	360.1	345.4	10.55
90193	222.2	445.6	417.7	12.49
90194	222.9	354.9	313.1	31.67
90195	223.6	403.4	352.5	28.31

Subcontracted Results  
Metals (Sub) Analyzed by Applied Speciation & Consulting

ARI Job ID: ZP15, ZP16



January 15, 2015

Cheronne Oreiro  
Analytical Resources Inc.  
4611 S. 134<sup>th</sup> Place Suite 100  
Tukwila, WA 98168  
(206) 695-6200

Re: SDG A141217S1 (Gas Works Park – Play Area Investigation)

Ms. Oreiro,

Attached is the report associated with six (6) aqueous samples submitted for arsenite, arsenate, monomethylarsonic acid, and dimethylarsinic acid quantitation on December 17, 2014. The samples were received on the day of submittal in a sealed cooler at 0.3°C. Arsenite, arsenate, monomethylarsonic acid, and dimethylarsinic acid quantitation was performed via ion chromatography inductively coupled plasma collision reaction cell mass spectrometry (IC-ICP-CRC-MS). Any issues associated with the analyses are addressed in the following report.

If you have any questions, please feel free to contact me at your convenience.

Sincerely,

Ben Wozniak  
Project Manager  
Applied Speciation and Consulting, LLC

Applied Speciation and Consulting, LLC

Report Prepared for:

Cheronne Oreiro  
Analytical Resources Inc.  
4611 S. 134th Place Suite 100  
Tukwila, WA 98168

January 15, 2015

### 1. Sample Reception

Six (6) aqueous samples were submitted for arsenite, arsenate, monomethylarsonic acid, and dimethylarsinic acid quantitation on December 17, 2014. The samples were received on the day of submittal in a sealed cooler at 0.3°C, as indicated on the attached chain of custody (COC) forms.

The samples were received in a laminar flow clean hood, void of trace metals contamination and ultra-violet radiation, and assigned discrete sample identifiers. All samples were then stored in a secure, monitored refrigerator (maintained at a temperature of 4°C) until the analyses could occur.

It should be noted that one aqueous sample, identified as PAI-10GW, was included in the sample shipment but was not originally listed on any of the COC forms submitted with the samples. The client was contacted and submitted a new COC for this sample. Both the original COCs and this additional COC have been included in this report.

### 2. Sample Preparation

All sample preparation is performed in laminar flow clean hoods known to be free from trace metals contamination. All applied water for dilutions and sample preservatives are also monitored for contamination to account for any biases associated with the sample results.

*Arsenic Speciation Analysis by IC-ICP-CRC-MS* Each sample submitted in an evacuated vial containing an EDTA preservative (provided by Applied Speciation and Consulting) was analyzed as received, without further chemical preservation.

### 3. Sample Analysis

All sample analysis is preceded by a minimum of a five-point calibration curve spanning the entire concentration range of interest. All calibration curves, associated with each species of interest, are standardized by linear regression resulting in a response factor. All sample results are **instrument blank corrected** to account for any operational biases.



Prior to sample analysis, all calibration curves are verified using second source standards which are identified as initial calibration verification standards (ICV).

Ongoing instrument performance is identified by the analysis of continuing calibration verification standards (CCV) and continuing calibration blanks (CCB) at a minimum interval of every ten analytical runs.

*Arsenic Speciation Analysis by IC-ICP-CRC-MS* All samples for arsenite, arsenate, monomethylarsonic acid, and dimethylarsinic acid quantitation were analyzed by ion chromatography inductively coupled plasma collision reaction cell mass spectrometry (IC-ICP-CRC-MS) on December 17-18, 2014. Aliquots of each sample are injected onto an anion exchange column and are mobilized by an alkaline (pH > 7) gradient. The eluting arsenic species are then introduced into a radio frequency (RF) plasma where energy-transfer processes cause desolvation, atomization, and ionization. The ions are extracted from the plasma through a differentially-pumped vacuum interface and travel through a pressurized chamber (CRC) containing a specific collision gas. Polyatomic interferences, due to their inherently larger size, collide more frequently with the collision gas and therefore may be separated from the analyte of interest via kinetic energy discrimination (KED). A solid-state detector detects ions transmitted through the mass analyzer on the basis of their mass-to-charge ratio (m/z), and the resulting current is processed by a data handling system.

Retention times for each eluting species are compared to known standards for species identification.

#### **4. Analytical Issues**

No significant analytical issues were encountered during the requested analyses. All quality control parameters associated with these samples were within acceptance limits.

It should be noted that several additional arsenic-containing species were detected in the submitted samples during the speciation analyses. While the identities of these species could not be confirmed at this time, the estimated concentration of arsenic associated with each detected species has been provided for each sample in the attached results table. Applied Speciation and Consulting may be able to pursue identification of these additional species upon client request.

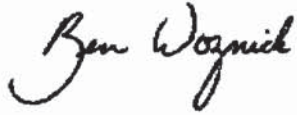
The estimated method detection limits (eMDLs) for arsenite, arsenate, and dimethylarsinic acid are generated from replicate analyses of the lowest standard in the calibration curve. Not all arsenic species are present in preparation blanks; therefore, eMDL calculations based on preparation blanks may be artificially biased low for individual arsenic species.

The eMDL for monomethylarsonic acid is calculated using the average eMDL of the species present in the calibration curve (*i.e.*, arsenite, arsenate, and dimethylarsinic acid). The ICP-CRC-MS determinative method exhibits a species-independent response, which is confirmed by the analysis of ICV standards for each species. Monomethylarsonic acid is not included

in the calibration standards, CCV standards, and matrix spikes due to impurities which would bias the results for other arsenic species.

If you have any questions regarding this report, please feel free to contact me.

Sincerely,

A handwritten signature in black ink that reads "Ben Wozniak". The signature is written in a cursive style with a large, looping initial "B".

Ben Wozniak  
Project Manager  
Applied Speciation and Consulting, LLC

Arsenic Speciation Results for ARI  
SDG: A141217S1

Contact: Cheronne Oreiro

Report Date: January 15, 2015

Report Generated by: Ben Wozniak  
Applied Speciation and Consulting, LLC

**Sample Results**

ARI Sample ID	Date Sampled	Dilution	As(III)	As(V)	MMAs	DMAs	Unk 8.4*	Unk 9.2*	Unk 10.2*	Unk 13.1*
14-27440-ZO83A	12/12/14	50	28.9	1.74	< 0.21 U	< 0.23 U	14.2	2.94	52.5	7.60
14-27441-ZO83B	12/12/14	500	702	141	< 2.1 U	< 2.3 U	144	7.9 J	120	19.4
14-27442-ZO83C	12/12/14	25000	42900	1200	< 110 U	< 120 U	1100	620	10500	680
14-27173-ZO53A	12/11/14	1000	482	15.4 J	< 4.2 U	< 4.6 U	102	128	2240	8.2 J
14-27518-ZP15A	12/15/14	50000	39700	790 J	< 210 U	< 230 U	3920	1210	19400	18700
14-27519-ZP15B	12/15/14	100	37.7	6.41	< 0.42 U	< 0.46 U	< 0.42 U	< 0.42 U	< 0.42 U	1.65 J

All results reflect the applied dilution and are reported in µg/L

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)

J = Sample concentration is between the eMDL and the Reporting Limit (RL)

\* Unknown arsenical species; please see narrative



Arsenic Speciation Results for ARI  
SDG: A141217S1

Contact: Cheronne Oreiro

Report Date: January 15, 2015

Report Generated by: Ben Wozniak  
Applied Speciation and Consulting, LLC

**Quality Control Summary - Preparation Blanks**

Analyte	Units	PBW1	PBW2	PBW3	PBW4	Mean	StdDev	eMDL* at			
								1x	1x	50x	
As(III)	µg/L	0.000	0.000	0.000	0.000	0.000	0.000	0.003	0.020	0.13	1.0
As(V)	µg/L	0.000	0.000	0.000	0.000	0.000	0.000	0.005	0.020	0.27	1.0
MMAs	µg/L	0.000	0.000	0.000	0.000	0.000	0.000	0.004	0.020	0.21	1.0
DMAs	µg/L	0.000	0.000	0.000	0.000	0.000	0.000	0.005	0.021	0.23	1.0

eMDL = Estimated Method Detection Limit; RL = Reporting Limit

\*Please see narrative regarding eMDL calculations

**Quality Control Summary - Certified Reference Materials**

Analyte	Units	CRM	True Value	Result	Recovery
As(III)	µg/L	ICV	5.000	5.079	101.6
As(V)	µg/L	ICV	5.000	4.736	94.7
MMAs	µg/L	ICV	4.610	4.680	101.5
DMAs	µg/L	ICV	3.625	3.682	101.6



Arsenic Speciation Results for ARI  
SDG: A141217S1

Contact: Cheronne Oreiro

Report Date: January 15, 2015  
Report Generated by: Ben Wozniak  
Applied Speciation and Consulting, LLC

**Quality Control Summary - Matrix Duplicate**

Analyte	Units	Sample ID	Rep 1	Rep 2	Mean	RPD
As(III)	µg/L	14-27519-ZP15B	37.72	37.36	37.54	1.0
As(V)	µg/L	14-27519-ZP15B	6.41	6.24	6.33	2.8
MMAs	µg/L	14-27519-ZP15B	< 0.42 U	< 0.42 U	NC	NC
DMAs	µg/L	14-27519-ZP15B	< 0.46 U	< 0.46 U	NC	NC

NC = Value was not calculated due to one or more concentrations below the eMDL

**Quality Control Summary - Matrix Spike/ Matrix Spike Duplicate**

Analyte	Units	Sample ID	MS Spike		MS		MSD Spike		MSD	
			Conc	Result	Result	Recovery	Conc	Result	Recovery	RPD
As(III)	µg/L	14-27519-ZP15B	100.0	139.5	102.0	102.0	100.0	138.2	100.7	0.9
As(V)	µg/L	14-27519-ZP15B	100.0	105.3	99.0	99.0	100.0	103.9	97.6	1.3
DMAs	µg/L	14-27519-ZP15B	104.9	106.9	101.9	101.9	104.9	105.3	100.4	1.5



January 20, 2015

Cheronne Oreiro  
Analytical Resources Inc.  
4611 S. 134<sup>th</sup> Place Suite 100  
Tukwila, WA 98168  
(206) 695-6200

Re: SDG A141217S2 (Gas Works Park – Play Area Investigation)

Ms. Oreiro,

Attached is the report associated with eight (8) soil core samples submitted for arsenic sequential extraction on December 17, 2014. The samples were received on the day of submittal in a sealed cooler at 0.3°C. Any issues associated with the analyses are addressed in the following report.

If you have any questions, please feel free to contact me at your convenience.

Sincerely,

Ben Wozniak  
Project Manager  
Applied Speciation and Consulting, LLC

Applied Speciation and Consulting, LLC

Report Prepared for:

Cheronne Oreiro  
Analytical Resources Inc.  
4611 S. 134th Place Suite 100  
Tukwila, WA 98168

January 20, 2015

### 1. Sample Reception

Eight (8) soil core samples were submitted for arsenic sequential extraction on December 17, 2014. The samples were received on the day of submittal in a sealed cooler at 0.3°C, as indicated on the attached chain of custody (COC) forms.

All samples were received in a laminar flow clean hood, void of trace metals contamination and ultra-violet radiation. Immediately upon reception all samples were designated discrete sample identifiers and then stored in a secure, monitored freezer (maintained at a temperature of  $\leq -10^{\circ}\text{C}$ ) until sample preparation could be performed.

### 2. Sample Preparation

All applied water for dilutions and sample reagents are monitored for contamination to account for any biases associated with the sample results.

Arsenic Sequential Extraction (Wenzel et al.) A sequential extraction method, based on Wenzel et al., was employed for correlation between arsenic and different substrate properties. In accordance with the client's request, all handling of the submitted soil core samples and preparation of the first four extracts (i.e., all extracts except the  $\text{HNO}_3 / \text{H}_2\text{O}_2$  extraction) was performed in a glove box maintained under anoxic conditions.

A known mass of each soil sample was transferred to polypropylene vial and an aliquot of 0.05 M  $(\text{NH}_4)_2\text{SO}_4$  was added to each vial. Each vial was capped and shaken on an inverting shaker for 4 hours at room temperature at 30 RPM.

The samples were removed from the shaker and centrifuged for 20 minutes at 3000RPM. After the supernatant was decanted into a separate vial for trace metals analysis, an aliquot of reagent water was added to each vial. The vials were shaken vigorously and centrifuged for 20 minutes at 3000RPM. The reagent water rinse was decanted and discarded.

An aliquot of 0.05 M  $\text{NH}_4\text{H}_2\text{PO}_4$  was added to each vial. Each vial was capped and shaken on an inverting shaker for 16 hours at room temperature at 30 RPM.



The samples were removed from the shaker and centrifuged for 20 minutes at 3000RPM. After the supernatant was decanted into a separate vial for trace metals analysis, an aliquot of reagent water was added to each vial. The vials were shaken vigorously and centrifuged for 20 minutes at 3000RPM. The reagent water rinse was decanted and discarded.

All sample vials were wrapped in aluminum foil to prevent photo-oxidation and then an aliquot of 0.2M ammonium oxalate buffer (pH=3.25) was added to each vial. Each vial was capped and shaken on an inverting shaker for 4 hours at room temperature at 30 RPM.

The samples were removed from the shaker and centrifuged for 20 minutes at 3000RPM. After the supernatant was decanted into a separate vial for trace metals analysis, an aliquot of 0.2M ammonium oxalate buffer was added to each vial. The vials were shaken vigorously and centrifuged for 20 minutes at 3000RPM. The supernatant was decanted and discarded.

An aliquot of 0.2M ammonium oxalate buffer with 0.1M ascorbic acid was added to each vial. The vials were then placed in a hotblock digestion apparatus at 96°C for 30 minutes.

The samples were removed from the shaker and centrifuged for 20 minutes at 3000RPM. After the supernatant was decanted into a separate vial for trace metals analysis, an aliquot of 0.2M ammonium oxalate buffer with 0.1M ascorbic acid was added to each vial. The vials were shaken vigorously and centrifuged for 20 minutes at 3000RPM. The supernatant was decanted and discarded.

The residual solid pellets remaining in the vials were then digested via with aliquots of concentrated HNO<sub>3</sub> and H<sub>2</sub>O<sub>2</sub> (in accordance with EPA Method 3050B).

In accordance with the client's request, all extracts were analyzed for iron in addition to arsenic.

### 3. Sample Analysis

All sample analysis is preceded by a minimum of a five-point calibration curve spanning the entire concentration range of interest. All calibration curves, associated with each species of interest, are standardized by linear regression resulting in a response factor. All sample results are instrument blank corrected to account for any operational biases associated with the analytical platform.

Prior to sample analysis, all calibration curves are verified using second source standards which are identified as initial calibration verification standards (ICV).

Ongoing instrument performance is identified by the analysis of continuing calibration verification standards (CCV) and continuing calibration blanks (CCB) at a minimum interval of every ten analytical runs.

Arsenic and Iron Quantitation by ICP-OAO-MS All sample fractions for arsenic and iron quantitation were analyzed by inductively coupled plasma triple quadrupole mass



spectrometry (ICP-QQQ-MS). An aliquot of each sample extract is introduced into a radio frequency (RF) plasma where energy-transfer processes cause desolvation, atomization, and ionization. The ions are extracted from the plasma through a differentially-pumped vacuum interface and travel through an initial quadrupole (Q1), which filters the target masses prior to their entrance into a second chamber. The second chamber contains specific reactive gasses or collision gasses that preferentially react either with interfering ions of the same target mass to charge ratios (m/z) or with the target analyte, producing an entirely different mass to charge ratio (m/z) which can then be differentiated from the initial interferences. The ions then exit the collision/reaction chamber into the mass analyzer (Q2). A solid-state detector detects ions transmitted through the mass analyzer, on the basis of their mass-to-charge ratio (m/z), and the resulting current is processed by a data handling system.

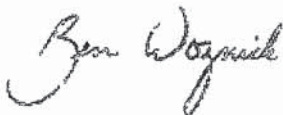
Total Solids Analysis A known mass of each sample was placed into a pre-weighed pan, and the combined mass of the sample and pan was recorded. All samples were then placed into a convection oven maintained at a temperature of 60°C. After drying for a minimum of twelve (12) hours, all samples were briefly cooled and reweighed. The total solids percentage of each sample was calculated by dividing the weight of the dried sample by the weight of the original sample.

#### 4. Analytical Discussion

No significant issues were encountered with the requested analyses. All quality control parameters associated with the samples were within acceptance limits.

If you have any questions or concerns regarding this report, please feel free to contact me.

Sincerely,



Ben Wozniak  
Project Manager  
Applied Speciation and Consulting, LLC

Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

Sample Results

14-27414-ZP16F

Fraction	Fraction Description	As	Fe
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	4.68	17.1
0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	1.60	159
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	1.11	1480
0.1M Ascorbic acid	Crystalline Fe oxyhydroxide	0.493	715
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	3.71	6640

All analyte results are reported in µg/g on a dry weight (DW) basis

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)

J = Sample concentration is between the eMDL and the Reporting Limit (RL)

Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

**Sample Results**

**14-27415-ZP16G**

<b>Fraction</b>	<b>Fraction Description</b>	<b>As</b>	<b>Fe</b>
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	0.556 J	208
0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	2.03	508
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	1.47	2430
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> +			
0.1M Ascorbic acid	Crystalline Fe oxyhydroxide	0.565 J	731
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	19.2	12700

All analyte results are reported in µg/g on a dry weight (DW) basis

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)

J = Sample concentration is between the eMDL and the Reporting Limit (RL)

Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

Sample Results

14-27416-ZP16H

Fraction	Fraction Description	As	Fe
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	1.93	36.8
0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	3.75	392
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	0.790	1170
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> +			
0.1M Ascorbic acid	Crystalline Fe oxyhydroxide	0.845	834
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	35.4	13800

All analyte results are reported in µg/g on a dry weight (DW) basis

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)

J = Sample concentration is between the eMDL and the Reporting Limit (RL)



Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

Sample Results

14-27447-ZO83H

Fraction	Fraction Description	As	Fe
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	18.5	35.0
0.05 M (NH <sub>4</sub> ) <sub>2</sub> H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	40.6	19.2 J
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	108	542
0.1M Ascorbic acid	Crystalline Fe oxyhydroxide	26.8	1380
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	93.0	1630

All analyte results are reported in µg/g on a dry weight (DW) basis

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)  
 J = Sample concentration is between the eMDL and the Reporting Limit (RL)

Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

Sample Results

14-27448-Z083I

Fraction	Fraction Description	As	Fe
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	20.3	5.6 J
0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	908	32.6
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	4790	19200
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> + Ascorbic acid	Crystalline Fe oxyhydroxide	2620	4300
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	750	4870

All analyte results are reported in µg/g on a dry weight (DW) basis

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)

J = Sample concentration is between the eMDL and the Reporting Limit (RL)

Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

Sample Results

14-27449-ZO83J

Fraction	Fraction Description	As	Fe
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	354	55.1
0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	180	255
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	24.7	718
0.1M Ascorbic acid	Crystalline Fe oxyhydroxide	78.5	532
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	9110	6120

All analyte results are reported in µg/g on a dry weight (DW) basis

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 J = Sample concentration is between the eMDL and the Reporting Limit (RL)

Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

Sample Results

14-27450-ZO83K

Fraction	Fraction Description	As	Fe
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	7.86	12.1 J
0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	8.07	180
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	3.10	799
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> + 0.1M Ascorbic acid	Crystalline Fe oxyhydroxide	2.89	491
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	37.6	11900

All analyte results are reported in µg/g on a dry weight (DW) basis

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)  
 J = Sample concentration is between the eMDL and the Reporting Limit (RL)



Arsenic Sequential Extraction Results for ARI  
 SDG. A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

Sample Results

14-27451-Z083L

Fraction	Fraction Description	As	Fe
0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	Non-specifically adsorbed arsenic	6.84	4.6 J
0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	Specifically-sorbed arsenic	0.832	338
0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	Amorphous Fe oxyhydroxide	0.721	1200
0.1M Ascorbic acid	Crystalline Fe oxyhydroxide	0.578	715
HNO <sub>3</sub> / H <sub>2</sub> O <sub>2</sub>	Residual	3.23	5970

All analyte results are reported in µg/g on a dry weight (DW) basis

U = Sample concentration is less than the estimated Method Detection Limit (eMDL)

J = Sample concentration is between the eMDL and the Reporting Limit (RL)

Arsenic Sequential Extraction Results for ARI  
SDG: A141217S2  
Contact: Cheronne Oreiro

Date: January 20, 2015  
Report Generated by: Ben Wozniak  
Applied Speciation and Consulting, LLC

**Quality Control Summary - Total Solids Results**

<b>Sample ID</b>	<b>Total Solids</b>	<b>Units</b>
14-27414-ZP16F	91.4	%
14-27415-ZP16G	67.7	%
14-27416-ZP16H	82.4	%
14-27447-ZO83H	77.2	%
14-27448-ZO83I	57.6	%
14-27449-ZO83J	70.8	%
14-27450-ZO83K	90.3	%
14-27451-ZO83L	89.5	%

Arsenic Sequential Extraction Results for ARI  
 SDG: A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

**Quality Control Summary - Preparation Blank Summary**

Analyte	Units	Fraction	PB1	PB2	PB3	PB4	Mean	StdDev	eMDL*	RL
As	µg/g DW	0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	0.028	0.042	0.033	0.025	0.032	0.007	0.022	0.40
As	µg/g DW	0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	0.054	0.032	0.017	0.028	0.033	0.015	0.046	0.40
As	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	0.039	0.031	0.049	0.083	0.051	0.023	0.069	0.40
As	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> + 0.1M Ascorbic acid	0.057	0.064	0.082	0.097	0.075	0.018	0.054	0.40
As	µg/g DW	HNO <sub>3</sub> /H <sub>2</sub> O <sub>2</sub>	0.06	0.09	0.05	0.25	0.11	0.09	0.28	0.50
Fe	µg/g DW	0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	0.9	0.7	-0.2	0.5	0.5	0.5	1.5	16
Fe	µg/g DW	0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	-0.14	-0.26	0.06	0.09	-0.06	0.17	0.50	16
Fe	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	1.6	1.8	2.8	2.9	2.3	0.7	2.0	16
Fe	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> + 0.1M Ascorbic acid	2.5	3.4	3.6	2.8	3.1	0.5	1.5	16
Fe	µg/g DW	HNO <sub>3</sub> /H <sub>2</sub> O <sub>2</sub>	0.4	-0.2	-0.1	1.6	0.4	0.8	2.5	20

eMDL = Estimated Method Detection Limit; RL = Reporting Limit  
 \*Please see narrative regarding eMDL calculations

Arsenic Sequential Extraction Results for ARI  
 SDG A141217S2  
 Contact: Cheronne Oreiro

Date: January 20, 2015  
 Report Generated by: Ben Wozniak  
 Applied Speciation and Consulting, LLC

**Quality Control Summary - Matrix Duplicate (MD)**

Analyte	Units	Fraction	Sample ID	Rep 1	Rep 2	Mean	RPD
As	µg/g DW	0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	14-27416-ZP16H	1.934	2.210	2.072	13.3
As	µg/g DW	0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	14-27416-ZP16H	3.749	3.744	3.746	0.1
As	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	14-27416-ZP16H	0.790	0.732	0.761	7.7
As	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> + 0.1M Ascorbic acid	14-27416-ZP16H	0.845	0.763	0.804	10.2
As	µg/g DW	HNO <sub>3</sub> /H <sub>2</sub> O <sub>2</sub>	14-27416-ZP16H	35.38	41.84	38.61	16.7
Fe	µg/g DW	0.05 M (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	14-27416-ZP16H	36.8	42.3	39.6	13.8
Fe	µg/g DW	0.05 M (NH <sub>4</sub> )H <sub>2</sub> PO <sub>4</sub>	14-27416-ZP16H	391.7	415.9	403.8	6.0
Fe	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	14-27416-ZP16H	1171	1480	1326	23.3
Fe	µg/g DW	0.2M (NH <sub>4</sub> ) <sub>2</sub> C <sub>2</sub> O <sub>4</sub> + 0.1M Ascorbic acid	14-27416-ZP16H	834.5	955.8	895.1	13.6
Fe	µg/g DW	HNO <sub>3</sub> /H <sub>2</sub> O <sub>2</sub>	14-27416-ZP16H	13810	13900	13860	0.6
Total Solids	%	-	14-27416-ZP16H	82.40	84.64	83.52	2.7

NC = Value was not calculated due to one or more concentrations below the eMDL





Laboratory: Applied Speciation & Consulting  
Lab Contact: Russell Gerads  
Lab Address: 18804 Northcreek Parkway  
Bothell, WA 98011  
Phone: 425-483-3300  
Fax: 425-483-9818

ARI Client: Geoengineers  
Project ID: Gas Works Park-Play Area Investigat  
ARI PM: Cheronne Oreiro  
Phone: 206-695-6214  
Fax: 206-695-6201  
Email: subdata@arilabs.com

Analytical Protocol: In-house  
Special Instructions:

Requested Turn Around:  
Email Results (Y/N): Yes

**Limits of Liability.** Subcontractor is expected to perform all requested services in accordance with appropriate methodology following Standard Operating Procedures that meet standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the negotiated amount for said services. The agreement by the Subcontractor to perform services requested by ARI releases ARI from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Subcontractor.

ARI ID	Client ID/ Add'l ID	Sampled	Matrix	Bottles	Analyses
C 14-27414-ZP16F	PAI-10-31.5-32.0	12/11/14 12:30	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As,Fe					
C 14-27415-ZP16G	PAI-11-12.0-12.5	12/11/14 13:50	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As,Fe					
C 14-27416-ZP16H	PAI-11-22-22.5	12/11/14 14:20	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As,Fe					

Carrier		Airbill		Date	
Relinquished by	Company	Date	Time		
<i>[Signature]</i>	ARI	12/17/14	1330		
Received by	Company	Date	Time		
<i>[Signature]</i>	ACS	12/17/14	13:54		

0.302



Laboratory: Applied Speciation & Consulting  
Lab Contact: Russell Gerads  
Lab Address: 18804 Northcreek Parkway  
Bothell, WA 98011  
Phone: 425-483-3300  
Fax: 425-483-9818

ARI Client: Geoenineers  
Project ID: Gas Works Park-Play Area Investigat  
ARI PM: Cheronne Oreirc  
Phone: 206-695-6214  
Fax: 206-695-6201  
Email: subdata@arilabs.com

Analytical Protocol: In-house  
Special Instructions:

Requested Turn Around:  
Email Results (Y/N): **email**

**Limits of Liability.** Subcontractor is expected to perform all requested services in accordance with appropriate methodology following Standard Operating Procedures that meet standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the negotiated amount for said services. The agreement by the Subcontractor to perform services requested by ARI releases ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Subcontractor.

ARI ID	Client ID/ Add'l ID	Sampled	Matrix	Bottles	Analyses
14-27440-Z083A	PAI-11GW	12/12/14 10:00	Water	1	Metals (Sub)
Special Instructions: Speciated As 6800					
14-27441-Z083B	PAI-12GW	12/12/14 12:30	Water	1	Metals (Sub)
Special Instructions: Speciated As 6800					
14-27442-Z083C	PAI-2GW	12/12/14 14:00	Water	1	Metals (Sub)
Special Instructions: Speciated As 6800					
14-27447-Z083H	PAI-12-8.5-9.0	12/12/14 09:55	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As, Fe					
14-27448-Z083I	PAI-12-13.5-14.0	12/12/14 10:15	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As, Fe					
14-27449-Z083J	PAI-2-17.5-18.0	12/12/14 12:00	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As, Fe					
14-27450-Z083K	PAI-2-19.0-19.5	12/12/14 11:50	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As, Fe					

Sample not included on CDC. *12/17/14*  
c PAI-10610

Carrier	Airbill	Date
Relinquished by <i>[Signature]</i>	Company <b>ARI</b>	Date <i>12/17/14</i> Time <i>13:30</i>
Received by <i>[Signature]</i>	Company <b>ACS</b>	Date <i>12/17/14</i> Time <i>13:34</i>

@ 0.3x



Laboratory: Applied Speciation & Consulting  
Lab Contact: Russell Gerads

ARI Client: Geoengineers  
Project ID: 0186-846-01 Task 1520

ARI Sample ID	Client Sample ID/ Add'l Sample ID	Sampled	Matrix	Bottles	Analyses
14-27451-Z083L	PAI-3-33.5-34.0	12/12/14 14:30	Soil	1	Metals (Sub)
Special Instructions: Sequential Extraction As, Fe					

Carrier		Airbill	Date	
Relinquished by <i>[Signature]</i>	Company <i>ARI</i>	Date <i>12/17/14</i>	Time <i>1330</i>	
Received by <i>[Signature]</i>	Company <i>ACS</i>	Date <i>12/12/14</i>	Time <i>13:54</i>	

@ 0.3<sup>2</sup>



Laboratory: Applied Speciation & Consulting  
 Lab Contact: Russell Gerads  
 Lab Address: 18804 Northcreek Parkway  
 Bothell, WA 98011  
 Phone: 425-483-3300  
 Fax: 425-483-9818

ARI Client: Geoengineers  
 Project ID: Gas Works Park-Paly Area Investigat  
 ARI PM: Cheronne Oreiro  
 Phone: 206-695-6214  
 Fax: 206-695-6201  
 Email: subdata@arilabs.com

Analytical Protocol: In-house  
 Special Instructions:

Requested Turn Around:  
 Email Results (Y/N): **email**

*Limits of Liability. Subcontractor is expected to perform all requested services in accordance with appropriate methodology following Standard Operating Procedures that meet standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the negotiated amount for said services. The agreement by the Subcontractor to perform services requested by ARI releases ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Subcontractor.*

ARI ID	Client ID/ Add'l ID	Sampled	Matrix	Bottles	Analyses
C 14-27518-ZP15A	MW-36D-141215	12/15/14 10:00	Water	1	Metals (Sub)
Special Instructions: Arsenic Speciation 6800					
C 14-27519-ZP15B	MW-36S-141215	12/15/14 11:30	Water	1	Metals (Sub)
Special Instructions: Arsenic Speciation 6800					

Carrier		Airbill		Date	
Relinquished by	Company	Date	Time		
	ARI	12/17/14	1330		
Received by	Company	Date	Time		
	ACS	12/17/14	13:34		

@0.32





**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants

January 23, 2015

Zanna Satterwhite  
GeoEngineers, Inc.  
Plaza 600 Building  
600 Stewart Street, Suite 1700  
Seattle, WA 98101

**RE: Client Project: Gas Works Park-Play Area Investigation, 0186-846-01 Task 1520**  
**ARI Job No.: ZP35**

Dear Zanna:

Please find enclosed the Chain of Custody records (COCs), sample receipt documentation, and the final data package for samples from the project referenced above.

Sample receipt and details of these analyses are discussed in the Case Narrative.

An electronic copy of this package will remain on file with ARI. Should you have any questions or problems, please feel free to contact me at your convenience.

Sincerely,

ANALYTICAL RESOURCES, INC.

A handwritten signature in black ink, appearing to read "Cheronne Oreiro".

Cheronne Oreiro  
Project Manager  
(206) 695-6214  
[cheronneo@arilabs.com](mailto:cheronneo@arilabs.com)  
[www.arilabs.com](http://www.arilabs.com)

cc: eFile: ZP35

Enclosures

## Chain of Custody Documentation

ARI Job ID: ZP35

# Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number: \_\_\_\_\_ Turn-around Requested: \_\_\_\_\_ Page: 1 of 2

ARI Client Company: Geo Engineers Phone: 206-239-3231 Ice Present?

Client Contact: Zanna Satterwhite No. of Coolers: \_\_\_\_\_ Cooler Temps: \_\_\_\_\_

Client Project Name: Bushwicks Park - Play Area Investigation

Client Project #: D186-846-01 Samplers: Robert Mijahira & Claudia DeLava

Analytical Resources, Incorporated  
Analytical Chemists and Consultants  
4611 South 134th Place, Suite 100  
Tukwila, WA 98168  
206-695-6200 206-695-6201 (fax)  
www.arilabs.com



Sample ID	Date	Time	Matrix	No. Containers	Analysis Requested				Notes/Comments
					Hold				
PAI-6-5.5-6.0	12/9	0910	Soil	6	X				
PAI-6-9.0-9.5	12/9	0905	Soil	6	X				
PAI-7-5.5-6.0	12/9	1000	Soil	6	X				
PAI-7-8.0-8.5	12/9	1010	Soil	6	X				
PAI-7-15.0-15.5	12/9	1100	Soil	6	X				
PAI-7-18.5-19.0	12/9	1110	Soil	6	X				
PAI-7-10.0-10.5	12/9	1040	Soil	6	X				
PAI-7-22.5-23.0	12/9	1145	Soil	6	X				
PAI-2-4.5-5.0	12/9	1300	Soil	6	X				
PAI-2-12.5-13.0	12/9	1400	Soil	4	X				
Comments/Special Instructions	Relinquished by: <u>[Signature]</u> (Signature) Printed Name: <u>Claudia DeLava</u> Company: <u>Geo Engineers</u> Date & Time: <u>12/10/14 1315</u>				Relinquished by: <u>[Signature]</u> (Signature) Printed Name: <u>Rich Nelson</u> Company: <u>ARI</u> Date & Time: <u>12/10/14 1315</u>				Received by: <u>[Signature]</u> (Signature) Printed Name: _____ Company: _____ Date & Time: _____

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.







AS per 17-10-17

Analytical Resources Incorporated  
1000 West 10th Street  
Vancouver, BC V6H 2G6  
Tel: 604-270-7700  
Fax: 604-270-7701



### Chain of Custody Record & Laboratory Analysis Request

LAB Agreement Number <b>1000</b>	Sub-Product Requested <b>Standard</b>	Page <b>1</b>	of <b>2</b>
ABI Client Company <b>Geo Energy Inc</b>	Phone <b>506-234-3031</b>	Date <b>12/1/14</b>	Time <b>15:00</b>
Client Contact <b>Z. ...</b>	Client Project Name <b>Bankfield's Park - Northern Investigation</b>	Client Address <b>1000 West 10th Street</b>	Client City <b>Vancouver</b>

Sample ID	Date	Time	Name	Accession	Analysis	Signature
PAT-6-55-60	12/1	09:10	SA	6		[Signature]
PAT-6-90-95	12/1	09:05	SA	6		[Signature]
PAT-7-55-60	12/1	10:00	SA	6		[Signature]
PAT-7-80-85	12/1	10:00	SA	6		[Signature]
PAT-7-150-155	12/1	10:00	SA	6		[Signature]
PAT-7-185-190	12/1	10:00	SA	6		[Signature]
PAT-7-190-195	12/1	10:00	SA	6		[Signature]
PAT-7-205-210	12/1	10:00	SA	6		[Signature]
PAT-7-45-50	12/1	12:00	SA	6		[Signature]
PAT-7-125-130	12/1	12:00	SA	6		[Signature]









# Cooler Receipt Form

ARI Client: GeoEngineer

Project Name: Gas Works Park

COC No(s): \_\_\_\_\_ (NA)

Delivered by: Fed-Ex UPS Courier Hand Delivered Other: \_\_\_\_\_

Assigned ARI Job No: Z0108

Tracking No: \_\_\_\_\_ (NA)

**Preliminary Examination Phase:**

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES (YES) NO (NO)

Were custody papers included with the cooler? ..... YES (YES) NO

Were custody papers properly filled out (ink, signed, etc.) ..... YES (YES) NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry)

Time: 1:10

4.0

If cooler temperature is out of compliance fill out form 00070F

Temp Gun ID#: 90877952

Cooler Accepted by: \_\_\_\_\_ Date: 12/10/14 Time: 1315

*Complete custody forms and attach all shipping documents*

**Log-In Phase:**

Was a temperature blank included in the cooler? ..... YES (NO) NO

What kind of packing material was used? ... Bubble Wrap Wet Ice (Gel Packs) Baggies (Foam Block) Paper Other: Box

Was sufficient ice used (if appropriate)? ..... NA YES (YES) NO

Were all bottles sealed in individual plastic bags? ..... YES (NO) NO

Did all bottles arrive in good condition (unbroken)? ..... YES (YES) NO

Were all bottle labels complete and legible? ..... YES (YES) NO

Did the number of containers listed on COC match with the number of containers received? ..... YES (YES) NO

Did all bottle labels and tags agree with custody papers? ..... YES (YES) NO

Were all bottles used correct for the requested analyses? ..... YES (YES) NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... (NA) YES NO

Were all VOC vials free of air bubbles? ..... (NA) YES NO

Was sufficient amount of sample sent in each bottle? ..... YES (YES) NO

Date VOC Trip Blank was made at ARI..... NA (Date)

Was Sample Split by ARI : (NA) YES Date/Time: \_\_\_\_\_ Equipment: \_\_\_\_\_ Split by: \_\_\_\_\_

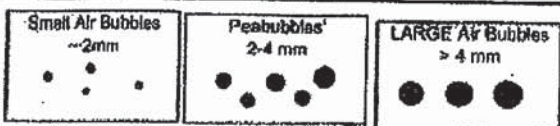
Samples Logged by: AN Date: 12/12/14 Time: 1040

**\*\* Notify Project Manager of discrepancies or concerns \*\***

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

**Additional Notes, Discrepancies, & Resolutions:**

By: \_\_\_\_\_ Date: \_\_\_\_\_



- Small → "sm" (< 2 mm)
- Peabubbles → "pb" (2 to < 4 mm)
- Large → "lg" (4 to < 6 mm)
- Headspace → "hs" (> 6 mm)



# Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number: \_\_\_\_\_ Turn-around Requested: \_\_\_\_\_  
 ARI Client Company: \_\_\_\_\_ Phone: \_\_\_\_\_  
 Client Contact: \_\_\_\_\_  
 Client Project Name: \_\_\_\_\_  
 Client Project #: \_\_\_\_\_  
 Analytical Resources, Incorporated  
 Analytical Chemists and Consultants  
 4611 South 134th Place, Suite 100  
 Tukwila, WA 98168  
 206-695-6200 206-695-6201 (fax)  
 www.arilabs.com



Page: 1 of 2  
 Date: 12/11/14  
 Ice Present?   
 Cooler Temps: \_\_\_\_\_  
 No. of Coolers: \_\_\_\_\_

Sample ID	Date	Time	Matrix	No. Containers	Analysis Requested		Notes/Comments
					Retinquired by:	Received by:	
PAI-9-3-3.5	12/11	0850	Soil	6			
PAI-9-9.5-18.0	12/11	0910	Soil	6			
PAI-9-11.5-12.0	12/11	0920	Soil	6			
PAI-9-12.5-13.0	12/11	0930	Soil	6			
PAI-9-12.5-13-DUP	12/11	0930	Soil	6			
PAI-9-17.5-18.0	12/11	0945	Soil	6			
RINSTATE-1412011	12/11	1010	Water	3			
PAI-10-9.5-10.0	12/11	1145	Soil	6			
PAI-10-14.5-15.0	12/11	1150	Soil	6			
PAI-10-19.5-20.0	12/11	1210	Soil	6			
Comments/Special Instructions					Retinquired by: (Signature) _____	Received by: (Signature) _____	
					Printed Name: _____	Printed Name: _____	
					Company: _____	Company: _____	
					Date & Time: _____	Date & Time: _____	

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.





**Subject:** RE: Gas Works COCs  
**From:** "Dan M. Baker" <dbaker@geoengineers.com>  
**Date:** 12/13/2014 12:54 PM  
**To:** "Zanna A. Satterwhite" <zsatterwhite@geoengineers.com>, Cheronne Oreiro <cheronneo@arilabs.com>  
**CC:** Sample Login <sample-receiving@arilabs.com>

Cheronne and sample log-in folks,

Attached are the revised chain of custody forms.

**Dan M. Baker**  
**Hydrogeologist/Principal | GeoEngineers, Inc.**  
**Phone:** 206.239.3232  
**Fax:** 206.728.2732  
**Mobile:** 206.930.9180  
**Email:** [dbaker@geoengineers.com](mailto:dbaker@geoengineers.com)

---

**From:** Zanna A. Satterwhite  
**Sent:** Saturday, December 13, 2014 11:58 AM  
**To:** Cheronne Oreiro  
**Cc:** Sample Login; Dan M. Baker  
**Subject:** RE: Gas Works COCs

Cheronne,  
We haven't forgotten the COCs for Gas Works. Dan is working on them.  
Thank you  
Zanna

---

**From:** Zanna A. Satterwhite  
**Sent:** Friday, December 12, 2014 6:07 PM  
**To:** 'Cheronne Oreiro'  
**Cc:** Sample Login; Dan M. Baker  
**Subject:** RE: Gas Works COCs

Okay, thanks. I'm ccing Dan.

---

**From:** Cheronne Oreiro [<mailto:cheronneo@arilabs.com>]  
**Sent:** Friday, December 12, 2014 6:06 PM  
**To:** Zanna A. Satterwhite  
**Cc:** Sample Login  
**Subject:** Gas Works COCs

Hi Zanna,

Please cc: [sample-receiving@arilabs.com](mailto:sample-receiving@arilabs.com) when you email the Gas Works COCs.  
Thanks!













# Cooler Receipt Form

ARI Client: GeoEngineers

Project Name: Gas Works Park - PAI

COC No(s): \_\_\_\_\_ (NA)

Delivered by: Fed-Ex UPS Courier Hand Delivered Other: \_\_\_\_\_

Assigned ARI Job No: ZP16

Tracking No: \_\_\_\_\_ NA

**Preliminary Examination Phase:**

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES  NO

Were custody papers included with the cooler? YES  NO

Were custody papers properly filled out (ink, signed, etc.) YES  NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry)  
Time: 11:33 (6.0) -18 S.S 416

If cooler temperature is out of compliance fill out form 00070F Temp Gun ID#: 90877952

Cooler Accepted by: A Date: 12/11/14 Time: 1525

**Complete custody forms and attach all shipping documents**

**Log-In Phase:**

Was a temperature blank included in the cooler? YES  NO

What kind of packing material was used? ... Bubble Wrap  Wet Ice  Gel Packs  Baggies  Foam Block Paper Other: DOX

Was sufficient ice used (if appropriate)? NA  YES  NO

Were all bottles sealed in individual plastic bags? YES  NO

Did all bottles arrive in good condition (unbroken)? YES  NO

Were all bottle labels complete and legible? YES  NO

Did the number of containers listed on COC match with the number of containers received? YES  NO

Did all bottle labels and tags agree with custody papers? YES  NO

Were all bottles used correct for the requested analyses? YES  NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA  YES  NO

Were all VOC vials free of air bubbles? NA  YES  NO

Was sufficient amount of sample sent in each bottle? YES  NO

Date VOC Trip Blank was made at ARI: NA 12/10/14

Was Sample Split by ARI: (NA) YES Date/Time: \_\_\_\_\_ Equipment: \_\_\_\_\_ Split by: \_\_\_\_\_

Samples Logged by: A Date: 12/15/14 Time: 1455

**\*\* Notify Project Manager of discrepancies or concerns \*\***

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

**Additional Notes, Discrepancies, & Resolutions:**  
PAI-11-22-22.5 both 8 & 4oz jars have cracked bottoms. Sulfur jar replaced, no volume lost.

By: A Date: 12/15/14

<b>Small Air Bubbles</b> ~2mm	<b>Peabubbles</b> 2-4 mm	<b>LARGE Air Bubbles</b> > 4 mm	Small → "sm" (< 2 mm)
			Peabubbles → "pb" (2 to 4 mm)
			Large → "lg" (4 to < 6 mm)
			Headspace → "hs" (> 6 mm)



# Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number: **2083** Turn-around Requested: \_\_\_\_\_ of \_\_\_\_\_  
 ARI Client Company: **GeoEngineers** Phone: **206-239-3231**  
 Client Contact: **Zanna Satterwhite**  
 Client Project Name: **Play Area Investigation**  
 Client Project #: **0186-846-01** Samplers: **Robert + Claudette Hannah**

Page: **1** of **1**  
 Date: **12/12/14**  
 No. of Coolers: \_\_\_\_\_  
 Ice Present? \_\_\_\_\_  
 Cooler Temps: \_\_\_\_\_

Analytical Resources, Incorporated  
 Analytical Chemists and Consultants  
 4611 South 134th Place, Suite 100  
 Tukwila, WA 98168  
 206-695-6200 206-695-6201 (fax)  
 www.arilabs.com



Sample ID	Date	Time	Matrix	No. Containers	Analysis Requested		Notes/Comments
PAI-11GW	12/12	1000	Water	13+0			Perous Iron
PAI-12-8.5-9.0	12/12	0955	Soil	3			↑ dry see cores (bottom cap is black)
PAI-12-13.5-14.0	12/12	1015	Soil	3			
PAI-12-19.5-20.0	12/12	1040	Soil	1			
PAI-2-17.5-18.0	12/12	1150	Soil	3			
PAI-2-19.0-19.5	12/12	1200	Soil	3			
PAI-12GW	12/12	1230	Water	13+1			Perous Iron
PAI-2GW	12/12	1400	Water	13+1			Perous Iron
PAI-3-33.5-34.0	12/12	1430	Soil	3			
TRIP BLANK			Water	3			
Comments/Special Instructions							
Relinquished by: <i>[Signature]</i> Printed Name: <b>Hannah McDonough</b> Company: <b>GeoEngineers</b>				Received by: <i>[Signature]</i> Printed Name: <b>C. OREIRO</b> Company: <b>ARI</b>			
Date & Time: <b>12/12/14 @ 1645</b>				Date & Time: <b>12/12/14 1045</b>			

**Limits of Liability:** ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

**Sample Retention Policy:** All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

**Subject:** RE: Gas Works COCs  
**From:** "Dan M. Baker" <dbaker@geoengineers.com>  
**Date:** 12/13/2014 12:54 PM  
**To:** "Zanna A. Satterwhite" <zsatterwhite@geoengineers.com>, Cheronne Oreiro <cheronneo@arilabs.com>  
**CC:** Sample Login <sample-receiving@arilabs.com>

Cheronne and sample log-in folks,

Attached are the revised chain of custody forms.

**Dan M. Baker**  
**Hydrogeologist/Principal | GeoEngineers, Inc.**  
**Phone:** 206.239.3232  
**Fax:** 206.728.2732  
**Mobile:** 206.930.9180  
**Email:** [dbaker@geoengineers.com](mailto:dbaker@geoengineers.com)

---

**From:** Zanna A. Satterwhite  
**Sent:** Saturday, December 13, 2014 11:58 AM  
**To:** Cheronne Oreiro  
**Cc:** Sample Login; Dan M. Baker  
**Subject:** RE: Gas Works COCs

Cheronne,  
We haven't forgotten the COCs for Gas Works. Dan is working on them.  
Thank you  
Zanna

---

**From:** Zanna A. Satterwhite  
**Sent:** Friday, December 12, 2014 6:07 PM  
**To:** 'Cheronne Oreiro'  
**Cc:** Sample Login; Dan M. Baker  
**Subject:** RE: Gas Works COCs

Okay, thanks. I'm ccing Dan.

---

**From:** Cheronne Oreiro [<mailto:cheronneo@arilabs.com>]  
**Sent:** Friday, December 12, 2014 6:06 PM  
**To:** Zanna A. Satterwhite  
**Cc:** Sample Login  
**Subject:** Gas Works COCs

Hi Zanna,

Please cc: [sample-receiving@arilabs.com](mailto:sample-receiving@arilabs.com) when you email the Gas Works COCs.  
Thanks!



ZAS REVISIONS 12-12-14

Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number: **206-239-323**  
 ARI Client Company: **State Street**  
 Client Contract: **206-239-323**  
 Client Project Name: **State Street**  
 Client Project #: **0186-846-01 1520**  
 Client Project Location: **Bay Area Investigation**  
 Client Project Description: **Robert + Charlotte Amtrak**

ARI Assigned Number: **206-239-323**  
 ARI Client Company: **State Street**  
 Client Contract: **206-239-323**  
 Client Project Name: **State Street**  
 Client Project #: **0186-846-01 1520**  
 Client Project Location: **Bay Area Investigation**  
 Client Project Description: **Robert + Charlotte Amtrak**

Sample ID	Date	Time	Matrix	Volume	Method	Notes	Notes/Comments
PAI-1160	12/12	1000	SOIL	3	AS-1000		
PAI-12-85-90	12/12	0955	SOIL	3	AS-1000		
PAI-12-13.5-140	12/12	0915	SOIL	3	AS-1000		
PAI-12-19.5-200	12/12	1040	SOIL	3	AS-1000		
PAI-2-17.5-180	12/12	1150	SOIL	3	AS-1000		
PAI-2-19.0-19.5	12/12	1200	SOIL	3	AS-1000		
PAI-1260	12/12	1230	Blank	3	AS-1000		
PAI-260	12/12	1400	Blank	3	AS-1000		
PAI-3-33.5-340	12/12	1430	SOIL	3	AS-1000		
Imp Blank			Blank	3	AS-1000		

ARI Analytical Resources, Incorporated  
 Analytical Chemists and Consultants  
 4617 South 134th Place, Suite 100  
 Tukwila, WA 98168  
 206-495-6200, 206-495-6201 fax  
 www.ariab.com

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 206-495-6200, 206-495-6201 fax  
 www.ariab.com





# Cooler Receipt Form

ARI Client: GeoEngineers  
 COC No(s): \_\_\_\_\_ (NA)  
 Assigned ARI Job No: 7083

Project Name: GasWorks Park - PAI  
 Delivered by: Fed-Ex UPS Courier (Hand Delivered) Other: \_\_\_\_\_  
 Tracking No: \_\_\_\_\_ (NA)

**Preliminary Examination Phase:**

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES (NO)  
 Were custody papers included with the cooler? YES NO  
 Were custody papers properly filled out (ink, signed, etc.) YES NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry)  
 Time: 10:45 10:29.1 11:3  
 If cooler temperature is out of compliance fill out form 00070F Temp Gun ID#: 70877952

Cooler Accepted by: CO (N) Date: 12/12/14 Time: 10:45

Complete custody forms and attach all shipping documents

**Log-In Phase:**

Was a temperature blank included in the cooler? YES (NO)  
 What kind of packing material was used? ... Bubble Wrap (Wet Ice) Gel Packs Baggies Foam Block Paper Other: BOX  
 Was sufficient ice used (if appropriate)? NA YES (NO)  
 Were all bottles sealed in individual plastic bags? YES (NO)  
 Did all bottles arrive in good condition (unbroken)? YES (NO)  
 Were all bottle labels complete and legible? YES (NO)  
 Did the number of containers listed on COC match with the number of containers received? YES (NO)  
 Did all bottle labels and tags agree with custody papers? YES (NO)  
 Were all bottles used correct for the requested analyses? YES (NO)  
 Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA YES (NO)  
 Were all VOC vials free of air bubbles? NA YES (NO)  
 Was sufficient amount of sample sent in each bottle? YES (NO)  
 Date VOC Trip Blank was made at ARI... NA 12/11/14  
 Was Sample Split by ARI: (NA) YES Date/Time: \_\_\_\_\_ Equipment: \_\_\_\_\_ Split by: \_\_\_\_\_

Samples Logged by: AV Date: 12/15/14 Time: 11:25

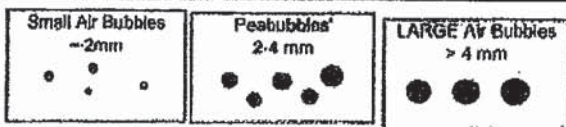
\*\* Notify Project Manager of discrepancies or concerns \*\*

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

**Additional Notes, Discrepancies, & Resolutions:**

PAI-12GW has 13 bottles PAI-11GW has 14 bottles  
 NO UNFILTERED CONV VOLUME ON 6W PAI-1GW EXCEPT SULFIDE  
 NO UNFILTERED CONV VOLUME FOR PAI-12GW 11GW + 2GW  
 FF SULFIDES placed on hold for

By: \_\_\_\_\_ Date: \_\_\_\_\_



Small → "sm" (< 2 mm)  
 Peabubbles → "pb" (2 to < 4 mm)  
 Large → "lg" (4 to < 6 mm)  
 Headspace → "hs" (> 6 mm)

NO UNFILTERED CONV VOLUME FOR PAI-2GW EXCEPT FOR SULFIDE

Case Narrative, Data Qualifiers, Control Limits

ARI Job ID: ZP35





## Case Narrative

**Client: GeoEngineers, Inc.**

**Project: Gas Works Park-Play Area Investigation, 0186-846-01 Task 1520**

**ARI Job No.: ZP35**

### Sample Receipt

Five soil samples were removed from archive on December 16, 2014 and logged under ARI job ZP35. The samples were analyzed for BTEX, SIM PAHs, and arsenic, as requested. For details regarding sample receipt, please refer to the Cooler Receipt Form.

Volume for sample **PAI-2-18.0-18.5** was received in a core sleeve. BTEX volume was removed from the core first. Remaining volume was split between the extraction laboratory and the metals laboratory.

### BTEX by SW8260C

The samples were analyzed within the method recommended holding times.

Initial and continuing calibrations were within method requirements. Internal standard areas were within limits.

The surrogate percent recoveries were within control limits.

The method blank was clean at the reporting limit. The LCS and LCSD percent recoveries were within control limits.

### PAHs by SW8270-SIM

The samples were extracted and analyzed within method recommended holding times.

Initial and continuing calibrations were within method requirements. Internal standard areas were within limits.

The surrogate percent recoveries were within control limits.

The method blank was clean at the reporting limit. The LCS percent recoveries were within control limits.

The matrix spike/matrix spike duplicate percent recoveries of Naphthalene were outside the control limits for sample **PAI-10-29.5-30.0**. The matrix spike duplicate percent recoveries of Fluorene and Phenanthrene were also outside the control limits. No corrective action is required for matrix QC.



**Arsenic by SW6010C**

The samples and associated laboratory QC were digested and analyzed within recommended holding times.

The method blank was clean at the reporting limit. The LCS percent recovery was within control limits.

The matrix spike percent recovery and the replicate RPD were within control limits.

# Sample ID Cross Reference Report



ARI Job No: ZP35  
Client: Geoengineers  
Project Event: 0186-846-01 Task 1570  
Project Name: Gas Works Park-Play Area Investigat

Sample ID	ARI Lab ID	ARI LIMS ID	Matrix	Sample Date/Time	VTSR
1. PAI-2-12.5-13.0	ZP35A	14-27575	Soil	12/09/14 14:00	12/12/14 16:45
2. PAI-2-18.0-18.5	ZP35B	14-27576	Soil	12/09/14 14:40	12/12/14 16:45
3. PAI-9-9.5-10.0	ZP35C	14-27577	Soil	12/11/14 09:10	12/12/14 16:45
4. PAI-10-29.5-30.0	ZP35D	14-27578	Soil	12/11/14 12:20	12/12/14 16:45
5. PAI-12-19.5-20.0	ZP35E	14-27579	Soil	12/12/14 10:40	12/12/14 16:45





## Data Reporting Qualifiers

Effective 12/31/13

### Inorganic Data

- U Indicates that the target analyte was not detected at the reported concentration
- \* Duplicate RPD is not within established control limits
- B Reported value is less than the CRDL but  $\geq$  the Reporting Limit
- N Matrix Spike recovery not within established control limits
- NA Not Applicable, analyte not spiked
- H The natural concentration of the spiked element is so much greater than the concentration spiked that an accurate determination of spike recovery is not possible
- L Analyte concentration is  $\leq 5$  times the Reporting Limit and the replicate control limit defaults to  $\pm 1$  RL instead of the normal 20% RPD

### Organic Data

- U Indicates that the target analyte was not detected at the reported concentration
- \* Flagged value is not within established control limits
- B Analyte detected in an associated Method Blank at a concentration greater than one-half of ARI's Reporting Limit or 5% of the regulatory limit or 5% of the analyte concentration in the sample.
- J Estimated concentration when the value is less than ARI's established reporting limits
- D The spiked compound was not detected due to sample extract dilution
- E Estimated concentration calculated for an analyte response above the valid instrument calibration range. A dilution is required to obtain an accurate quantification of the analyte.



- Q Indicates a detected analyte with an initial or continuing calibration that does not meet established acceptance criteria (<20%RSD, <20%Drift or minimum RRF).
- S Indicates an analyte response that has saturated the detector. The calculated concentration is not valid; a dilution is required to obtain valid quantification of the analyte
- NA The flagged analyte was not analyzed for
- NR Spiked compound recovery is not reported due to chromatographic interference
- NS The flagged analyte was not spiked into the sample
- M Estimated value for an analyte detected and confirmed by an analyst but with low spectral match parameters. This flag is used only for GC-MS analyses
- N The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification"
- Y The analyte is not detected at or above the reported concentration. The reporting limit is raised due to chromatographic interference. The Y flag is equivalent to the U flag with a raised reporting limit.
- EMPC Estimated Maximum Possible Concentration (EMPC) defined in EPA Statement of Work DLM02.2 as a value "calculated for 2,3,7,8-substituted isomers for which the quantitation and /or confirmation ion(s) has signal to noise in excess of 2.5, but does not meet identification criteria" **(Dioxin/Furan analysis only)**
- C The analyte was positively identified on only one of two chromatographic columns. Chromatographic interference prevented a positive identification on the second column
- P The analyte was detected on both chromatographic columns but the quantified values differ by  $\geq 40\%$  RPD with no obvious chromatographic interference
- X Analyte signal includes interference from polychlorinated diphenyl ethers. **(Dioxin/Furan analysis only)**
- Z Analyte signal includes interference from the sample matrix or perfluorokerosene ions. **(Dioxin/Furan analysis only)**



## **Geotechnical Data**

- A The total of all fines fractions. This flag is used to report total fines when only sieve analysis is requested and balances total grain size with sample weight.
- F Samples were frozen prior to particle size determination
- SM Sample matrix was not appropriate for the requested analysis. This normally refers to samples contaminated with an organic product that interferes with the sieving process and/or moisture content, porosity and saturation calculations
- SS Sample did not contain the proportion of "fines" required to perform the pipette portion of the grain size analysis
- W Weight of sample in some pipette aliquots was below the level required for accurate weighting



## Analytical Method Information

Analyte	MDL	Reporting Limit	Surrogate %R	Duplicate RPD	Matrix Spike %R	RPD	Blank Spike / LCS %R	RPD
<b>8260C VOA Solid in Solid (EPA 8260C)</b>								
Preservation: NaHSO <sub>4</sub> , MeOH, Cool <6°C								
Container: VOA Vial, Clear, 40 mL, Amount Required: 15 g Hold Time: 14 days								
Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub>								
Dichlorodifluoromethane	0.207	1.00 ug/kg		30	67 - 142	30	67 - 142	30
Chloromethane	0.263	1.00 ug/kg		30	65 - 129	30	65 - 129	30
Vinyl Chloride	0.235	1.00 ug/kg		30	74 - 134	30	74 - 134	30
Bromomethane	0.187	1.00 ug/kg		30	40 - 172	30	40 - 172	30
Chloroethane	0.462	1.00 ug/kg		30	53 - 154	30	53 - 154	30
Trichlorofluoromethane	0.266	1.00 ug/kg		30	57 - 161	30	57 - 161	30
Acrolein	3.81	50.0 ug/kg		30	60 - 130	30	60 - 130	30
Acetone	0.482	5.00 ug/kg		30	48 - 132	30	48 - 132	30
1,1,2-Trichloro-1,2,2-Trifluoroethane	0.287	2.00 ug/kg		30	72 - 142	30	72 - 142	30
1,1-Dichloroethene	0.336	1.00 ug/kg		30	73 - 138	30	73 - 138	30
Bromoethane	0.440	2.00 ug/kg		30	74 - 132	30	74 - 132	30
Iodomethane	0.215	1.00 ug/kg		30	34 - 181	30	34 - 181	30
Methylene Chloride	0.635	2.00 ug/kg		30	61 - 128	30	61 - 128	30
Carbon Disulfide	0.559	1.00 ug/kg		30	72 - 146	30	72 - 146	30
Acrylonitrile	1.03	5.00 ug/kg		30	59 - 124	30	59 - 124	30
Methyl tert-butyl Ether	0.231	1.00 ug/kg		30	68 - 124	30	68 - 124	30
trans-1,2-Dichloroethene	0.266	1.00 ug/kg		30	73 - 131	30	73 - 131	30
Vinyl Acetate	0.381	5.00 ug/kg		30	54 - 138	30	54 - 138	30
1,1-Dichloroethane	0.203	1.00 ug/kg		30	65 - 139	30	65 - 139	30
2-Butanone	0.513	5.00 ug/kg		30	64 - 120	30	64 - 120	30
2,2-Dichloropropane	0.292	1.00 ug/kg		30	77 - 137	30	77 - 137	30
cis-1,2-Dichloroethene	0.240	1.00 ug/kg		30	75 - 124	30	75 - 124	30
Chloroform	0.234	1.00 ug/kg		30	75 - 126	30	75 - 126	30
Bromochloromethane	0.323	1.00 ug/kg		30	69 - 133	30	69 - 133	30
1,1,1-Trichloroethane	0.226	1.00 ug/kg		30	78 - 133	30	78 - 133	30
1,1-Dichloropropene	0.312	1.00 ug/kg		30	80 - 123	30	80 - 123	30
Carbon tetrachloride	0.213	1.00 ug/kg		30	76 - 136	30	76 - 136	30
1,2-Dichloroethane	0.191	1.00 ug/kg		30	77 - 120	30	77 - 120	30
Benzene	0.296	1.00 ug/kg		30	80 - 120	30	80 - 120	30
Trichloroethene	0.212	1.00 ug/kg		30	80 - 120	30	80 - 120	30
1,2-Dichloropropane	0.162	1.00 ug/kg		30	74 - 120	30	74 - 120	30
Bromodichloromethane	0.254	1.00 ug/kg		30	80 - 122	30	80 - 122	30
Dibromomethane	0.147	1.00 ug/kg		30	80 - 120	30	80 - 120	30
2-Chloroethyl vinyl ether	0.276	5.00 ug/kg		30	20 - 157	30	20 - 157	30
4-Methyl-2-Pentanone	0.420	5.00 ug/kg		30	70 - 124	30	70 - 124	30
cis-1,3-Dichloropropene	0.226	1.00 ug/kg		30	80 - 124	30	80 - 124	30
Toluene	0.151	1.00 ug/kg		30	78 - 120	30	78 - 120	30
trans-1,3-Dichloropropene	0.216	1.00 ug/kg		30	80 - 126	30	80 - 126	30
1,1,2-Trichloroethane	0.286	1.00 ug/kg		30	77 - 120	30	77 - 120	30
1,2-Dibromoethane	0.176	1.00 ug/kg		30	79 - 120	30	79 - 120	30
2-Hexanone	0.439	5.00 ug/kg		30	62 - 128	30	62 - 128	30
1,3-Dichloropropane	0.209	1.00 ug/kg		30	77 - 120	30	77 - 120	30
Tetrachloroethene	0.257	1.00 ug/kg		30	76 - 131	30	76 - 131	30
Dibromochloromethane	0.266	1.00 ug/kg		30	77 - 123	30	77 - 123	30
Chlorobenzene	0.219	1.00 ug/kg		30	80 - 120	30	80 - 120	30
1,1,1,2-Tetrachloroethane	0.233	1.00 ug/kg		30	80 - 120	30	80 - 120	30

## Analytical Method Information

Analyte	MDL	Reporting	Surrogate	Duplicate	Matrix Spike		Blank Spike / LCS	
		Limit	%R	RPD	%R	RPD	%R	RPD
Ethylbenzene	0.202	1.00 ug/kg		30	80 - 120	30	80 - 120	30
m,p-Xylene	0.392	1.00 ug/kg		30	80 - 123	30	80 - 123	30
o-Xylene	0.224	1.00 ug/kg		30	80 - 120	30	80 - 120	30
Styrene	0.138	1.00 ug/kg		30	80 - 122	30	80 - 122	30
Bromoform	0.297	1.00 ug/kg		30	63 - 120	30	63 - 120	30
Isopropyl Benzene	0.233	1.00 ug/kg		30	77 - 127	30	77 - 127	30
1,1,2,2-Tetrachloroethane	0.253	1.00 ug/kg		30	71 - 120	30	71 - 120	30
1,2,3-Trichloropropane	0.517	2.00 ug/kg		30	75 - 120	30	75 - 120	30
trans-1,4-Dichloro 2-Butene	0.437	5.00 ug/kg		30	62 - 127	30	62 - 127	30
n-Propylbenzene	0.272	1.00 ug/kg		30	76 - 126	30	76 - 126	30
Bromobenzene	0.153	1.00 ug/kg		30	75 - 120	30	75 - 120	30
1,3,5-Trimethylbenzene	0.254	1.00 ug/kg		30	77 - 126	30	77 - 126	30
2-Chlorotoluene	0.300	1.00 ug/kg		30	76 - 120	30	76 - 120	30
4-Chlorotoluene	0.277	1.00 ug/kg		30	75 - 121	30	75 - 121	30
t-Butylbenzene	0.306	1.00 ug/kg		30	77 - 125	30	77 - 125	30
1,2,4-Trimethylbenzene	0.230	1.00 ug/kg		30	77 - 125	30	77 - 125	30
s-Butylbenzene	0.240	1.00 ug/kg		30	77 - 127	30	77 - 127	30
4-Isopropyl Toluene	0.236	1.00 ug/kg		30	78 - 131	30	78 - 131	30
1,3-Dichlorobenzene	0.227	1.00 ug/kg		30	76 - 120	30	76 - 120	30
1,4-Dichlorobenzene	0.232	1.00 ug/kg		30	75 - 120	30	75 - 120	30
n-Butylbenzene	0.262	1.00 ug/kg		30	75 - 134	30	75 - 134	30
1,2-Dichlorobenzene	0.293	1.00 ug/kg		30	77 - 120	30	77 - 120	30
1,2-Dibromo-3-Chloropropane	0.586	5.00 ug/kg		30	61 - 128	30	61 - 128	30
1,2,4-Trichlorobenzene	0.332	5.00 ug/kg		30	75 - 130	30	75 - 130	30
Hexachloro-1,3-Butadiene	0.410	5.00 ug/kg		30	72 - 135	30	72 - 135	30
Naphthalene	0.429	5.00 ug/kg		30	71 - 122	30	71 - 122	30
1,2,3-Trichlorobenzene	0.305	5.00 ug/kg		30	76 - 122	30	76 - 122	30
surr: 1,2-Dichloroethane-d4			80 - 149					
surr: 1,2-Dichlorobenzene-d4			80 - 120					
surr: Toluene-d8			77 - 120					
surr: 4-Bromofluorobenzene			80 - 120					
surr: Dibromofluoromethane			80 - 120					
Pentafluorobenzene								
Chlorobenzene-d5								
1,4-Difluorobenzene								
1,4-Dichlorobenzene-d4								



## Analytical Method Information

Analyte	MDL	Reporting	Surrogate	Duplicate	Matrix Spike		Blank Spike / LCS	
		Limit	%R	RPD	%R	RPD	%R	RPD
<b>8270D-SIM PAH (5 ug/kg) in Solid (EPA 8270D-SIM)</b>								
Preservation: Cool <6°C								
Container: Glass WM, Clear, 8 oz								
Amount Required: 150 g								
Hold Time: 14 days								
Naphthalene	2.26	5.00 ug/kg		30	36 - 120	30	36 - 120	30
2-Methylnaphthalene	1.69	5.00 ug/kg		30	35 - 120	30	35 - 120	30
1-Methylnaphthalene	1.61	5.00 ug/kg		30	39 - 120	30	39 - 120	30
Biphenyl	1.44	5.00 ug/kg		30	30 - 160	30	30 - 160	30
2,6-Dimethylnaphthalene	0.750	5.00 ug/kg		30	30 - 160	30	30 - 160	30
Acenaphthylene	1.61	5.00 ug/kg		30	35 - 120	30	35 - 120	30
Acenaphthene	1.49	5.00 ug/kg		30	39 - 120	30	39 - 120	30
Dibenzofuran	1.41	5.00 ug/kg		30	38 - 120	30	38 - 120	30
2,3,5-Trimethylnaphthalene	0.419	5.00 ug/kg		30		30		30
Fluorene	1.47	5.00 ug/kg		30	41 - 120	30	41 - 120	30
Dibenzothiophene	0.425	5.00 ug/kg		30		30		30
Phenanthrene	1.58	5.00 ug/kg		30	46 - 120	30	46 - 120	30
Anthracene	1.78	5.00 ug/kg		30	36 - 120	30	36 - 120	30
Carbazole	0.189	5.00 ug/kg		30	30 - 160	30	30 - 160	30
1-Methylphenanthrene	0.700	5.00 ug/kg		30	30 - 160	30	30 - 160	30
Fluoranthene	1.87	5.00 ug/kg		30	46 - 120	30	46 - 120	30
Pyrene	2.26	5.00 ug/kg		30	49 - 120	30	49 - 120	30
Benzo(a)anthracene	2.22	5.00 ug/kg		30	42 - 120	30	42 - 120	30
Chrysene	1.92	5.00 ug/kg		30	48 - 120	30	48 - 120	30
Benzo(b)fluoranthene	2.11	5.00 ug/kg		30	35 - 127	30	35 - 127	30
Benzo(k)fluoranthene	2.28	5.00 ug/kg		30	37 - 129	30	37 - 129	30
Benzo(j)fluoranthene	1.75	5.00 ug/kg		30	40 - 120	30	40 - 120	30
Benzo(e)pyrene	0.647	5.00 ug/kg		30	30 - 160	30	30 - 160	30
Benzo(a)pyrene	2.38	5.00 ug/kg		30	36 - 120	30	36 - 120	30
Perylene	3.56	5.00 ug/kg		30	44 - 120	30	44 - 120	30
Indeno(1,2,3-cd)pyrene	3.01	5.00 ug/kg		30	40 - 120	30	40 - 120	30
Dibenzo(a,h)anthracene	2.56	5.00 ug/kg		30	38 - 120	30	38 - 120	30
Benzo(g,h,i)perylene	2.79	5.00 ug/kg		30	38 - 120	30	38 - 120	30
surr: 2-Methylnaphthalene-d10								32 - 120
surr: Dibenzo[a,h]anthracene-d14								21 - 133
surr: Fluoranthene-d10								36 - 134
Naphthalene-d8								
Acenaphthene-d10								
Phenanthrene-d10								
Chrysene-d12								
Perylene-d12								



## Analytical Method Information

Analyte	MDL	Reporting	Surrogate	Duplicate	Matrix Spike		Blank Spike / LCS	
		Limit	%R	RPD	%R	RPD	%R	RPD
<b>Met 6010C in Solid (EPA 6010C)</b>								
Preservation: Cool <6°C								
Container: Glass WM, Clear, 4 oz								
Amount Required: 100 g								
Hold Time: 180 days								
Aluminum	0.757	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Antimony	0.628	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Arsenic	0.333	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Barium	0.133	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Beryllium	0.0160	0.100 mg/kg		20	75 - 125	20	80 - 120	20
Boron	0.739	2.00 mg/kg		20	75 - 125	20	80 - 120	20
Cadmium	0.0180	0.200 mg/kg		20	75 - 125	20	80 - 120	20
Calcium	1.13	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Chromium	0.124	0.500 mg/kg		20	75 - 125	20	80 - 120	20
Cobalt	0.0270	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Copper	0.0920	0.200 mg/kg		20	75 - 125	20	80 - 120	20
Iron	0.750	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Lead	0.155	2.00 mg/kg		20	75 - 125	20	80 - 120	20
Magnesium	0.961	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Manganese	0.0280	0.100 mg/kg		20	75 - 125	20	80 - 120	20
Molybdenum	0.0790	0.500 mg/kg		20	75 - 125	20	80 - 120	20
Nickel	0.386	1.00 mg/kg		20	75 - 125	20	80 - 120	20
Potassium	6.57	50.0 mg/kg		20	75 - 125	20	80 - 120	20
Selenium	0.499	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Silver	0.0430	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Sodium	1.14	50.0 mg/kg		20	75 - 125	20	80 - 120	20
Sodium-1	114	5000 mg/kg		20	75 - 125	20	80 - 120	20
Strontium	0.00900	0.100 mg/kg		20	75 - 125	20	80 - 120	20
Thallium	0.310	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Tin	0.141	1.00 mg/kg		20	75 - 125	20	80 - 120	20
Titanium	0.211	0.500 mg/kg		20	75 - 125	20	80 - 120	20
Vanadium	0.0270	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Zinc	0.145	1.00 mg/kg		20	75 - 125	20	80 - 120	20

Volatile Analysis  
Report and Summary QC Forms

ARI Job ID: ZP35

**ORGANICS ANALYSIS DATA SHEET**

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 1

Sample ID: PAI-2-18.0-18.5  
SAMPLE

Lab Sample ID: ZP35B  
LIMS ID: 14-27576  
Matrix: Soil  
Data Release Authorized: *MW*  
Reported: 01/05/15

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: 12/09/14  
Date Received: 12/12/14

Instrument/Analyst: NT5/PAB  
Date Analyzed: 12/19/14 19:09

Sample Amount: 30.6 mg-dry-wt  
Purge Volume: 5.0 mL  
Moisture: 30.8%

CAS Number	Analyte	DL	LOQ	Result
71-43-2	Benzene	48	160	11,000
108-88-3	Toluene	25	160	7,100
100-41-4	Ethylbenzene	33	160	14,000
179601-23-1	m,p-Xylene	64	160	11,000
95-47-6	o-Xylene	37	160	6,400

Reported in µg/kg (ppb)

**Volatile Surrogate Recovery**

d4-1,2-Dichloroethane	93.0%
d8-Toluene	104%
Bromofluorobenzene	99.8%
d4-1,2-Dichlorobenzene	100%

Results corrected for soil moisture content per Section 11.10.5 of EPA Method 8000C.



**ORGANICS ANALYSIS DATA SHEET**

Volatiles by Purge & Trap GC/MS-Method SW8260C

Page 1 of 1

Sample ID: PAI-10-29.5-30.0

SAMPLE

Lab Sample ID: ZP35D

LIMS ID: 14-27578

Matrix: Soil

Data Release Authorized: *MMW*

Reported: 01/05/15

QC Report No: ZP35-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1570

Date Sampled: 12/11/14

Date Received: 12/12/14

Instrument/Analyst: NT5/PAB

Date Analyzed: 12/19/14 19:34

Sample Amount: 81.1 mg-dry-wt

Purge Volume: 5.0 mL

Moisture: 7.8%

CAS Number	Analyte	DL	LOQ	Result
71-43-2	Benzene	18	62	670
108-88-3	Toluene	9.3	62	180
100-41-4	Ethylbenzene	12	62	260
179601-23-1	m,p-Xylene	24	62	380
95-47-6	o-Xylene	14	62	120

Reported in µg/kg (ppb)

**Volatile Surrogate Recovery**

d4-1,2-Dichloroethane	89.6%
d8-Toluene	105%
Bromofluorobenzene	98.8%
d4-1,2-Dichlorobenzene	97.9%

Results corrected for soil moisture content per Section 11.10.5 of EPA Method 8000C.

**ORGANICS ANALYSIS DATA SHEET**

Volatiles by Purge & Trap GC/MS-Method SW8260C

Sample ID: PAI-12-19.5-20.0

Page 1 of 1

SAMPLE

Lab Sample ID: ZP35E

QC Report No: ZP35-Geoengineers

LIMS ID: 14-27579

Project: Gas Works Park-Play Area Investigat

Matrix: Soil

0186-846-01 Task 1570

Data Release Authorized: *MW*

Date Sampled: 12/12/14

Reported: 01/05/15

Date Received: 12/12/14

Instrument/Analyst: NT5/PAB

Sample Amount: 2.81 mg-dry-wt

Date Analyzed: 12/19/14 19:59

Purge Volume: 5.0 mL

Moisture: 28.4%

CAS Number	Analyte	DL	LOQ	Result
71-43-2	Benzene	530	1800	2,000
108-88-3	Toluene	270	1800	1,900
100-41-4	Ethylbenzene	360	1800	2,700
179601-23-1	m,p-Xylene	700	1800	6,300
95-47-6	o-Xylene	400	1800	4,600

Reported in µg/kg (ppb)

**Volatile Surrogate Recovery**

d4-1,2-Dichloroethane	92.3%
d8-Toluene	103%
Bromofluorobenzene	98.6%
d4-1,2-Dichlorobenzene	99.6%

Results corrected for soil moisture content per Section 11.10.5 of EPA Method 8000C.

VOA SURROGATE RECOVERY SUMMARY



Matrix: Soil

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570

ARI ID	Client ID	Level	DCE	TOL	BFB	DCB	TOT OUT
MB-121914A	Method Blank	Med	95.3%	104%	97.2%	99.1%	0
LCS-121914A	Lab Control	Med	96.6%	107%	99.4%	99.2%	0
LCSD-121914A	Lab Control Dup	Med	99.7%	102%	101%	99.9%	0
ZP35B	PAI-2-18.0-18.5	Med	93.0%	104%	99.8%	100%	0
ZP35D	PAI-10-29.5-30.0	Med	89.6%	105%	98.8%	97.9%	0
ZP35E	PAI-12-19.5-20.0	Med	92.3%	103%	98.6%	99.6%	0

SW8260C	LCS/MB LIMITS		QC LIMITS	
	Low	Med	Low	Med
(DCE) = d4-1,2-Dichloroethane	80-149	80-124	80-149	80-124
(TOL) = d8-Toluene	77-120	80-120	77-120	80-120
(BFB) = Bromofluorobenzene	80-120	80-120	80-120	80-120
(DCB) = d4-1,2-Dichlorobenzene	80-120	80-120	80-120	80-120

Log Number Range: 14-27576 to 14-27579



**ORGANICS ANALYSIS DATA SHEET**

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 1

Sample ID: LCS-121914A  
LAB CONTROL SAMPLE

Lab Sample ID: LCS-121914A  
LIMS ID: 14-27576  
Matrix: Soil  
Data Release Authorized: *MW*  
Reported: 01/05/15

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: NA  
Date Received: NA

Instrument/Analyst LCS: NT5/PAB  
LCSD: NT5/PAB  
Date Analyzed LCS: 12/19/14 10:05  
LCSD: 12/19/14 10:30

Sample Amount LCS: 100 mg-dry-wt  
LCSD: 100 mg-dry-wt  
Purge Volume LCS: 5.0 mL  
LCSD: 5.0 mL  
Moisture: NA

Analyte	LCS	Spike Added-LCS	LCS Recovery	LCSD	Spike Added-LCSD	LCSD Recovery	RPD
Benzene	2460	2500	98.4%	2390	2500	95.6%	2.9%
Toluene	2770	2500	111%	2590	2500	104%	6.7%
Ethylbenzene	2550	2500	102%	2500	2500	100%	2.0%
m,p-Xylene	5100	5000	102%	4950	5000	99.0%	3.0%
o-Xylene	2540	2500	102%	2460	2500	98.4%	3.2%

Reported in µg/kg (ppb)

RPD calculated using sample concentrations per SW846.

**Volatile Surrogate Recovery**

	LCS	LCSD
d4-1,2-Dichloroethane	96.6%	99.7%
d8-Toluene	107%	102%
Bromofluorobenzene	99.4%	101%
d4-1,2-Dichlorobenzene	99.2%	99.9%

4A  
VOLATILE METHOD BLANK SUMMARY

Method Blank ID.

MB1219

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY ARE

Lab File ID: MB1219

Lab Sample ID: MB1219

Date Analyzed: 12/19/14

Time Analyzed: 1054

Instrument ID: NT5

Heated Purge: (Y/N) Y

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
	=====	=====	=====	=====
01	LCS1219	LCS1219	LCS1219	1005
02	LCS1219	LCS1219	LCS1219A	1030
03	PAI-2-18.0-1	ZP35B	ZP35B	1909
04	PAI-10-29.5-	ZP35D	ZP35D	1934
05	PAI-12-19.5-	ZP35E	ZP35E	1959
06				
07				
08				
09				
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30				

COMMENTS:

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ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C  
Page 1 of 1

Sample ID: MB-121914A  
METHOD BLANK

Lab Sample ID: MB-121914A  
LIMS ID: 14-27576  
Matrix: Soil  
Data Release Authorized: *mmw*  
Reported: 01/05/15

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: NA  
Date Received: NA

Instrument/Analyst: NT5/PAB  
Date Analyzed: 12/19/14 10:54

Sample Amount: 100 mg-dry-wt  
Purge Volume: 5.0 mL  
Moisture: NA

CAS Number	Analyte	DL	LOQ	Result
71-43-2	Benzene	15	50	< 50 U
108-88-3	Toluene	7.6	50	< 50 U
100-41-4	Ethylbenzene	10	50	< 50 U
179601-23-1	m,p-Xylene	20	50	< 50 U
95-47-6	o-Xylene	11	50	< 50 U

Reported in µg/kg (ppb)

**Volatile Surrogate Recovery**

d4-1,2-Dichloroethane	95.3%
d8-Toluene	104%
Bromofluorobenzene	97.2%
d4-1,2-Dichlorobenzene	99.1%



5A  
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC      Contract: GEOENGINEERS  
 Lab Code: ARI                      Case No.: GAS WORKS PARK-PLAY AREA      SDG No.: ZP35  
 Lab File ID: BFB1120X                                      BFB Injection Date: 11/20/14  
 Instrument ID: NT5    BFB Injection Time: 1217  
 GC Column: RTXVMS              ID: 0.18 (mm)                      Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	18.5
75	30.0 - 60.0% of mass 95	46.4
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.1
173	Less than 2.0% of mass 174	0.1 ( 0.2)1
174	50.0 - 100.0% of mass 95	77.3
175	5.0 - 9.0% of mass 174	5.9 ( 7.6)1
176	95.0 - 101.0% of mass 174	76.1 ( 98.5)1
177	5.0 - 9.0% of mass 176	5.2 ( 6.8)2

1-Value is % mass 174                                      2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	1	SCL0002-CAL1	0011120	11/20/14	1321
02	2	SCL0002-CAL2	0021120	11/20/14	1346
03	5	SCL0002-CAL3	0051120	11/20/14	1410
04	10	SCL0002-CAL4	0101120	11/20/14	1435
05	50	SCL0002-CAL5	0501120	11/20/14	1500
06	100	SCL0002-CAL6	1001120	11/20/14	1525
07	150	SCL0002-CAL7	1501120	11/20/14	1550
08	200	SCL0002-CAL8	2001120	11/20/14	1614
09	ICV1120	SCL0002-SCV1	ICV1120	11/20/14	1639
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					





FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

LAB FILE ID: RF1: 0011120

RF2: 0021120

RF5: 0051120

RF10: 0101120

RF50: 0501120

COMPOUND	RF1	RF2	RF5	RF10	RF50
Chloromethane	0.949	1.008	0.921	1.124	1.051
Vinyl Chloride	0.867	0.917	0.810	1.046	1.016
Bromomethane	0.494	0.469	0.373	0.467	0.349
Chloroethane	0.578	0.777	0.578	0.685	0.600
Trichlorofluoromethane	1.172	1.211	1.065	1.204	1.146
Acrolein	0.135	0.143	0.127	0.148	0.142
1,1,1-Trichloroethane	0.684	0.665	0.612	0.675	0.605
Acetone	0.269	0.238	0.176	0.199	0.180
1,1-Dichloroethene	0.651	0.636	0.590	0.682	0.601
Bromoethane	0.491	0.452	0.394	0.505	0.406
Iodomethane	0.165	0.143	0.148	0.267	0.384
Methylene Chloride		0.954	0.756	0.805	0.712
Acrylonitrile	0.255	0.288	0.282	0.305	0.302
Carbon Disulfide	2.309	2.143	1.980	2.280	1.989
Trans-1,2-Dichloroethene	0.747	0.804	0.745	0.804	0.778
Vinyl Acetate	0.363	0.366	0.366	0.362	0.364
1,1-Dichloroethane	1.416	1.460	1.412	1.505	1.466
2-Butanone	0.076	0.075	0.081	0.088	0.082
2,2-Dichloropropane	1.057	1.097	1.090	1.162	1.135
Cis-1,2-Dichloroethene	0.797	0.805	0.814	0.830	0.803
Chloroform	1.279	1.258	1.206	1.284	1.231
Bromochloromethane	0.319	0.372	0.366	0.356	0.350
1,1,1-Trichloroethane	1.148	1.152	1.134	1.201	1.162
1,1-Dichloropropene	0.419	0.427	0.429	0.433	0.415
Carbon Tetrachloride	0.392	0.401	0.378	0.407	0.394
1,2-Dichloroethane	0.339	0.393	0.377	0.392	0.364
Benzene	1.171	1.263	1.207	1.275	1.157
Trichloroethene	0.307	0.314	0.306	0.330	0.298
1,2-Dichloropropane	0.350	0.309	0.329	0.334	0.318
Bromodichloromethane	0.362	0.382	0.375	0.391	0.371
Dibromomethane	0.151	0.158	0.163	0.163	0.156
2-Chloroethyl Vinyl Ether	0.163	0.184	0.189	0.184	0.186
4-Methyl-2-Pentanone	0.110	0.140	0.130	0.139	0.142
Cis 1,3-dichloropropene	0.435	0.481	0.455	0.478	0.490
Toluene	0.795	0.958	0.793	0.811	0.897
Trans 1,3-Dichloropropene	0.382	0.385	0.388	0.417	0.440
2-Hexanone	0.180	0.187	0.193	0.208	0.201

FORM VI VOA

ZP35: 00041



FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

LAB FILE ID: RF1: 0011120

RF2: 0021120

RF5: 0051120

RF10: 0101120

RF50: 0501120

COMPOUND	RF1	RF2	RF5	RF10	RF50
1,1,2-Trichloroethane	0.210	0.271	0.240	0.279	0.269
1,3-Dichloropropane	0.373	0.391	0.369	0.425	0.397
Tetrachloroethene	0.315	0.301	0.267	0.308	0.283
Chlorodibromomethane	0.205	0.234	0.219	0.253	0.245
1,2-Dibromoethane	0.216	0.253	0.246	0.258	0.254
Chlorobenzene	0.774	0.815	0.818	0.844	0.790
Ethyl Benzene	1.466	1.465	1.389	1.522	1.396
1,1,1,2-Tetrachloroethane	0.246	0.260	0.253	0.265	0.261
m,p-xylene	0.598	0.584	0.559	0.591	0.553
o-Xylene	0.566	0.540	0.534	0.552	0.532
Styrene	0.881	0.841	0.851	0.915	0.859
Bromoform	0.267	0.263	0.248	0.272	0.280
1,1,2,2-Tetrachloroethane	0.466	0.534	0.519	0.534	0.526
1,2,3-Trichloropropane	0.164	0.176	0.173	0.190	0.170
Trans-1,4-Dichloro 2-Butene	0.133	0.142	0.128	0.150	0.152
N-Propyl Benzene	3.163	3.265	3.103	3.166	2.974
Bromobenzene	0.592	0.621	0.599	0.608	0.576
Isopropyl Benzene	2.783	2.756	2.710	2.785	2.659
2-Chloro Toluene	1.916	1.893	1.744	1.842	1.746
4-Chloro Toluene	1.932	1.999	1.885	1.885	1.832
T-Butyl Benzene	2.049	2.022	1.994	2.055	1.930
1,3,5-Trimethyl Benzene	2.325	2.267	2.224	2.315	2.195
1,2,4-Trimethylbenzene	2.245	2.277	2.261	2.292	2.179
S-Butyl Benzene	3.059	3.086	2.987	3.069	2.905
4-Isopropyl Toluene	2.404	2.503	2.435	2.480	2.384
1,3-Dichlorobenzene	1.229	1.250	1.188	1.197	1.138
1,4-Dichlorobenzene	1.372	1.342	1.237	1.214	1.158
N-Butyl Benzene	2.406	2.262	2.301	2.316	2.177
1,2-Dichlorobenzene	1.254	1.162	1.117	1.140	1.076
1,2-Dibromo 3-Chloropropane	0.086	0.075	0.086	0.090	0.096
1,2,4-Trichlorobenzene	0.958	0.809	0.751	0.760	0.718
Hexachloro 1,3-Butadiene	0.413	0.414	0.387	0.408	0.376
Naphthalene	2.262	1.943	1.927	1.946	1.839
1,2,3-Trichlorobenzene	0.793	0.726	0.716	0.701	0.667
Dichlorodifluoromethane	0.557	0.510	0.610	0.833	0.781
Methyl tert butyl ether	1.960	1.988	2.006	2.102	2.122

FORM VI VOA

ZP35: 00042

FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

LAB FILE ID: RF1: 0011120

RF2: 0021120

RF5: 0051120

RF10: 0101120

RF50: 0501120

COMPOUND	RF1	RF2	RF5	RF10	RF50
d4-1,2-Dichloroethane	0.726	0.738	0.728	0.735	0.750
d8-Toluene	1.240	1.408	1.252	1.247	1.338
4-Bromofluorobenzene	0.529	0.508	0.498	0.510	0.514
d4-1,2-Dichlorobenzene	0.954	0.952	0.940	0.933	0.951
Dibromofluoromethane	0.687	0.658	0.677	0.682	0.701

FORM VI VOA

ZP35: 00043



FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

LAB FILE ID: RF100: 1001120

RF150: 1501120

RF200: 2001120

COMPOUND	RF100	RF150	RF200
Chloromethane	1.043	1.006	1.011
Vinyl Chloride	0.975	0.965	0.940
Bromomethane	0.307	0.306	0.298
Chloroethane	0.571	0.556	0.538
Trichlorofluoromethane	1.067	1.034	0.976
Acrolein	0.118	0.112	0.103
1,1,2-Trichloro-2,2-Trifluoroethane	0.593	0.610	0.564
Acetone	0.146		
1,1-Dichloroethene	0.592	0.597	0.550
Bromoethane	0.366	0.361	0.336
Iodomethane	0.415	0.433	0.395
Methylene Chloride	0.552		
Acrylonitrile	0.303	0.282	0.278
Carbon Disulfide	1.956	2.007	1.852
Trans-1,2-Dichloroethene	0.723	0.567	0.539
Vinyl Acetate	0.379	0.365	0.360
1,1-Dichloroethane	1.469	1.419	1.405
2-Butanone	0.085	0.076	0.075
2,2-Dichloropropane	1.130	1.086	1.076
Cis-1,2-Dichloroethene	0.807	0.776	0.784
Chloroform	1.248	1.203	1.192
Bromochloromethane	0.356	0.334	0.334
1,1,1-Trichloroethane	1.188	1.138	1.123
1,1-Dichloropropene	0.429	0.406	0.433
Carbon Tetrachloride	0.398	0.381	0.398
1,2-Dichloroethane	0.377	0.354	0.372
Benzene	1.208	1.146	1.212
Trichloroethene	0.301	0.293	0.299
1,2-Dichloropropane	0.327	0.314	0.325
Bromodichloromethane	0.379	0.366	0.372
Dibromomethane	0.159	0.149	0.154
2-Chloroethyl Vinyl Ether	0.199	0.209	0.197
4-Methyl-2-Pentanone	0.133	0.137	0.140
Cis 1,3-dichloropropene	0.489	0.513	0.475
Toluene	0.817	0.868	0.839
Trans 1,3-Dichloropropene	0.429	0.449	0.456
2-Hexanone	0.213	0.206	0.209

FORM VI VOA

ZP35 : 00044



FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

LAB FILE ID: RF100: 1001120

RF150: 1501120

RF200: 2001120

COMPOUND	RF100	RF150	RF200
1,1,2-Trichloroethane	0.244	0.262	0.264
1,3-Dichloropropane	0.382	0.405	0.412
Tetrachloroethene	0.262	0.289	0.295
Chlorodibromomethane	0.244	0.256	0.262
1,2-Dibromoethane	0.238	0.255	0.261
Chlorobenzene	0.809	0.774	0.793
Ethyl Benzene	1.405	1.339	1.332
1,1,1,2-Tetrachloroethane	0.271	0.267	0.270
m,p-xylene	0.550	0.529	0.524
o-Xylene	0.546	0.530	0.537
Styrene	0.893	0.848	0.868
Bromoform	0.294	0.294	0.296
1,1,2,2-Tetrachloroethane	0.556	0.542	0.558
1,2,3-Trichloropropane	0.176	0.173	0.175
Trans-1,4-Dichloro 2-Butene	0.159	0.160	0.163
N-Propyl Benzene	2.939	2.795	2.767
Bromobenzene	0.589	0.579	0.595
Isopropyl Benzene	2.623	2.549	2.519
2-Chloro Toluene	1.758	1.727	1.735
4-Chloro Toluene	1.822	1.789	1.796
T-Butyl Benzene	1.940	1.870	1.882
1,3,5-Trimethyl Benzene	2.192	2.115	2.119
1,2,4-Trimethylbenzene	2.177	2.092	2.082
S-Butyl Benzene	2.859	2.737	2.696
4-Isopropyl Toluene	2.364	2.273	2.251
1,3-Dichlorobenzene	1.152	1.129	1.149
1,4-Dichlorobenzene	1.157	1.137	1.154
N-Butyl Benzene	2.168	2.086	2.064
1,2-Dichlorobenzene	1.086	1.067	1.095
1,2-Dibromo 3-Chloropropane	0.103	0.102	0.102
1,2,4-Trichlorobenzene	0.741	0.740	0.750
Hexachloro 1,3-Butadiene	0.384	0.384	0.392
Naphthalene	1.956	1.887	1.889
1,2,3-Trichlorobenzene	0.695	0.691	0.707
Dichlorodifluoromethane	0.760	0.743	0.719
Methyl tert butyl ether	2.137	1.776	1.540

FORM VI VOA

ZP35: 00045

FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

LAB FILE ID: RF100: 1001120 RF150: 1501120 RF200: 2001120

COMPOUND	RF100	RF150	RF200
d4-1,2-Dichloroethane	0.746	0.711	0.700
d8-Toluene	1.248	1.311	1.195
4-Bromofluorobenzene	0.515	0.516	0.518
d4-1,2-Dichlorobenzene	0.944	0.946	0.937
Dibromofluoromethane	0.713	0.691	0.686

FORM VI VOA

ZP35 : 00046

FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R <sup>2</sup>
Chloromethane	AVRG	1.014	6.2
Vinyl Chloride	AVRG	0.942	8.2
Bromomethane	LINR		0.9975
Chloroethane	AVRG	0.610	13.2
Trichlorofluoromethane	AVRG	1.109	7.8
Acrolein	AVRG	0.128	12.6
1,1,2-Trichloro-1,2,2-Trifluoroethane	AVRG	0.626	6.9
Acetone	2ORDR		0.9994
1,1-Dichloroethene	AVRG	0.612	6.8
Bromoethane	AVRG	0.414	15.1
Iodomethane	LINR		0.9962
Methylene Chloride	AVRG	0.756	19.3
Acrylonitrile	AVRG	0.287	5.8
Carbon Disulfide	AVRG	2.064	7.9
Trans-1,2-Dichloroethene	AVRG	0.714	14.5
Vinyl Acetate	AVRG	0.366	1.6
1,1-Dichloroethane	AVRG	1.444	2.5
2-Butanone	AVRG	0.080	6.4
2,2-Dichloropropane	AVRG	1.104	3.2
Cis-1,2-Dichloroethene	AVRG	0.802	2.1
Chloroform	AVRG	1.238	2.9
Bromochloromethane	AVRG	0.348	5.2
1,1,1-Trichloroethane	AVRG	1.156	2.3
1,1-Dichloropropene	AVRG	0.424	2.3
Carbon Tetrachloride	AVRG	0.394	2.5
1,2-Dichloroethane	AVRG	0.371	5.0
Benzene	AVRG	1.205	3.9
Trichloroethene	AVRG	0.306	3.8
1,2-Dichloropropane	AVRG	0.326	3.9
Bromodichloromethane	AVRG	0.375	2.4
Dibromomethane	AVRG	0.157	3.3
2-Chloroethyl Vinyl Ether	AVRG	0.189	7.2
4-Methyl-2-Pentanone	AVRG	0.134	7.8
Cis 1,3-dichloropropene	AVRG	0.477	4.9
Toluene	AVRG	0.847	6.8
Trans 1,3-Dichloropropene	AVRG	0.418	7.1
2-Hexanone	AVRG	0.200	5.8

<- Indicates value outside QC limits:  
(%RSD < 20% or R<sup>2</sup> > 0.990)



FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R <sup>2</sup>
1,1,2-Trichloroethane	AVRG	0.255	8.8
1,3-Dichloropropane	AVRG	0.394	4.9
Tetrachloroethene	AVRG	0.290	6.5
Chlorodibromomethane	AVRG	0.240	8.1
1,2-Dibromoethane	AVRG	0.248	5.8
Chlorobenzene	AVRG	0.802	3.0
Ethyl Benzene	AVRG	1.414	4.6
1,1,1,2-Tetrachloroethane	AVRG	0.262	3.4
m,p-xylene	AVRG	0.561	4.9
o-Xylene	AVRG	0.542	2.2
Styrene	AVRG	0.869	2.9
Bromoform	AVRG	0.277	6.2
1,1,2,2-Tetrachloroethane	AVRG	0.529	5.4
1,2,3-Trichloropropane	AVRG	0.175	4.2
Trans-1,4-Dichloro 2-Butene	AVRG	0.148	8.6
N-Propyl Benzene	AVRG	3.022	6.0
Bromobenzene	AVRG	0.595	2.5
Isopropyl Benzene	AVRG	2.673	3.9
2-Chloro Toluene	AVRG	1.795	4.3
4-Chloro Toluene	AVRG	1.868	3.9
T-Butyl Benzene	AVRG	1.968	3.7
1,3,5-Trimethyl Benzene	AVRG	2.219	3.6
1,2,4-Trimethylbenzene	AVRG	2.201	3.7
S-Butyl Benzene	AVRG	2.925	5.2
4-Isopropyl Toluene	AVRG	2.387	3.8
1,3-Dichlorobenzene	AVRG	1.179	3.8
1,4-Dichlorobenzene	AVRG	1.221	7.4
N-Butyl Benzene	AVRG	2.223	5.3
1,2-Dichlorobenzene	AVRG	1.125	5.5
1,2-Dibromo 3-Chloropropane	AVRG	0.092	10.9
1,2,4-Trichlorobenzene	AVRG	0.778	9.9
Hexachloro 1,3-Butadiene	AVRG	0.395	3.7
Naphthalene	AVRG	1.956	6.6
1,2,3-Trichlorobenzene	AVRG	0.712	5.2
Dichlorodifluoromethane	AVRG	0.689	16.8
Methyl tert butyl ether	AVRG	1.954	10.4

<- Indicates value outside QC limits:  
(%RSD < 20% or R<sup>2</sup> > 0.990)

FORM 6  
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Calibration Date: 11/20/14

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R <sup>2</sup>
d4-1,2-Dichloroethane	AVRG	0.729	2.3
d8-Toluene	AVRG	1.280	5.3
4-Bromofluorobenzene	AVRG	0.513	1.8
d4-1,2-Dichlorobenzene	AVRG	0.945	0.8
Dibromofluoromethane	AVRG	0.687	2.4

<- Indicates value outside QC limits:  
(%RSD < 20% or R<sup>2</sup> > 0.990)



## VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Cont. Calib. Date: 11/20/14

Init. Calib. Date: 11/20/14

Cont. Calib. Time: 1410

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
Chloromethane	1.014	1.0512	0.100	AVRG	3.7
Vinyl Chloride	0.942	1.0155	0.010	AVRG	7.8
Bromomethane	50.000	6.138	0.010	LINR	-87.7 <-
Chloroethane	0.610	0.6000	0.010	AVRG	-1.6
Trichlorofluoromethane	1.109	1.1456	0.010	AVRG	3.3
Acrolein	0.128	0.1416	0.010	AVRG	10.6
1,1,2-Trichloro-2,2-Trifluoroethane	0.626	0.6054	0.010	AVRG	-3.3
Acetone	250.00	17.629	0.010	2ORDR	-92.9 <-
1,1-Dichloroethene	0.612	0.6008	0.010	AVRG	-1.8
Bromoethane	0.414	0.4056	0.010	AVRG	-2.0
Iodomethane	50.000	1.810	0.010	LINR	-96.4 <-
Methylene Chloride	0.756	0.7124	0.010	AVRG	-5.8
Acrylonitrile	0.287	0.3016	0.010	AVRG	5.1
Carbon Disulfide	2.064	1.9888	0.010	AVRG	-3.6
Trans-1,2-Dichloroethene	0.713	0.7784	0.010	AVRG	9.2
Vinyl Acetate	0.366	0.3643	0.010	AVRG	-0.5
1,1-Dichloroethane	1.444	1.4664	0.100	AVRG	1.6
2-Butanone	0.080	0.0820	0.010	AVRG	2.5
2,2-Dichloropropane	1.104	1.1350	0.010	AVRG	2.8
Cis-1,2-Dichloroethene	0.802	0.8030	0.010	AVRG	0.1
Chloroform	1.238	1.2306	0.010	AVRG	-0.6
Bromochloromethane	0.348	0.3501	0.010	AVRG	0.6
1,1,1-Trichloroethane	1.156	1.1625	0.010	AVRG	0.6
1,1-Dichloropropene	0.424	0.4152	0.010	AVRG	-2.1
Carbon Tetrachloride	0.394	0.3937	0.010	AVRG	-0.1
1,2-Dichloroethane	0.371	0.3644	0.010	AVRG	-1.8
Benzene	1.205	1.1573	0.010	AVRG	-4.0
Trichloroethene	0.306	0.2980	0.010	AVRG	-2.6
1,2-Dichloropropane	0.326	0.3182	0.010	AVRG	-2.4
Bromodichloromethane	0.375	0.3715	0.010	AVRG	-0.9
Dibromomethane	0.157	0.1562	0.010	AVRG	-0.5
2-Chloroethyl Vinyl Ether	0.189	0.1856	0.010	AVRG	-1.8
4-Methyl-2-Pentanone	0.134	0.1418	0.010	AVRG	5.8
Cis 1,3-dichloropropene	0.477	0.4896	0.010	AVRG	2.6
Toluene	0.847	0.8968	0.010	AVRG	5.9
Trans 1,3-Dichloropropene	0.418	0.4396	0.010	AVRG	5.2
2-Hexanone	0.200	0.2008	0.010	AVRG	0.4

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF



7A  
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Cont. Calib. Date: 11/20/14

Init. Calib. Date: 11/20/14

Cont. Calib. Time: 1410

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
1,1,2-Trichloroethane	0.255	0.2686	0.010	AVRG	5.3
1,3-Dichloropropane	0.394	0.3967	0.010	AVRG	0.7
Tetrachloroethene	0.290	0.2831	0.010	AVRG	-2.4
Chlorodibromomethane	0.240	0.2449	0.010	AVRG	2.0
1,2-Dibromoethane	0.248	0.2544	0.010	AVRG	2.6
Chlorobenzene	0.802	0.7901	0.300	AVRG	-1.5
Ethyl Benzene	1.414	1.3961	0.010	AVRG	-1.3
1,1,1,2-Tetrachloroethane	0.262	0.2612	0.010	AVRG	-0.3
m,p-xylene	0.561	0.5531	0.010	AVRG	-1.4
o-Xylene	0.542	0.5318	0.010	AVRG	-1.9
Styrene	0.870	0.8595	0.010	AVRG	-1.2
Bromoform	0.277	0.2803	0.100	AVRG	1.2
1,1,2,2-Tetrachloroethane	0.529	0.5260	0.300	AVRG	-0.6
1,2,3-Trichloropropane	0.175	0.1703	0.010	AVRG	-2.7
Trans-1,4-Dichloro 2-Butene	0.148	0.1522	0.010	AVRG	2.8
N-Propyl Benzene	3.022	2.9739	0.010	AVRG	-1.6
Bromobenzene	0.595	0.5758	0.010	AVRG	-3.2
Isopropyl Benzene	2.673	2.6587	0.010	AVRG	-0.5
2-Chloro Toluene	1.795	1.7460	0.010	AVRG	-2.7
4-Chloro Toluene	1.868	1.8323	0.010	AVRG	-1.9
T-Butyl Benzene	1.968	1.9304	0.010	AVRG	-1.9
1,3,5-Trimethyl Benzene	2.219	2.1952	0.010	AVRG	-1.1
1,2,4-Trimethylbenzene	2.201	2.1787	0.010	AVRG	-1.0
S-Butyl Benzene	2.925	2.9052	0.010	AVRG	-0.7
4-Isopropyl Toluene	2.387	2.3838	0.010	AVRG	-0.1
1,3-Dichlorobenzene	1.179	1.1379	0.010	AVRG	-3.5
1,4-Dichlorobenzene	1.221	1.1579	0.010	AVRG	-5.2
N-Butyl Benzene	2.222	2.1774	0.010	AVRG	-2.0
1,2-Dichlorobenzene	1.125	1.0764	0.010	AVRG	-4.3
1,2-Dibromo 3-Chloropropane	0.092	0.0965	0.010	AVRG	4.9
1,2,4-Trichlorobenzene	0.778	0.7183	0.010	AVRG	-7.7
Hexachloro 1,3-Butadiene	0.395	0.3761	0.010	AVRG	-4.8
Naphthalene	1.956	1.8389	0.010	AVRG	-6.0
1,2,3-Trichlorobenzene	0.712	0.6667	0.010	AVRG	-6.4
Dichlorodifluoromethane	0.689	0.7813	0.010	AVRG	13.4
Methyl tert butyl ether	1.954	2.1226	0.010	AVRG	8.6

<- Exceeds QC limit of 20% D  
\* RF less than minimum RF

7A  
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Cont. Calib. Date: 11/20/14

Init. Calib. Date: 11/20/14

Cont. Calib. Time: 1410

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
=====	=====	=====	=====	=====	=====
d4-1,2-Dichloroethane_____	0.729	0.7497	0.010	AVRG	2.8
d8-Toluene_____	1.280	1.3384	0.010	AVRG	4.6
4-Bromofluorobenzene_____	0.514	0.5144	0.010	AVRG	0.1
d4-1,2-Dichlorobenzene_____	0.945	0.9515	0.010	AVRG	0.7
Dibromofluoromethane_____	0.687	0.7009	0.010	AVRG	2.0

<- Exceeds QC limit of 20% D

\* RF less than minimum RF



7A  
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC  
ARI Job No: ZP35  
Instrument ID: NT5  
Init. Calib. Date: 11/20/14

Client: GEOENGINEERS  
Project: GAS WORKS PARK-PLAY AREA  
Cont. Calib. Date: 12/19/14  
Cont. Calib. Time: 0928

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift	
Chloromethane	1.014	0.7222	0.100	AVRG	-28.8	<-
Vinyl Chloride	0.942	0.7454	0.010	AVRG	-20.9	<-
Bromomethane	50.000	31.949	0.010	LINR	-36.1	<-
Chloroethane	0.610	0.4814	0.010	AVRG	-21.1	<-
Trichlorofluoromethane	1.109	1.0807	0.010	AVRG	-2.6	
Acrolein	0.128	0.1084	0.010	AVRG	-15.3	
1,1,2-Trichloro-2,2-Trifluoroethane	0.626	0.6233	0.010	AVRG	-0.4	
Acetone	250.00	234.81	0.010	2ORDR	-6.1	
1,1-Dichloroethene	0.612	0.5675	0.010	AVRG	-7.3	
Bromoethane	0.414	0.3999	0.010	AVRG	-3.4	
Iodomethane	50.000	36.813	0.010	LINR	-26.4	<-
Methylene Chloride	0.756	0.6431	0.010	AVRG	-14.9	
Acrylonitrile	0.287	0.2406	0.010	AVRG	-16.2	
Carbon Disulfide	2.064	1.9389	0.010	AVRG	-6.1	
Trans-1,2-Dichloroethene	0.713	0.7065	0.010	AVRG	-0.9	
Vinyl Acetate	0.366	0.3364	0.010	AVRG	-8.1	
1,1-Dichloroethane	1.444	1.3832	0.100	AVRG	-4.2	
2-Butanone	0.080	0.0737	0.010	AVRG	-7.9	
2,2-Dichloropropane	1.104	1.1863	0.010	AVRG	7.4	
Cis-1,2-Dichloroethene	0.802	0.7568	0.010	AVRG	-5.6	
Chloroform	1.238	1.2474	0.010	AVRG	0.8	
Bromochloromethane	0.348	0.3486	0.010	AVRG	0.2	
1,1,1-Trichloroethane	1.156	1.2281	0.010	AVRG	6.2	
1,1-Dichloropropene	0.424	0.4561	0.010	AVRG	7.6	
Carbon Tetrachloride	0.394	0.4757	0.010	AVRG	20.7	<-
1,2-Dichloroethane	0.371	0.4192	0.010	AVRG	13.0	
Benzene	1.205	1.2426	0.010	AVRG	3.1	
Trichloroethene	0.306	0.3193	0.010	AVRG	4.3	
1,2-Dichloropropane	0.326	0.3200	0.010	AVRG	-1.8	
Bromodichloromethane	0.375	0.4167	0.010	AVRG	11.1	
Dibromomethane	0.157	0.1651	0.010	AVRG	5.2	
2-Chloroethyl Vinyl Ether	0.189	0.1472	0.010	AVRG	-22.1	<-
4-Methyl-2-Pentanone	0.134	0.1377	0.010	AVRG	2.8	
Cis 1,3-dichloropropene	0.477	0.5065	0.010	AVRG	6.2	
Toluene	0.847	0.8280	0.010	AVRG	-2.2	
Trans 1,3-Dichloropropene	0.418	0.4842	0.010	AVRG	15.8	
2-Hexanone	0.200	0.1913	0.010	AVRG	-4.4	

<- Exceeds QC limit of 20% D  
\* RF less than minimum RF



7A  
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Cont. Calib. Date: 12/19/14

Init. Calib. Date: 11/20/14

Cont. Calib. Time: 0928

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
1,1,2-Trichloroethane	0.255	0.2801	0.010	AVRG	9.8
1,3-Dichloropropane	0.394	0.4097	0.010	AVRG	4.0
Tetrachloroethene	0.290	0.2997	0.010	AVRG	3.3
Chlorodibromomethane	0.240	0.2766	0.010	AVRG	15.2
1,2-Dibromoethane	0.248	0.2819	0.010	AVRG	13.7
Chlorobenzene	0.802	0.8354	0.300	AVRG	4.2
Ethyl Benzene	1.414	1.4608	0.010	AVRG	3.3
1,1,1,2-Tetrachloroethane	0.262	0.2954	0.010	AVRG	12.7
m,p-xylene	0.561	0.5756	0.010	AVRG	2.6
o-Xylene	0.542	0.5576	0.010	AVRG	2.9
Styrene	0.870	0.9084	0.010	AVRG	4.4
Bromoform	0.277	0.3060	0.100	AVRG	10.5
1,1,2,2-Tetrachloroethane	0.529	0.5281	0.300	AVRG	-0.2
1,2,3-Trichloropropane	0.175	0.1796	0.010	AVRG	2.6
Trans-1,4-Dichloro 2-Butene	0.148	0.1506	0.010	AVRG	1.8
N-Propyl Benzene	3.022	3.0621	0.010	AVRG	1.3
Bromobenzene	0.595	0.6177	0.010	AVRG	3.8
Isopropyl Benzene	2.673	2.7599	0.010	AVRG	3.2
2-Chloro Toluene	1.795	1.8351	0.010	AVRG	2.2
4-Chloro Toluene	1.868	1.9102	0.010	AVRG	2.2
T-Butyl Benzene	1.968	2.0498	0.010	AVRG	4.2
1,3,5-Trimethyl Benzene	2.219	2.3068	0.010	AVRG	4.0
1,2,4-Trimethylbenzene	2.201	2.2826	0.010	AVRG	3.7
S-Butyl Benzene	2.925	3.0088	0.010	AVRG	2.9
4-Isopropyl Toluene	2.387	2.4932	0.010	AVRG	4.4
1,3-Dichlorobenzene	1.179	1.2271	0.010	AVRG	4.1
1,4-Dichlorobenzene	1.221	1.2345	0.010	AVRG	1.1
N-Butyl Benzene	2.222	2.2529	0.010	AVRG	1.4
1,2-Dichlorobenzene	1.125	1.1291	0.010	AVRG	0.4
1,2-Dibromo 3-Chloropropane	0.092	0.0963	0.010	AVRG	4.7
1,2,4-Trichlorobenzene	0.778	0.7634	0.010	AVRG	-1.9
Hexachloro 1,3-Butadiene	0.395	0.4025	0.010	AVRG	1.9
Naphthalene	1.956	1.8658	0.010	AVRG	-4.6
1,2,3-Trichlorobenzene	0.712	0.6929	0.010	AVRG	-2.7
Dichlorodifluoromethane	0.689	0.3525	0.010	AVRG	-48.8
Methyl tert butyl ether	1.954	2.0252	0.010	AVRG	3.6

<- Exceeds QC limit of 20% D  
\* RF less than minimum RF

## VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Instrument ID: NT5

Cont. Calib. Date: 12/19/14

Init. Calib. Date: 11/20/14

Cont. Calib. Time: 0928

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
=====	=====	=====	=====	=====	=====
d4-1,2-Dichloroethane	0.729	0.7189	0.010	AVRG	-1.4
d8-Toluene	1.280	1.2160	0.010	AVRG	-5.0
4-Bromofluorobenzene	0.514	0.5141	0.010	AVRG	0.0
d4-1,2-Dichlorobenzene	0.945	0.9645	0.010	AVRG	2.1
Dibromofluoromethane	0.687	0.6533	0.010	AVRG	-4.9

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF



8A  
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 0101120

Ical Date: 11/20/14

Instrument ID: NT5

Project Run Date: 11/20/14

	IS1 (PFB) AREA #	RT #	IS2 (DFB) AREA #	RT #	IS3 (CLB) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	278101	5.12	699590	5.56	830842	8.00
UPPER LIMIT	556202	5.62	1399180	6.06	1661684	8.50
LOWER LIMIT	139050	4.62	349795	5.06	415421	7.50
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 ICV1120	274481	5.11	695211	5.55	823689	8.00
02						
03						
04						
05						
06						
07						
08						
09						
10						
11						
12						
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						

IS1 (PFB) = Pentafluorobenzene  
 IS2 (DFB) = 1,4-Difluorobenzene  
 IS3 (CLB) = d5-Chlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

\* Values outside of QC limits.



8A  
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY AREA

Ical Midpoint ID: 0101120

Ical Date: 11/20/14

Instrument ID: NT5

Project Run Date: 11/20/14

	IS4 (DCB) AREA #	RT #	AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	450241	10.09				
UPPER LIMIT	900482	10.59				
LOWER LIMIT	225120	9.59				
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 ICV1120	439864	10.10				
02						
03						
04						
05						
06						
07						
08						
09						
10						
11						
12						
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						

IS4 (DCB) = d4-1,4-Dichlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

\* Values outside of QC limits.

8A  
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC  
ARI Job No: ZP35  
Ical Midpoint ID: 0101120  
Instrument ID: NT5

Client: GEOENGINEERS  
Project: GAS WORKS PARK-PLAY AREA  
Ical Date: 11/20/14  
Project Run Date: 12/19/14

	IS1 (PFB) AREA #	RT #	IS2 (DFB) AREA #	RT #	IS3 (CLB) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	278101	5.12	699590	5.56	830842	8.00
UPPER LIMIT	556202	5.62	1399180	6.06	1661684	8.50
LOWER LIMIT	139050	4.62	349795	5.06	415421	7.50
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS1219	337612	5.11	779608	5.55	966808	8.00
02 LCS1219	342947	5.11	793503	5.55	968460	8.00
03 MB1219	325612	5.12	761187	5.56	910835	8.00
04 PAI-2-18.0-1	348239	5.11	795333	5.55	975408	8.00
05 PAI-10-29.5-	357025	5.11	802587	5.55	978386	8.00
06 PAI-12-19.5-	338685	5.11	773757	5.55	927662	8.00
07						
08						
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20						
21						
22						

IS1 (PFB) = Pentafluorobenzene  
IS2 (DFB) = 1,4-Difluorobenzene  
IS3 (CLB) = d5-Chlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint  
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

\* Values outside of QC limits.



8A  
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC  
ARI Job No: ZP35  
Ical Midpoint ID: 0101120  
Instrument ID: NT5

Client: GEOENGINEERS  
Project: GAS WORKS PARK-PLAY AREA  
Ical Date: 11/20/14  
Project Run Date: 12/19/14

	IS4 (DCB) AREA #	RT #	AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	450241	10.09				
UPPER LIMIT	900482	10.59				
LOWER LIMIT	225120	9.59				
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS1219	520111	10.09				
02 LCS1219	529836	10.09				
03 MB1219	484034	10.09				
04 PAI-2-18.0-1	533394	10.10				
05 PAI-10-29.5-	526713	10.10				
06 PAI-12-19.5-	498437	10.09				
07						
08						
09						
10						
11						
12						
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						

IS4 (DCB) = d4-1,4-Dichlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

\* Values outside of QC limits.



SIM PAH Analysis  
Report and Summary QC Forms

ARI Job ID: ZP35

ORGANICS ANALYSIS DATA SHEET  
PNAs by SIM SW8270D-SIM GC/MS  
Extraction Method: SW3546  
Page 1 of 1

Sample ID: PAI-2-18.0-18.5  
SAMPLE

Lab Sample ID: ZP35B  
LIMS ID: 14-27576  
Matrix: Soil  
Data Release Authorized:  
Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: 12/09/14  
Date Received: 12/12/14

Date Extracted: 12/23/14  
Date Analyzed: 01/19/15 22:43  
Instrument/Analyst: NT8/JZ  
GPC Cleanup: No

Sample Amount: 10.4 g-dry-wt  
Final Extract Volume: 0.50 mL  
Dilution Factor: 10.0  
Percent Moisture: 30.8 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	22	48	64,000 ES
208-96-8	Acenaphthylene	15	48	3,600
83-32-9	Acenaphthene	14	48	9,600 E
86-73-7	Fluorene	14	48	8,300 E
85-01-8	Phenanthrene	15	48	13,000 ES
120-12-7	Anthracene	17	48	5,900 E
206-44-0	Fluoranthene	18	48	8,200 E
129-00-0	Pyrene	22	48	7,000 E
56-55-3	Benzo (a) anthracene	21	48	2,100
218-01-9	Chrysene	18	48	2,200
205-99-2	Benzo (b) fluoranthene	20	48	840
207-08-9	Benzo (k) fluoranthene	22	48	530
50-32-8	Benzo (a) pyrene	23	48	850
193-39-5	Indeno (1,2,3-cd) pyrene	29	48	360
53-70-3	Dibenz (a,h) anthracene	25	48	110
191-24-2	Benzo (g,h,i) perylene	27	48	430
TOTBFA	Total Benzofluoranthenes	22	48	1,900


Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene 40.0%  
d10-2-Methylnaphthalene 56.7%  
d14-Dibenzo (a,h) anthracen D

**ORGANICS ANALYSIS DATA SHEET**  
**PNAs by SIM SW8270D-SIM GC/MS**  
**Extraction Method: SW3546**  
 Page 1 of 1

**Sample ID: PAI-2-18.0-18.5**  
**DILUTION**

Lab Sample ID: ZP35B  
 LIMS ID: 14-27576  
 Matrix: Soil  
 Data Release Authorized:   
 Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570  
 Date Sampled: 12/09/14  
 Date Received: 12/12/14

Date Extracted: 12/23/14  
 Date Analyzed: 01/20/15 11:35  
 Instrument/Analyst: NT8/JZ  
 GPC Cleanup: No

Sample Amount: 10.4 g-dry-wt  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1000  
 Percent Moisture: 30.8 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	2200	4,800	620,000 E
208-96-8	Acenaphthylene	1500	4,800	4,200 J
83-32-9	Acenaphthene	1400	4,800	14,000
86-73-7	Fluorene	1400	4,800	11,000
85-01-8	Phenanthrene	1500	4,800	28,000
120-12-7	Anthracene	1700	4,800	7,500
206-44-0	Fluoranthene	1800	4,800	12,000
129-00-0	Pyrene	2200	4,800	9,600
56-55-3	Benzo (a) anthracene	2100	4,800	2,400 J
218-01-9	Chrysene	1800	4,800	< 4,800 U
205-99-2	Benzo (b) fluoranthene	2000	4,800	< 4,800 U
207-08-9	Benzo (k) fluoranthene	2200	4,800	< 4,800 U
50-32-8	Benzo (a) pyrene	2300	4,800	< 4,800 U
193-39-5	Indeno (1,2,3-cd) pyrene	2900	4,800	< 4,800 U
53-70-3	Dibenz (a,h) anthracene	2500	4,800	< 4,800 U
191-24-2	Benzo (g,h,i) perylene	2700	4,800	< 4,800 U
TOTBFA	Total Benzofluoranthenes	2200	4,800	< 4,800 U

Reported in µg/kg (ppb)


**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene D  
 d10-2-Methylnaphthalene D  
 d14-Dibenzo (a,h) anthracen D



**ORGANICS ANALYSIS DATA SHEET**  
**PNAs by SIM SW8270D-SIM GC/MS**  
**Extraction Method: SW3546**  
 Page 1 of 1

**Sample ID: PAI-2-18.0-18.5**  
**DILUTION2**

Lab Sample ID: ZP35B  
 LIMS ID: 14-27576  
 Matrix: Soil  
 Data Release Authorized:   
 Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570  
 Date Sampled: 12/09/14  
 Date Received: 12/12/14

Date Extracted: 12/23/14  
 Date Analyzed: 01/20/15 16:54  
 Instrument/Analyst: NT8/JZ  
 GPC Cleanup: No

Sample Amount: 10.4 g-dry-wt  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 2000  
 Percent Moisture: 30.8 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	4400	9,600	640,000
208-96-8	Acenaphthylene	3100	9,600	< 9,600 U
83-32-9	Acenaphthene	2900	9,600	14,000
86-73-7	Fluorene	2800	9,600	11,000
85-01-8	Phenanthrene	3000	9,600	29,000
120-12-7	Anthracene	3400	9,600	6,900 J
206-44-0	Fluoranthene	3600	9,600	12,000
129-00-0	Pyrene	4400	9,600	8,800 J
56-55-3	Benzo(a)anthracene	4300	9,600	< 9,600 U
218-01-9	Chrysene	3700	9,600	< 9,600 U
205-99-2	Benzo(b)fluoranthene	4100	9,600	< 9,600 U
207-08-9	Benzo(k)fluoranthene	4400	9,600	< 9,600 U
50-32-8	Benzo(a)pyrene	4600	9,600	< 9,600 U
193-39-5	Indeno(1,2,3-cd)pyrene	5800	9,600	< 9,600 U
53-70-3	Dibenz(a,h)anthracene	4900	9,600	< 9,600 U
191-24-2	Benzo(g,h,i)perylene	5400	9,600	< 9,600 U
TOTBFA	Total Benzofluoranthenes	4400	9,600	< 9,600 U


Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene D  
 d10-2-Methylnaphthalene D  
 d14-Dibenzo(a,h)anthracen D

ORGANICS ANALYSIS DATA SHEET  
 PNAs by SIM SW8270D-SIM GC/MS  
 Extraction Method: SW3546  
 Page 1 of 1

Sample ID: PAI-10-29.5-30.0  
 SAMPLE

Lab Sample ID: ZP35D  
 LIMS ID: 14-27578  
 Matrix: Soil  
 Data Release Authorized:   
 Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570  
 Date Sampled: 12/11/14  
 Date Received: 12/12/14

Date Extracted: 12/23/14  
 Date Analyzed: 01/19/15 23:08  
 Instrument/Analyst: NT8/JZ  
 GPC Cleanup: No

Sample Amount: 10.1 g-dry-wt  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00  
 Percent Moisture: 7.8 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	2.2	4.9	3,300 ES
208-96-8	Acenaphthylene	1.6	4.9	67
83-32-9	Acenaphthene	1.5	4.9	91
86-73-7	Fluorene	1.4	4.9	89
85-01-8	Phenanthrene	1.6	4.9	330
120-12-7	Anthracene	1.8	4.9	84
206-44-0	Fluoranthene	1.8	4.9	310
129-00-0	Pyrene	2.2	4.9	300
56-55-3	Benzo (a) anthracene	2.2	4.9	120
218-01-9	Chrysene	1.9	4.9	140
205-99-2	Benzo (b) fluoranthene	2.1	4.9	110
207-08-9	Benzo (k) fluoranthene	2.2	4.9	59
50-32-8	Benzo (a) pyrene	2.3	4.9	130
193-39-5	Indeno (1,2,3-cd) pyrene	3.0	4.9	110
53-70-3	Dibenz (a,h) anthracene	2.5	4.9	22
191-24-2	Benzo (g,h,i) perylene	2.8	4.9	150
TOTBFA	Total Benzofluoranthenes	2.2	4.9	220


Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene	62.0%
d10-2-Methylnaphthalene	42.3%
d14-Dibenzo (a,h) anthracen	66.0%

**ORGANICS ANALYSIS DATA SHEET**  
**PNAs by SIM SW8270D-SIM GC/MS**  
**Extraction Method: SW3546**  
 Page 1 of 1

**Sample ID: PAI-10-29.5-30.0**  
**DILUTION**

Lab Sample ID: ZP35D  
 LIMS ID: 14-27578  
 Matrix: Soil  
 Data Release Authorized:   
 Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570  
 Date Sampled: 12/11/14  
 Date Received: 12/12/14

Date Extracted: 12/23/14  
 Date Analyzed: 01/20/15 12:00  
 Instrument/Analyst: NT8/JZ  
 GPC Cleanup: No

Sample Amount: 10.1 g-dry-wt  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 30.0  
 Percent Moisture: 7.8 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	67	150	7,600
208-96-8	Acenaphthylene	48	150	< 150 U
83-32-9	Acenaphthene	44	150	87 J
86-73-7	Fluorene	43	150	< 150 U
85-01-8	Phenanthrene	47	150	310
120-12-7	Anthracene	53	150	83 J
206-44-0	Fluoranthene	55	150	310
129-00-0	Pyrene	67	150	270
56-55-3	Benzo (a) anthracene	66	150	110 J
218-01-9	Chrysene	57	150	120 J
205-99-2	Benzo (b) fluoranthene	62	150	100 J
207-08-9	Benzo (k) fluoranthene	67	150	< 150 U
50-32-8	Benzo (a) pyrene	70	150	120 J
193-39-5	Indeno (1,2,3-cd) pyrene	89	150	98 J
53-70-3	Dibenz (a,h) anthracene	76	150	< 150 U
191-24-2	Benzo (g,h,i) perylene	83	150	150 J
TOTBFA	Total Benzofluoranthenes	67	150	220

Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene D  
 d10-2-Methylnaphthalene D  
 d14-Dibenzo (a,h) anthracen D



ORGANICS ANALYSIS DATA SHEET  
PNAs by SIM SW8270D-SIM GC/MS  
Extraction Method: SW3546  
Page 1 of 1

Sample ID: PAI-12-19.5-20.0  
SAMPLE

Lab Sample ID: ZP35E  
LIMS ID: 14-27579  
Matrix: Soil  
Data Release Authorized: *AB*  
Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: 12/12/14  
Date Received: 12/12/14

Date Extracted: 12/23/14  
Date Analyzed: 01/20/15 00:23  
Instrument/Analyst: NT8/JZ  
GPC Cleanup: No

Sample Amount: 0.74 g-dry-wt  
Final Extract Volume: 0.50 mL  
Dilution Factor: 10.0  
Percent Moisture: 28.4 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	310	680	340,000 ES
208-96-8	Acenaphthylene	220	680	13,000
83-32-9	Acenaphthene	200	680	6,900
86-73-7	Fluorene	200	680	31,000
85-01-8	Phenanthrene	210	680	90,000 E
120-12-7	Anthracene	240	680	27,000
206-44-0	Fluoranthene	250	680	48,000
129-00-0	Pyrene	310	680	52,000
56-55-3	Benzo (a) anthracene	300	680	17,000
218-01-9	Chrysene	260	680	20,000
205-99-2	Benzo (b) fluoranthene	290	680	9,400
207-08-9	Benzo (k) fluoranthene	310	680	5,400
50-32-8	Benzo (a) pyrene	320	680	16,000
193-39-5	Indeno (1,2,3-cd) pyrene	410	680	7,400
53-70-3	Dibenz (a,h) anthracene	350	680	2,000
191-24-2	Benzo (g,h,i) perylene	380	680	10,000
TOTBFA	Total Benzofluoranthenes	310	680	21,000


Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene	63.3%
d10-2-Methylnaphthalene	63.3%
d14-Dibenzo (a,h) anthracen	63.3%

ORGANICS ANALYSIS DATA SHEET  
PNAs by SIM SW8270D-SIM GC/MS  
Extraction Method: SW3546  
Page 1 of 1

Sample ID: PAI-12-19.5-20.0  
DILUTION

Lab Sample ID: ZP35E  
LIMS ID: 14-27579  
Matrix: Soil  
Data Release Authorized:   
Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: 12/12/14  
Date Received: 12/12/14

Date Extracted: 12/23/14  
Date Analyzed: 01/20/15 12:26  
Instrument/Analyst: NT8/JZ  
GPC Cleanup: No

Sample Amount: 0.74 g-dry-wt  
Final Extract Volume: 0.50 mL  
Dilution Factor: 300  
Percent Moisture: 28.4 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	9200	20,000	700,000
208-96-8	Acenaphthylene	6500	20,000	14,000 J
83-32-9	Acenaphthene	6000	20,000	< 20,000 U
86-73-7	Fluorene	6000	20,000	32,000
85-01-8	Phenanthrene	6400	20,000	110,000
120-12-7	Anthracene	7200	20,000	29,000
206-44-0	Fluoranthene	7600	20,000	52,000
129-00-0	Pyrene	9200	20,000	57,000
56-55-3	Benzo (a) anthracene	9000	20,000	16,000 J
218-01-9	Chrysene	7800	20,000	19,000 J
205-99-2	Benzo (b) fluoranthene	8600	20,000	9,700 J
207-08-9	Benzo (k) fluoranthene	9200	20,000	< 20,000 U
50-32-8	Benzo (a) pyrene	9600	20,000	16,000 J
193-39-5	Indeno (1,2,3-cd) pyrene	12000	20,000	< 20,000 U
53-70-3	Dibenz (a,h) anthracene	10000	20,000	< 20,000 U
191-24-2	Benzo (g,h,i) perylene	11000	20,000	11,000 J
TOTBFA	Total Benzofluoranthenes	9200	20,000	22,000

Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene D  
d10-2-Methylnaphthalene D  
d14-Dibenzo (a,h) anthracen D

**SIM SW8270 SURROGATE RECOVERY SUMMARY**

Matrix: Soil

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570

<u>Client ID</u>	<u>FLN</u>	<u>MNP</u>	<u>DBA</u>	<u>TOT OUT</u>
PAI-2-18.0-18.5	40.0%	56.7%	D D	0
PAI-2-18.0-18.5 DL	D D	D D	D D	0
PAI-2-18.0-18.5 DL2	D D	D D	D D	0
MB-122314	73.3%	55.3%	82.7%	0
LCS-122314	71.3%	52.0%	80.3%	0
PAI-10-29.5-30.0	62.0%	42.3%	66.0%	0
PAI-10-29.5-30.0 DL	D D	D D	D D	0
PAI-10-29.5-30.0 MS	77.3%	51.7%	78.7%	0
PAI-10-29.5-30.0 MSD	76.3%	50.7%	78.7%	0
PAI-12-19.5-20.0	63.3%	63.3%	63.3%	0
PAI-12-19.5-20.0 DL	D D	D D	D D	0

	<b>LCS/MB LIMITS</b>	<b>QC LIMITS</b>
(FLN) = d10-Fluoranthene	(36-134)	(36-134)
(MNP) = d10-2-Methylnaphthalene	(32-120)	(32-120)
(DBA) = d14-Dibenzo(a,h)anthracene	(21-133)	(21-133)

Prep Method: SW3546  
Log Number Range: 14-27576 to 14-27579



ORGANICS ANALYSIS DATA SHEET  
PNAs by SW8270D-SIM GC/MS  
Page 1 of 1

Sample ID: PAI-10-29.5-30.0  
MATRIX SPIKE

Lab Sample ID: ZP35D  
LIMS ID: 14-27578  
Matrix: Soil  
Data Release Authorized: *AS*  
Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
Event: 0186-846-01 Task 1570  
Date Sampled: 12/11/14  
Date Received: 12/12/14

Date Extracted MS/MSD: 12/23/14  
Date Analyzed MS: 01/19/15 23:33  
MSD: 01/19/15 23:58  
Instrument/Analyst MS: NT8/JZ  
MSD: NT8/JZ

Sample Amount MS: 10.19 g-dry-wt  
MSD: 10.19 g-dry-wt  
Final Extract Volume MS: 0.50 mL  
MSD: 0.50 mL  
Dilution Factor MS: 1.00  
MSD: 1.00

Analyte	Sample	MS	Spike Added-MS	MS Recovery	MSD	Spike Added-MSD	MSD Recovery	RPD
Naphthalene	3300 ES	4940 ES	147	NA	4770 ES	147	NA	3.5%
Acenaphthylene	67	170	147	70.1%	201	147	91.2%	16.7%
Acenaphthene	91	190	147	67.3%	250	147	108%	27.3%
Fluorene	89	198	147	74.1%	330	147	164%	50.0%
Phenanthrene	330	423	147	63.3%	678 E	147	237%	46.3%
Anthracene	84	188	147	70.7%	246	147	110%	26.7%
Fluoranthene	310	433	147	83.7%	474	147	112%	9.0%
Pyrene	300	417	147	79.6%	444	147	98.0%	6.3%
Benzo(a)anthracene	120	239	147	81.0%	242	147	83.0%	1.2%
Chrysene	140	255	147	78.2%	260	147	81.6%	1.9%
Benzo(b)fluoranthene	110	230	147	81.6%	233	147	83.7%	1.3%
Benzo(k)fluoranthene	59	174	147	78.2%	174	147	78.2%	0.0%
Benzo(a)pyrene	130	254	147	84.4%	255	147	85.0%	0.4%
Indeno(1,2,3-cd)pyrene	110	234	147	84.4%	233	147	83.7%	0.4%
Dibenz(a,h)anthracene	22	137	147	78.2%	136	147	77.6%	0.7%
Benzo(g,h,i)perylene	150	283	147	90.5%	279	147	87.8%	1.4%
Total Benzofluoranthenes	220	557	442	76.2%	559	442	76.7%	0.4%

Reported in µg/kg (ppb)

NA-No recovery due to high concentration (> 4X) of analyte in original sample, calculated negative recovery, or undetected spike.  
RPD calculated using sample concentrations per SW846.

**ORGANICS ANALYSIS DATA SHEET**  
**PNAs by SIM SW8270D-SIM GC/MS**  
**Extraction Method: SW3546**  
 Page 1 of 1

**Sample ID: PAI-10-29.5-30.0**  
**MATRIX SPIKE**

Lab Sample ID: ZP35D  
 LIMS ID: 14-27578  
 Matrix: Soil  
 Data Release Authorized: *AS*  
 Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570  
 Date Sampled: 12/11/14  
 Date Received: 12/12/14

Date Extracted: 12/23/14  
 Date Analyzed: 01/19/15 23:33  
 Instrument/Analyst: NT8/JZ  
 GPC Cleanup: No

Sample Amount: 10.2 g-dry-wt  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00  
 Percent Moisture: 7.8 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	2.2	4.9	---
208-96-8	Acenaphthylene	1.6	4.9	---
83-32-9	Acenaphthene	1.5	4.9	---
86-73-7	Fluorene	1.4	4.9	---
85-01-8	Phenanthrene	1.6	4.9	---
120-12-7	Anthracene	1.7	4.9	---
206-44-0	Fluoranthene	1.8	4.9	---
129-00-0	Pyrene	2.2	4.9	---
56-55-3	Benzo(a)anthracene	2.2	4.9	---
218-01-9	Chrysene	1.9	4.9	---
205-99-2	Benzo(b)fluoranthene	2.1	4.9	---
207-08-9	Benzo(k)fluoranthene	2.2	4.9	---
50-32-8	Benzo(a)pyrene	2.3	4.9	---
193-39-5	Indeno(1,2,3-cd)pyrene	3.0	4.9	---
53-70-3	Dibenz(a,h)anthracene	2.5	4.9	---
191-24-2	Benzo(g,h,i)perylene	2.7	4.9	---
TOTBFA	Total Benzofluoranthenes	2.2	4.9	---

Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene	77.3%
d10-2-Methylnaphthalene	51.7%
d14-Dibenzo(a,h)anthracen	78.7%

**ORGANICS ANALYSIS DATA SHEET**  
**PNA's by SIM SW8270D-SIM GC/MS**  
**Extraction Method: SW3546**  
 Page 1 of 1

**Sample ID: PAI-10-29.5-30.0**  
**MATRIX SPIKE DUP**

Lab Sample ID: ZP35D  
 LIMS ID: 14-27578  
 Matrix: Soil  
 Data Release Authorized:  
 Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570  
 Date Sampled: 12/11/14  
 Date Received: 12/12/14

Date Extracted: 12/23/14  
 Date Analyzed: 01/19/15 23:58  
 Instrument/Analyst: NT8/JZ  
 GPC Cleanup: No

Sample Amount: 10.2 g-dry-wt  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00  
 Percent Moisture: 7.8 %

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	2.2	4.9	---
208-96-8	Acenaphthylene	1.6	4.9	---
83-32-9	Acenaphthene	1.5	4.9	---
86-73-7	Fluorene	1.4	4.9	---
85-01-8	Phenanthrene	1.6	4.9	---
120-12-7	Anthracene	1.7	4.9	---
206-44-0	Fluoranthene	1.8	4.9	---
129-00-0	Pyrene	2.2	4.9	---
56-55-3	Benzo(a)anthracene	2.2	4.9	---
218-01-9	Chrysene	1.9	4.9	---
205-99-2	Benzo(b)fluoranthene	2.1	4.9	---
207-08-9	Benzo(k)fluoranthene	2.2	4.9	---
50-32-8	Benzo(a)pyrene	2.3	4.9	---
193-39-5	Indeno(1,2,3-cd)pyrene	3.0	4.9	---
53-70-3	Dibenz(a,h)anthracene	2.5	4.9	---
191-24-2	Benzo(g,h,i)perylene	2.7	4.9	---
TOTBFA	Total Benzofluoranthenes	2.2	4.9	---

Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene	76.3%
d10-2-Methylnaphthalene	50.7%
d14-Dibenzo(a,h)anthracen	78.7%



**ORGANICS ANALYSIS DATA SHEET**

**PNAs by SW8270D-SIM GC/MS**

Page 1 of 1


**Sample ID: LCS-122314**

**LAB CONTROL SAMPLE**

Lab Sample ID: LCS-122314

LIMS ID: 14-27578

Matrix: Soil

Data Release Authorized: 

Reported: 01/21/15

QC Report No: ZP35-Geoengineers

Project: Gas Works Park-Play Area Investigat

Event: 0186-846-01 Task 1570

Date Sampled: NA

Date Received: NA

Date Extracted: 12/23/14

Date Analyzed LCS: 01/19/15 22:18

Instrument/Analyst LCS: NT8/JZ

Sample Amount LCS: 10.00 g-dry-wt

Final Extract Volume LCS: 0.50 mL

Dilution Factor LCS: 1.00

Analyte	LCS	Spike Added	Recovery
Naphthalene	80.4	150	53.6%
Acenaphthylene	84.8	150	56.5%
Acenaphthene	79.2	150	52.8%
Fluorene	89.6	150	59.7%
Phenanthrene	96.2	150	64.1%
Anthracene	93.0	150	62.0%
Fluoranthene	105	150	70.0%
Pyrene	97.5	150	65.0%
Benzo(a)anthracene	104	150	69.3%
Chrysene	101	150	67.3%
Benzo(b)fluoranthene	108	150	72.0%
Benzo(k)fluoranthene	102	150	68.0%
Benzo(a)pyrene	96.9	150	64.6%
Indeno(1,2,3-cd)pyrene	107	150	71.3%
Dibenz(a,h)anthracene	109	150	72.7%
Benzo(g,h,i)perylene	109	150	72.7%
Total Benzofluoranthenes	289	450	64.2%

Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene	71.3%
d10-2-Methylnaphthalene	52.0%
d14-Dibenzo(a,h)anthracene	80.3%

4B  
SEMIVOLATILE METHOD BLANK SUMMARY

BLANK NO.

ZP35MBS1

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY

Lab File ID: 01191530

Date Extracted: 12/23/14

Instrument ID: NT8

Date Analyzed: 01/19/15

Matrix: SOLID

Time Analyzed: 2153

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
01	ZP35LCSS1	ZP35LCSS1	01191531	01/19/15
02	PAI-2-18.0-18.5	ZP35B	01191532	01/19/15
03	PAI-10-29.5-30.0	ZP35D	01191533	01/19/15
04	PAI-10-29.5-30.	ZP35DMS	01191534	01/19/15
05	PAI-10-29.5-30.	ZP35DMSD	01191535	01/19/15
06	PAI-12-19.5-20.0	ZP35E	01191536	01/20/15
07	PAI-2-18.0-18.5	ZP35B	01201506	01/20/15
08	PAI-10-29.5-30.0	ZP35D	01201507	01/20/15
09	PAI-12-19.5-20.0	ZP35E	01201508	01/20/15
10	PAI-2-18.0-18.5	ZP35B	01201514	01/20/15
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**ORGANICS ANALYSIS DATA SHEET**  
**PNAs by SIM SW8270D-SIM GC/MS**  
**Extraction Method: SW3546**  
 Page 1 of 1

**Sample ID: MB-122314**  
**METHOD BLANK**

Lab Sample ID: MB-122314  
 LIMS ID: 14-27578  
 Matrix: Soil  
 Data Release Authorized:  
 Reported: 01/21/15

QC Report No: ZP35-Geoengineers  
 Project: Gas Works Park-Play Area Investigat  
 0186-846-01 Task 1570  
 Date Sampled: NA  
 Date Received: NA

Date Extracted: 12/23/14  
 Date Analyzed: 01/19/15 21:53  
 Instrument/Analyst: NT8/JZ  
 GPC Cleanup: No

Sample Amount: 10.0 g-dry-wt  
 Final Extract Volume: 0.50 mL  
 Dilution Factor: 1.00  
 Percent Moisture: NA

CAS Number	Analyte	DL	LOQ	Result
91-20-3	Naphthalene	2.3	5.0	< 5.0 U
208-96-8	Acenaphthylene	1.6	5.0	< 5.0 U
83-32-9	Acenaphthene	1.5	5.0	< 5.0 U
86-73-7	Fluorene	1.5	5.0	< 5.0 U
85-01-8	Phenanthrene	1.6	5.0	< 5.0 U
120-12-7	Anthracene	1.8	5.0	< 5.0 U
206-44-0	Fluoranthene	1.9	5.0	< 5.0 U
129-00-0	Pyrene	2.3	5.0	< 5.0 U
56-55-3	Benzo(a)anthracene	2.2	5.0	< 5.0 U
218-01-9	Chrysene	1.9	5.0	< 5.0 U
205-99-2	Benzo(b)fluoranthene	2.1	5.0	< 5.0 U
207-08-9	Benzo(k)fluoranthene	2.3	5.0	< 5.0 U
50-32-8	Benzo(a)pyrene	2.4	5.0	< 5.0 U
193-39-5	Indeno(1,2,3-cd)pyrene	3.0	5.0	< 5.0 U
53-70-3	Dibenz(a,h)anthracene	2.6	5.0	< 5.0 U
191-24-2	Benzo(g,h,i)perylene	2.8	5.0	< 5.0 U
TOTBFA	Total Benzofluoranthenes	2.3	5.0	< 5.0 U

Reported in µg/kg (ppb)

**SIM Semivolatile Surrogate Recovery**

d10-Fluoranthene	73.3%
d10-2-Methylnaphthalene	55.3%
d14-Dibenzo(a,h)anthracen	82.7%



5B  
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

Instrument ID: NT8

Project: GAS WORKS PARK-PLAY

DFTPP Injection Date: 01/05/15

DFTPP Injection Time: 1510

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	25.9
68	Less than 2.0% of mass 69	1.0 ( 1.9)1
69	Mass 69 relative abundance	50.7
70	Less than 2.0% of mass 69	0.3 ( 0.5)1
127	10.0 - 80.0% of mass 198	51.2
197	Less than 2.0% of mass 198	0.7
198	Base Peak, 100% relative abundance	100.0
199	5.0 to 9.0% of mass 198	7.7
275	10.0 - 60.0% of mass 198	31.2
365	Greater than 1.0% of mass 198	4.12
441	0.0 - 24.0% of mass 442	8.7 ( 14.6)2
442	50.0 - 200.0% of mass 198	59.9
443	15.0 - 24.0% of mass 442	12.1 ( 20.2)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	IC250105	SDA0002-CAL4	01051502	01/05/15	1523
02	IC010105	SDA0002-CAL1	01051503	01/05/15	1548
03	IC050105	SDA0002-CAL2	01051504	01/05/15	1614
04	IC10105	SDA0002-CAL3	01051505	01/05/15	1639
05	IC50105	SDA0002-CAL5	01051506	01/05/15	1704
06	IC100105	SDA0002-CAL6	01051507	01/05/15	1730
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5B  
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

Instrument ID: NT8

Project: GAS WORKS PARK-PLAY

DFTPP Injection Date: 01/19/15

DFTPP Injection Time: 1844

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	19.2
68	Less than 2.0% of mass 69	0.0 ( 0.0)1
69	Mass 69 relative abundance	52.8
70	Less than 2.0% of mass 69	0.3 ( 0.6)1
127	10.0 - 80.0% of mass 198	50.3
197	Less than 2.0% of mass 198	0.0
198	Base Peak, 100% relative abundance	100.0
199	5.0 to 9.0% of mass 198	7.5
275	10.0 - 60.0% of mass 198	31.8
365	Greater than 1.0% of mass 198	3.51
441	0.0 - 24.0% of mass 442	10.1 ( 16.6)2
442	50.0 - 200.0% of mass 198	61.0
443	15.0 - 24.0% of mass 442	13.1 ( 21.5)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	ICV0119A	ICV0119A	01191523	01/19/15	1857
02	ZP35MBS1	ZP35MBS1	01191530	01/19/15	2153
03	ZP35LCSS1	ZP35LCSS1	01191531	01/19/15	2218
04	PAI-2-18.0-18.5	ZP35B	01191532	01/19/15	2243
05	PAI-10-29.5-30.0	ZP35D	01191533	01/19/15	2308
06	PAI-10-29.5-30.	ZP35DMS	01191534	01/19/15	2333
07	PAI-10-29.5-30.	ZP35DMSD	01191535	01/19/15	2358
08	PAI-12-19.5-20.0	ZP35E	01191536	01/20/15	0023
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5B  
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

Instrument ID: NT8

Project: GAS WORKS PARK-PLAY

DFTPP Injection Date: 01/20/15

DFTPP Injection Time: 0936

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
51	10.0 - 80.0% of mass 198	18.9
68	Less than 2.0% of mass 69	0.0 ( 0.0)1
69	Mass 69 relative abundance	53.6
70	Less than 2.0% of mass 69	0.3 ( 0.7)1
127	10.0 - 80.0% of mass 198	49.4
197	Less than 2.0% of mass 198	0.0
198	Base Peak, 100% relative abundance	100.0
199	5.0 to 9.0% of mass 198	7.5
275	10.0 - 60.0% of mass 198	32.4
365	Greater than 1.0% of mass 198	3.75
441	0.0 - 24.0% of mass 442	9.7 ( 15.6)2
442	50.0 - 200.0% of mass 198	62.3
443	15.0 - 24.0% of mass 442	12.5 ( 20.1)2

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	ICV0120	ICV0120	01201502	01/20/15	0950
02	PAI-2-18.0-18.5	ZP35B	01201506	01/20/15	1135
03	PAI-10-29.5-30.0	ZP35D	01201507	01/20/15	1200
04	PAI-12-19.5-20.0	ZP35E	01201508	01/20/15	1226
05	PAI-2-18.0-18.5	ZP35B	01201514	01/20/15	1654
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## SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY

Instrument ID: NT8

Cont. Calib. Date: 01/19/15

Init. Calib. Date: 01/05/15

Cont. Calib. Time: 1857

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
=====	=====	=====	=====	=====	=====
Naphthalene	1.150	1.050	0.700	AVRG	-8.7
2-Methylnaphthalene	0.673	0.705	0.400	AVRG	4.8
Acenaphthylene	1.724	1.827	0.900	AVRG	6.0
Acenaphthene	1.171	1.171	0.900	AVRG	0.0
Dibenzofuran	1.645	1.631	0.800	AVRG	-0.8
Fluorene	1.333	1.358	0.900	AVRG	1.9
Phenanthrene	1.088	1.081	0.700	AVRG	-0.6
Anthracene	1.078	1.098	0.700	AVRG	1.8
Fluoranthene	1.316	1.334	0.600	AVRG	1.4
Pyrene	1.231	1.133	0.600	AVRG	-8.0
Benzo(a)anthracene	1.218	1.164	0.800	AVRG	-4.4
Chrysene	1.182	1.116	0.700	AVRG	-5.6
Benzo(b)fluoranthene	1.133	1.151	0.700	AVRG	1.6
Benzo(k)fluoranthene	1.178	1.165	0.700	AVRG	-1.1
Benzo(j)fluoranthene	1.160	1.111	0.010	AVRG	-4.2
Benzo(a)pyrene	1.116	1.134	0.700	AVRG	1.6
Indeno(1,2,3-cd)pyrene	1.241	1.322	0.500	AVRG	6.5
Dibenzo(a,h)anthracene	1.024	1.106	0.400	AVRG	8.0
Benzo(g,h,i)perylene	1.073	1.133	0.500	AVRG	5.6
1-methylnaphthalene	0.622	0.653	0.010	AVRG	5.0
Perylene	1.122	1.113	0.010	AVRG	-0.8
=====	=====	=====	=====	=====	=====
2-Methylnaphthalene-d10	0.707	0.719	0.010	AVRG	1.7
Dibenzo(a,h)anthracene-d14	0.820	0.936	0.010	AVRG	14.1
Fluoranthene-d10	1.185	1.209	0.010	AVRG	2.0

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF



## SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY

Instrument ID: NT8

Cont. Calib. Date: 01/20/15

Init. Calib. Date: 01/05/15

Cont. Calib. Time: 0950

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
=====	=====	=====	=====	=====	=====
Naphthalene	1.150	1.065	0.700	AVRG	-7.4
2-Methylnaphthalene	0.673	0.696	0.400	AVRG	3.4
Acenaphthylene	1.724	1.821	0.900	AVRG	5.6
Acenaphthene	1.171	1.194	0.900	AVRG	2.0
Dibenzofuran	1.645	1.625	0.800	AVRG	-1.2
Fluorene	1.333	1.392	0.900	AVRG	4.4
Phenanthrene	1.088	1.086	0.700	AVRG	-0.2
Anthracene	1.078	1.108	0.700	AVRG	2.8
Fluoranthene	1.316	1.336	0.600	AVRG	1.5
Pyrene	1.231	1.137	0.600	AVRG	-7.6
Benzo(a)anthracene	1.218	1.199	0.800	AVRG	-1.6
Chrysene	1.182	1.132	0.700	AVRG	-4.2
Benzo(b)fluoranthene	1.133	1.129	0.700	AVRG	-0.4
Benzo(k)fluoranthene	1.178	1.167	0.700	AVRG	-0.9
Benzo(j)fluoranthene	1.160	1.179	0.010	AVRG	1.6
Benzo(a)pyrene	1.116	1.120	0.700	AVRG	0.4
Indeno(1,2,3-cd)pyrene	1.241	1.216	0.500	AVRG	-2.0
Dibenzo(a,h)anthracene	1.024	1.059	0.400	AVRG	3.4
Benzo(g,h,i)perylene	1.073	1.081	0.500	AVRG	0.7
1-methylnaphthalene	0.622	0.656	0.010	AVRG	5.5
Perylene	1.122	1.124	0.010	AVRG	0.2
=====	=====	=====	=====	=====	=====
2-Methylnaphthalene-d10	0.707	0.697	0.010	AVRG	-1.4
Dibenzo(a,h)anthracene-d14	0.820	0.865	0.010	AVRG	5.5
Fluoranthene-d10	1.185	1.211	0.010	AVRG	2.2

&lt;- Exceeds QC limit of 20% D

\* RF less than minimum RF



8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/19/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
CCAL	462048	4.55	296335	6.81	537194	8.82
UPPER LIMIT		5.05		7.31		9.32
LOWER LIMIT		4.05		6.31		8.32
01 ZP35MBS1	410219	4.55	273837	6.81	488450	8.82
02 ZP35LCSS1	436987	4.55	276604	6.81	495657	8.82
03 PAI-2-18.0-1	554754	4.58	336136	6.82	552321	8.83
04 PAI-10-29.5-	401613	4.56	261542	6.81	467211	8.82
05 PAI-10-29.5-	459388	4.56	269142	6.81	486475	8.82
06 PAI-10-29.5-	474262	4.57	279375	6.81	493770	8.82
07 PAI-12-19.5-	431703	4.56	284357	6.81	503149	8.82
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IS1 = Naphthalene-d8  
IS2 = Acenaphthene-d10  
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/19/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
=====	=====	=====	=====	=====	=====	=====
CCAL	652608	13.46	652482	17.20		
UPPER LIMIT		13.96		17.70		
LOWER LIMIT		12.96		16.70		
01 ZP35MBS1	602025	13.46	628969	17.21		
02 ZP35LCSS1	594239	13.45	600570	17.21		
03 PAI-2-18.0-1	658596	13.46	677235	17.21		
04 PAI-10-29.5-	554891	13.45	567695	17.21		
05 PAI-10-29.5-	576771	13.45	580697	17.21		
06 PAI-10-29.5-	590772	13.45	585604	17.21		
07 PAI-12-19.5-	606747	13.45	610228	17.21		
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IS4 = Chrysene-d12

IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.



8B  
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/20/15

	IS1 (NPT) AREA #	RT #	IS2 (ANT) AREA #	RT #	IS3 (PHN) AREA #	RT #
ICAL MIDPT	387555	4.71	258447	6.97	456948	8.98
UPPER LIMIT	775110		516894		913896	
LOWER LIMIT	193778		129224		228474	
CCAL	389201	4.55	259938	6.81	477075	8.82
UPPER LIMIT		5.05		7.31		9.32
LOWER LIMIT		4.05		6.31		8.32
01 PAI-2-18.0-1	367872	4.55	239912	6.81	416451	8.82
02 PAI-10-29.5-	386963	4.55	252604	6.81	444303	8.82
03 PAI-12-19.5-	402737	4.56	258461	6.81	455206	8.82
04 PAI-2-18.0-1	334827	4.55	218873	6.81	390968	8.82
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IS1 = Naphthalene-d8  
IS2 = Acenaphthene-d10  
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.



SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: GEOENGINEERS

ARI Job No: ZP35

Project: GAS WORKS PARK-PLAY

Ical Midpoint ID: 01051502

Ical Date: 01/05/15

Instrument ID: NT8

Cont. Cal Date: 01/20/15

	IS4 (CRY) AREA #	RT #	IS5 (PRY) AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	492500	13.72	485582	17.52		
UPPER LIMIT	985000		971164			
LOWER LIMIT	246250		242791			
=====	=====	=====	=====	=====	=====	=====
CCAL	577051	13.45	560055	17.20		
UPPER LIMIT		13.95		17.70		
LOWER LIMIT		12.95		16.70		
01 PAI-2-18.0-1	531889	13.45	558571	17.21		
02 PAI-10-29.5-	551919	13.45	567725	17.20		
03 PAI-12-19.5-	550147	13.45	552530	17.20		
04 PAI-2-18.0-1	491438	13.46	466538	17.21		
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IS4 = Chrysene-d12  
 IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint  
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint  
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal  
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

\* Values outside of QC limits.

Metals Analysis  
Report and Summary QC Forms

ARI Job ID: ZP35

# Cover Page

## INORGANIC ANALYSIS DATA PACKAGE



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

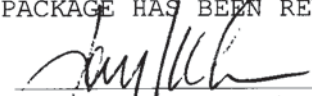
SDG: ZP35

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
PAI-2-12.5-13.0	ZP35A	14-27575	
PAI-2-12.5-13.0D	ZP35ADUP	14-27575	
PAI-2-12.5-13.0S	ZP35ASPK	14-27575	
PAI-9-9.5-10.0	ZP35C	14-27577	
PBS	ZP35MB1	14-27577	
LCSS	ZP35MB1SPK	14-27577	

Were ICP interelement corrections applied ?                      Yes/No    YES  
Were ICP background corrections applied ?                      Yes/No    YES  
If yes - were raw data generated before  
application of background corrections ?                      Yes/No    NO

Comments: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature:                       Name: Jay Kuhn  
Date: 12/30/14                      Title: Inorganics Director



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: PAI-2-12.5-13.0

SAMPLE

Lab Sample ID: ZP35A

LIMS ID: 14-27575

Matrix: Soil

Data Release Authorized:

Reported: 12/30/14

QC Report No: ZP35-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01 Task 1570

Date Sampled: 12/09/14

Date Received: 12/12/14

Percent Total Solids: 56.9%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/29/14	7440-38-2	Arsenic	2.0	20	80

U-Analyte undetected at given DL

J-Analyte detected between DL and LOQ

DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: PAI-9-9.5-10.0  
SAMPLE

Lab Sample ID: ZP35C  
LIMS ID: 14-27577  
Matrix: Soil  
Data Release Authorized:  
Reported: 12/30/14

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: 12/11/14  
Date Received: 12/12/14

Percent Total Solids: 89.1%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/29/14	7440-38-2	Arsenic	0.48	5	160

U-Analyte undetected at given DL  
J-Analyte detected between DL and LOQ  
DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: PAI-2-12.5-13.0

MATRIX SPIKE

Lab Sample ID: ZP35A  
LIMS ID: 14-27575  
Matrix: Soil  
Data Release Authorized:  
Reported: 12/30/14

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: 12/09/14  
Date Received: 12/12/14

**MATRIX SPIKE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Arsenic	6010C	80	430	343	102%	

Reported in mg/kg-dry

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: PAI-2-12.5-13.0  
DUPLICATE

Lab Sample ID: ZP35A  
LIMS ID: 14-27575  
Matrix: Soil  
Data Release Authorized:  
Reported: 12/30/14

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: 12/09/14  
Date Received: 12/12/14

**MATRIX DUPLICATE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010C	80	80	0.0%	+/- 20	L

Reported in mg/kg-dry

\*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: ZP35LCS  
LIMS ID: 14-27577  
Matrix: Soil  
Data Release Authorized:  
Reported: 12/30/14

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: NA  
Date Received: NA

**BLANK SPIKE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Arsenic	6010C	220	200	110%	

Reported in mg/kg-dry

N-Control limit not met  
NA-Not Applicable, Analyte Not Spiked  
Control Limits: 80-120%

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

**Sample ID: METHOD BLANK**

Page 1 of 1

Lab Sample ID: ZP35MB  
LIMS ID: 14-27577  
Matrix: Soil  
Data Release Authorized:  
Reported: 12/30/14

QC Report No: ZP35-Geoengineers  
Project: Gas Works Park-Play Area Investigat  
0186-846-01 Task 1570  
Date Sampled: NA  
Date Received: NA

Percent Total Solids: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	DL	LOQ	mg/kg Q
3050B	12/23/14	6010C	12/29/14	7440-38-2	Arsenic	0.46	5	5 U

U-Analyte undetected at given DL  
J-Analyte detected between DL and LOQ  
DL-Method Detection Limit

Results reported below the LOQ are for statistical purposes only and have not been evaluated by either an analyst or data reviewer.



# Calibration Verification



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

UNITS: ug/L

SDG: ZP35

ANALYTE	AS	ICP	EL	M	RUN	ICV	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic		IP122971				2000.0	2046.13	102.3		2000.0	2031.70	101.6	2021.54	101.1	2098.61	104.9	2122.52	106.1	2127.04	106.4

Control Limits: Mercury 80-120; Other Metals 90-110

# Calibration Verification



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP35

UNITS: ug/L

ANALYTE	EL	M	RUN	CCVTV	CCV6	%R	CCV7	%R	CCV8	%R	CCV9	%R	CCV10	%R	CCV11	%R
Arsenic	AS	ICP	IP122971	2000.0	2148.74	107.4										

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

ZP35 : 00094

# CRDL Standard



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP35

UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Arsenic	AS	ICP	IPI22971	50.0		50.05	100.1										

Control Limits: no control limits have been established by the EPA at this time.

FORM II (2)

ZP35 : 00005



# Calibration Blanks



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP35

UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Arsenic	AS	ICP	IP122971	10.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U

# Calibration Blanks

CLIENT: Geoenigneers

PROJECT: Gas Works Park-Play

SDG: ZP35



UNITS: ug/L

ANALYTE	AS	ICP	EL	METH	RUN	CRDL	IDL	CCB6	CCB7	CCB8	CCB9	CCB10	CCB11	C
Arsenic					IP122971	10.0	50.0	50.0						

# ICP Interference Check Sample



CLIENT: Geoenineers

ICS SOURCE: I.V.

PROJECT: Gas Works Park-Play

RUNID: IP122971

SDG: ZP35

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSAB1	%R	ICSA2	ICSAB2	%R	ICSA3	ICSAB3	%R
Aluminum	200000	200000	205572.8	211758.7	105.9						
Antimony	1000	1000	16.6	1069.0	106.9						
Arsenic	1000	1000	18.7	1067.1	106.7						
Barium	1000	1000	-3.2	1025.0	102.5						
Beryllium	1000	1000	0.1	1017.4	101.7						
Boron			-3.6	-4.4							
Cadmium	1000	1000	0.9	1021.4	102.1						
Calcium	100000	100000	103567.3	103318.8	103.3						
Chromium	1000	1000	-0.5	1033.1	103.3						
Cobalt	1000	1000	2.1	970.3	97.0						
Copper	1000	1000	0.1	1059.9	106.0						
Iron	200000	200000	202170.5	208089.7	104.0						
Lead	1000	1000	-11.7	1012.0	101.2						
Magnesium	100000	100000	106670.6	105141.9	105.1						
Manganese	1000	1000	0.0	968.3	96.8						
Molybdenum			3.2	2.9							
Nickel	1000	1000	0.3	1006.1	100.6						
Potassium			-18.4	-26.3							
Selenium	1000	1000	32.4	1083.4	108.3						
Silicon			1.3	3.0							
Silver	1000	1000	-1.1	1115.0	111.5						
Sodium			-68.8	-72.5							
Strontium			3.0	3.0							
Thallium	1000	1000	0.9	971.9	97.2						
Tin			-16.9	-20.3							
Titanium			1.3	1.9							
Vanadium	1000	1000	-2.3	990.0	99.0						
Zinc	1000	1000	3.5	990.4	99.0						

ZP35 : 00000



# IDLs and ICP Linear Ranges



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP35

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
Arsenic	AS	ICP	OPTIMA ICP 2	197.20		10	50.0	4/1/2012	30000.0	6/10/2014

# ICP Interement Correction Factors



CLIENT: Geoenigneers

PROJECT: Gas Works Park-Play

SDG: ZP35

IEC DATE: 10/27/2014

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	AL	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Antimony	206.84	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	14.0198100	0.0000000	0.0000000
Arsenic	188.98	0.0000000	0.0000000	0.0000000	0.0000000	0.0764100	0.0000000	-1.0860250	1.6016350	0.0000000	0.0000000
Barium	233.53	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.1911190	0.0000000	0.0000000	0.1532030
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0098370	0.0000000	0.0000000
Boron	249.67	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	2.1785010	0.0000000	0.0000000	0.0000000
Cadmium	228.80	0.0000000	5.7686950	0.0000000	0.0000000	0.0000000	0.0000000	0.1134910	0.0000000	0.0000000	0.0000000
Calcium	317.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Chromium	267.72	0.0000000	0.0000000	0.0000000	0.0000000	0.0105610	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Cobalt	228.62	0.0000000	0.0000000	0.0853470	0.0000000	0.0000000	0.0000000	0.0000000	-0.0421260	0.0000000	-0.0413720
Copper	324.75	0.0000000	0.0000000	0.0000000	0.0000000	0.0021980	0.0000000	-0.1622490	-0.0152760	0.0000000	-0.0447290
Iron	273.96	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.9508650	0.0000000	0.0000000
Lead	220.35	-0.2096680	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-1.7516130	1.5325290	0.0499030
Magnesium	279.08	0.0000000	0.0000000	0.0000000	0.0000000	0.1119060	0.0000000	-1.6746270	-1.1710960	0.0000000	0.6610200
Manganese	257.61	0.0060150	0.0000000	0.0000000	0.0000000	0.0041750	0.0000000	0.0137770	0.0000000	0.0000000	-0.0046620
Molybdenum	202.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0156160	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Nickel	231.60	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Selenium	196.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.5290750	0.0000000	0.0000000	0.0000000
Silicon	288.16	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-3.5634780	0.0000000	-0.5863590	0.0000000	0.0000000
Silver	328.07	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Sodium	589.59	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium	190.80	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	6.1573050	0.4310930	0.0000000	-0.1254120
Tin	189.93	0.0000000	0.0000000	0.0000000	0.0000000	-0.1475120	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Titanium	334.90	0.0000000	0.0000000	0.0000000	0.0000000	0.0797880	0.0000000	0.0000000	0.19655610	0.0000000	0.0000000
Vanadium	292.40	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-3.7864850	0.0000000	0.0538250
Zinc	206.20	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

FORM XI



# ICP Interelement Correction Factors



CLIENT: Geoenigneers

PROJECT: Gas Works Park-Play

SDG: ZP35

IEC DATE: 10/27/2014

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PB	SB	TI	TL	V	ZN
Aluminum	308.22	0.000000	0.000000	15.2105350	0.000000	0.000000	0.000000	1.9181250	0.000000	14.9692830	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.6574750	0.000000	-4.0497020	0.000000
Arsenic	188.98	0.000000	0.000000	3.6569730	0.000000	0.000000	0.000000	-26.2017890	0.000000	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.1112510	0.000000	0.000000	0.000000	0.000000	0.2137080	0.000000
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0122940	0.000000	0.2849470	0.000000
Boron	249.67	0.000000	0.000000	-1.1347080	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	0.000000	0.000000	-0.9151470	0.000000	0.000000	0.000000	0.000000	0.0642140	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.1073910	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.3457620	0.000000
Cobalt	228.62	0.000000	0.000000	-0.1331780	0.1621790	0.000000	0.000000	1.7359160	0.000000	0.000000	0.000000
Copper	324.75	0.0055170	0.000000	0.3194440	0.000000	0.000000	0.000000	0.1761040	0.000000	0.000000	0.000000
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	7.5009230	0.000000
Lead	220.35	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Magnesium	279.08	0.000000	0.000000	-5.2575470	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Manganese	257.61	0.000000	0.000000	0.000000	0.000000	-0.1944900	0.000000	0.000000	0.000000	-0.0205160	0.000000
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	-0.5357680	0.000000	0.4509940	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.6217450	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.2780320	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Thallium	190.80	0.000000	0.000000	-1.4387970	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.0427690	0.000000	0.000000	0.000000	3.7649150	0.000000
Titanium	334.90	0.000000	0.000000	0.9643590	0.000000	0.000000	0.000000	-0.2280960	0.000000	0.000000	0.000000
Vanadium	292.40	0.000000	-0.1516390	-0.4437580	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Zinc	206.20	0.000000	0.000000	0.2610130	0.000000	-0.0494060	0.000000	0.5348410	0.000000	0.000000	0.000000



# Preparation Log



CLIENT: Geoengineers  
PROJECT: Gas Works Park-Play  
SDG: ZP35

ANALYSIS METHOD: ICP  
ARI PREP CODE: SWC  
PREPDATE: 12/23/2014

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
PAI-2-12.5-13.0	ZP35A	1.026	0.0	50.0
PAI-2-12.5-13.0D	ZP35ADUP	1.027	0.0	50.0
PAI-2-12.5-13.0S	ZP35ASPK	1.024	0.0	50.0
PAI-9-9.5-10.0	ZP35C	1.079	0.0	50.0
PBS	ZP35MB1	1.000	0.0	50.0
LCSS	ZP35MB1SPK	1.000	0.0	50.0







# Analysis Run Log



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZP35

INSTRUMENT ID: OPTIMA ICP 2

RUNID: IP122971 METHOD: ICP

START DATE: 12/29/2014

END DATE: 12/29/2014

CLIENT ID	ARI ID	DIL. TIME	%R	AG	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN		
CCB	CCB6	1.00	14582																															

Total Solids

ARI Job ID: ZP35

Extractions Total Solids-extts  
Data By: Susan D. Dunning  
Created: 12/18/14

Worklist: 2601  
Analyst: SDRD  
Comments:

Oven ID: \_\_\_\_\_

Balance ID: \_\_\_\_\_

Samples In:            Date: \_\_\_\_\_ Time: \_\_\_\_\_ Temp: \_\_\_\_\_ Analyst: \_\_\_\_\_

Samples Out:           Date: \_\_\_\_\_ Time: \_\_\_\_\_ Temp: \_\_\_\_\_ Analyst: \_\_\_\_\_

	ARI ID CLIENT ID	Tare Wt (g)	Wet Wt (g)	Dry Wt (g)	% TS	Dcnt	5g	10g	12.5g
1.	ZP35B 14-27576 PAI-2-18.0-18.5	1.18	9.08	6.65	69.2	No	7.23	14.45	18.06
2.	ZP35D 14-27578 PAI-10-29.5-30.0	1.20	10.49	9.77	92.2	No	5.42	10.85	13.56
3.	ZP35E 14-27579 PAI-12-19.5-20.0	1.19	9.88	7.41	71.6	No	6.98	13.97	17.46



Extractions Total Solids-exttts  
Data By: Susan D. Dunning  
Created: 12/18/14

Worklist: 2601  
Analyst: SDRD  
Comments:

Oven ID: 815

Balance ID: B334705934

Samples In: Date: 12/19/14 Time: 14:47 Temp: 102 Analyst: TH

Samples Out: Date: 12/21/14 Time: 14:35 Temp: 105 Analyst: TH

ARI ID CLIENT ID	Tare Wt (g)	Wet Wt (g)	Dry Wt (g)	% TS	Dcnt	5g	10g	12.5g
1. ZP35B 14-27576 PAI-2-18.0-18.5	<u>1.18</u>	<u>9.08</u>	<u>5.65</u> <u>6.65</u>		No			
2. ZP35D 14-27578 PAI-10-29.5-30.0	<u>1.20</u>	<u>10.49</u>	<u>9.77</u>		No			
3. ZP35E 14-27579 PAI-12-19.5-20.0	<u>1.19</u>	<u>9.88</u>	<u>7.41</u>		No			

Solids Data Entry Report  
Date: 12/23/14

Checked by: CS Date: 12/23/14  
Data Analyst: DM

Solids Determination performed on 12/22/14 by DM

JOB	SAMPLE	CLIENTID	TAREWEIGHT	SAMPDISH	DRYWEIGHT	SOLIDS
ZP35	A	PAI-2-12.5-13.0	1.042	10.426	6.380	56.88
ZP35	C	PAI-9-9.5-10.0	1.013	10.786	9.721	89.10

ZP35: 00109



# Total Solids Bench Sheet

Laboratory Section Mx1015

Oven Identification: 07 Balance ID: 068755

Samples in Oven: Date: 12-22-14 Time: 1340 Temp: 105°C Analyst: DM

Removed from Oven: Date: 12-23-14 Time: 0635 Temp: 104°C Analyst: DM

ARI Sample ID	Tare Weight (g)	Tare + Sample Wet (g)	Tare + Sample Dry (g)	Date & Time Last Weight	Final Weighting >12 hrs <sup>1</sup>
ZP14 A	0.974	10.557	9.361	-	✓
" B	1.029	10.390	9.864	-	✓
" C	1.043	10.690	10.025	-	✓
" D	1.049	10.727	10.038	-	✓
ZP06 A	1.011	5.918	1.884	-	✓
ZP35 A	1.042	10.426	6.980	-	✓
" C	1.013	10.786	9.721	-	✓
ZP16 A	1.019	10.177	8.635	-	✓
" B	1.010	10.041	8.436	-	✓
" C	1.002	10.597	6.866	-	✓
" E	0.979	10.429	6.318	-	✓
" I	1.012	10.142	8.446	-	✓
ZP06 A	0.997	10.545	7.063	-	✓
ZP11 A	1.049	10.848	7.924	-	✓
" B	0.989	10.831	7.615	-	✓
" C	1.019	10.500	8.671	-	✓
" D	1.022	10.407	6.775	-	✓
" E	0.981	10.717	8.184	-	✓
" F	1.068	10.351	9.475	-	✓
" G	1.052	10.209	9.191	-	✓
" H	1.012	10.216	8.563	-	✓
" I	1.058	10.241	9.220	-	✓
" J	1.022	10.533	9.534	-	✓

1) Place a check mark in this column if samples have dried > 12 but < 24 hours. When samples have been at 104°C < 12 hours, constant weight must be verified as described in SOP 10023S. Use a 2<sup>nd</sup> bench sheet for additional weightings.



Volatile Raw Data  
Preparation Log

ARI Job ID: ZP35



Analytical Resources,  
Incorporated  
Analytical Chemists and  
Consultants

# VOA Method 5035 Extraction Bench Sheet

(8260C, 8260C-SIM, 8021B, NWTPH-GX, AK-101, TPH-G, VPH, TCLP-ZHE)

ARI Project No.

Client ID

Prep/Extraction Date

MeOH Lot No.

Analyst

2015

1-19-17

1/1

Lab ID	Vial No.	Preservative		Method 5035 Sample Weight					MeOH Spilt Volume (µL)	Comments
		NaHSO <sub>3</sub>	CH <sub>3</sub> OH	Lot #	Vial Weight (g)	Tare (from vial) (g)	Sample Weight (g)	Extract Volume (mL)		
1	2-P150		-	05279	31.07	28.507	2.567	5ml	100µL	
2	2-P150		-	F	33.50	28.675	4.725	↓	↓	
3	2-P150		-				5.06	↓	5µL	
4										
5										
6										
7										
8										
9										
10										
11										
12										
13										
14										
15										
16										
17										
18										
19										
20										
21										
22										
23										
				Balance ID:						

2015 : 00112



**Analytical Resources, Incorporated**  
Analytical Chemists and Consultants

January 16, 2015

Zanna Satterwhite  
GeoEngineers, Inc.  
Plaza 600 Building  
600 Stewart Street, Suite 1700  
Seattle, WA 98101

**RE: Client Project: Gas Works Park-Play Area Investigation, 0186-846-01 Task 1520**  
**ARI Job No.: ZR94**

Dear Zanna:

Please find enclosed the Chain of Custody record (COC), sample receipt documentation, and the final data package for samples from the project referenced above.

Sample receipt and details of these analyses are discussed in the Case Narrative.

An electronic copy of this package will remain on file with ARI. Should you have any questions or problems, please feel free to contact me at your convenience.

Sincerely,

ANALYTICAL RESOURCES, INC.

Cheronne Oreiro  
Project Manager  
(206) 695-6214  
[cheronneo@arilabs.com](mailto:cheronneo@arilabs.com)  
[www.arilabs.com](http://www.arilabs.com)

cc: eFile: ZR94

Enclosures



Chain of Custody Documentation

ARI Job ID: ZR94







# Cooler Receipt Form

ARI Client: GeoEngineers  
COC No(s): \_\_\_\_\_ (NA)  
Assigned ARI Job No: 215

Project Name: Gas Works Park - PAI  
Delivered by: Fed-Ex UPS Courier Hand Delivered Other: \_\_\_\_\_  
Tracking No: \_\_\_\_\_ (NA)

**Preliminary Examination Phase:**

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES (NO)  
Were custody papers included with the cooler? YES NO  
Were custody papers properly filled out (ink, signed, etc.) YES NO  
Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry) \_\_\_\_\_  
Time: 1350 5.2 5.6 \_\_\_\_\_  
If cooler temperature is out of compliance fill out form 00070F Temp Gun ID#: 9087795

Cooler Accepted by: AV Date: 12/15/14 Time: 1350

**Complete custody forms and attach all shipping documents**

**Log-In Phase:**

Was a temperature blank included in the cooler? YES (NO)  
What kind of packing material was used? ... Bubble Wrap (Wet Ice) Gel Packs (Baggies) Foam Block Paper Other: \_\_\_\_\_  
Was sufficient ice used (if appropriate)? NA YES NO  
Were all bottles sealed in individual plastic bags? YES (NO)  
Did all bottles arrive in good condition (unbroken)? YES NO  
Were all bottle labels complete and legible? YES NO  
Did the number of containers listed on COC match with the number of containers received? YES NO  
Did all bottle labels and tags agree with custody papers? YES NO  
Were all bottles used correct for the requested analyses? YES NO  
Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA YES NO  
Were all VOC vials free of air bubbles? NA YES (NO)  
Was sufficient amount of sample sent in each bottle? YES NO  
Date VOC Trip Blank was made at ARI... NA 12/11/14  
Was Sample Split by ARI: (NA) YES Date/Time: \_\_\_\_\_ Equipment: \_\_\_\_\_ Split by: \_\_\_\_\_

Samples Logged by: AV Date: 12/16/14 Time: 1250

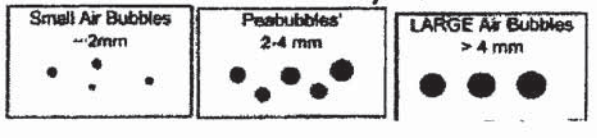
**\*\* Notify Project Manager of discrepancies or concerns \*\***

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

**Additional Notes, Discrepancies, & Resolutions:**

~~FB1 = MW-365-14~~ TBI = MW360-141215 TBI2 = MW-365-141215  
TBI2 = 1Lg MW-360-141215 = 2Lg  
Placed FF sulfides on hold.  
all other COMV requests are FF.

By: AV Date: 12/16/14



Small → "sm" (< 2 mm)  
Peabubbles → "pb" (2 to < 4 mm)  
Large → "lg" (4 to < 6 mm)  
Headspace → "hs" (> 6 mm)

TOC + DOC requested on COC, no bottles received.





ARI Job No: ZP15

Inquiry Number: NONE  
 Analysis Requested: 12/16/14  
 Contact: Satterwhite, Zanna  
 Client: Geoenigneers  
 Logged by: AV  
 Sample Set Used: Yes-481  
 Validatable Package: Lv4  
 Deliverables:

PC: Cheronne  
 VTSR: 12/15/14

Project #: 0186-846-01 Task 1520  
 Project: Gas Works Park-Paly Area Investigat  
 Sample Site:  
 SDG No:  
 Analytical Protocol: In-house

LOGNUM	ARI ID	CLIENT ID	CN	WAD	NH3	COD	FOG	MET	PHEN	PHOS	TKN	NO23	TOC	S2	TPHD	Fe2+	DMET DOC	ADJUSTED LOT	AMOUNT	DATE/BY	
			>12	>12	<2	<2	<2	<2	<2	<2	<2	<2	<2	>9	<2	<2	FLT	TO	ADDED		
14-27518	ZP15A	MW-36D-141215						DIS					Fail	Fail		*	Y				
14-27519	ZP15B	MW-36S-141215						DIS					Fail	Fail		*	Y				

\* Lab to determine ferrous iron preservation  
 \*\* Sulfide preserved with ZnOAC, lab to adjust pH.

Checked By AV Date 12/16/14

Case Narrative, Data Qualifiers, Control Limits

ARI Job ID: ZR94



## Case Narrative

**Client: GeoEngineers, Inc.**

**Project: Gas Works Park-Play Area Investigation, 0186-846-01 Task 1520**

**ARI Job Nos.: ZO41 & ZO68**

### Sample Receipt

The preserved metals volume for one water sample was removed from archive on January 8, 2015. The preserved volume was analyzed for TSS in duplicate, as requested. The TSS filters were retained and digested for metals. For details regarding sample receipt, please refer to the Cooler Receipt Form.

### Metals by SW6010C (Filters)

The sample and associated laboratory QC were digested and analyzed within method recommended holding times.

Iron was present in the filter blank at a level that was greater than the reporting limit. No corrective action was taken.

The duplicate RPD of iron was outside the control limit for the filter sample **MW-36D-141215**. All relevant data have been flagged with a "\*" qualifier on the Form VI. No further corrective action was taken.

### TSS by SM2540D

The water sample was prepared and analyzed outside the method recommended holding time of seven days.

The method blank was clean at the reporting limit. The LCS percent recovery was within control limits.

The replicate RPD was within control limits.



# Sample ID Cross Reference Report



ARI Job No: ZR94  
Client: Geoengineers  
Project Event: 0186-846-01  
Project Name: Gas Works Park-Play Area Investigat

Sample ID	ARI Lab ID	ARI LIMS ID	Matrix	Sample Date/Time	VTSR
1. MW-36D-141215	ZR94A	15-234	Water	12/15/14 10:00	12/15/14 13:50
2. MW-36D-141215	ZR94B	15-235	Filter	12/15/14 10:00	12/15/14 13:50

## Analytical Method Information

Analyte	MDL	Reporting	Surrogate	Duplicate	Matrix Spike		Blank Spike / LCS	
		Limit	%R	RPD	%R	RPD	%R	RPD
<b>Met 6010C in Solid (EPA 6010C)</b>								
Preservation: Cool <6°C								
Container: Glass WM, Clear, 4 oz								
Amount Required: 100 g								
Hold Time: 180 days								
Aluminum	0.757	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Antimony	0.628	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Arsenic	0.333	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Barium	0.133	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Beryllium	0.0160	0.100 mg/kg		20	75 - 125	20	80 - 120	20
Boron	0.739	2.00 mg/kg		20	75 - 125	20	80 - 120	20
Cadmium	0.0180	0.200 mg/kg		20	75 - 125	20	80 - 120	20
Calcium	1.13	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Chromium	0.124	0.500 mg/kg		20	75 - 125	20	80 - 120	20
Cobalt	0.0270	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Copper	0.0920	0.200 mg/kg		20	75 - 125	20	80 - 120	20
Iron	0.750	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Lead	0.155	2.00 mg/kg		20	75 - 125	20	80 - 120	20
Magnesium	0.961	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Manganese	0.0280	0.100 mg/kg		20	75 - 125	20	80 - 120	20
Molybdenum	0.0790	0.500 mg/kg		20	75 - 125	20	80 - 120	20
Nickel	0.386	1.00 mg/kg		20	75 - 125	20	80 - 120	20
Potassium	6.57	50.0 mg/kg		20	75 - 125	20	80 - 120	20
Selenium	0.499	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Silver	0.0430	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Sodium	1.14	50.0 mg/kg		20	75 - 125	20	80 - 120	20
Sodium-1	114	5000 mg/kg		20	75 - 125	20	80 - 120	20
Strontium	0.00900	0.100 mg/kg		20	75 - 125	20	80 - 120	20
Thallium	0.310	5.00 mg/kg		20	75 - 125	20	80 - 120	20
Tin	0.141	1.00 mg/kg		20	75 - 125	20	80 - 120	20
Titanium	0.211	0.500 mg/kg		20	75 - 125	20	80 - 120	20
Vanadium	0.0270	0.300 mg/kg		20	75 - 125	20	80 - 120	20
Zinc	0.145	1.00 mg/kg		20	75 - 125	20	80 - 120	20

Analytical Method Information

Analyte	MDL	Reporting Limit	Surrogate %R	Duplicate RPD	Matrix Spike %R	Matrix Spike RPD	Blank Spike / LCS %R	Blank Spike / LCS RPD
<b>Solids, Total Suspended SM 2540 D-97 in Water (SM 2540 D-97)</b>								
Preservation: Cool <6°C								
Container: HDPE NM, 1000 mL			Amount Required: 1000 mL			Hold Time: 7 days		
Suspended Solids		1.00 mg/L		20				



Metals Analysis  
Report and Summary QC Forms

ARI Job ID: ZR94

# Cover Page

INORGANIC ANALYSIS DATA PACKAGE



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZR94

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
MW-36D-141215	ZR94B	15-235	
MW-36D-141215D	ZR94BDUP	15-235	
PBS	ZR94MB1	15-235	

Were ICP interelement corrections applied ?                      Yes/No    YES  
Were ICP background corrections applied ?                      Yes/No    YES  
If yes - were raw data generated before  
application of background corrections ?                      Yes/No    NO

Comments: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature: 

Name: Jay Kuhn

Date: 1/14/15

Title: Inorganics Director

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**


Page 1 of 1

Sample ID: MW-36D-141215  
SAMPLE

Lab Sample ID: ZR94B

LIMS ID: 15-235

Matrix: Filter

Data Release Authorized: 

Reported: 01/14/15

QC Report No: ZR94-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01

Date Sampled: 12/15/14

Date Received: 12/15/14

Percent Total Solids: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	µg/Sample	Q
6010C	01/12/15	6010C	01/13/15	7440-38-2	Arsenic	20	7,500	
6010C	01/12/15	6010C	01/13/15	7439-89-6	Iron	20	50	

U-Analyte undetected at given RL

RL-Reporting Limit



**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

Page 1 of 1

Sample ID: MW-36D-141215

DUPLICATE

Lab Sample ID: ZR94B

LIMS ID: 15-235

Matrix: Filter

Data Release Authorized: 

Reported: 01/14/15

QC Report No: ZR94-Geoengineers

Project: Gas Works Park-Play Area Investigat

0186-846-01

Date Sampled: 12/15/14

Date Received: 12/15/14

**MATRIX DUPLICATE QUALITY CONTROL REPORT**

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010C	7,500	7,640	1.8%	+/- 20%	
Iron	6010C	50	130	88.9%	+/- 20	L*

Reported in µg/Sample

\*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

**INORGANICS ANALYSIS DATA SHEET**

**TOTAL METALS**

**Sample ID: METHOD BLANK**

Page 1 of 1

Lab Sample ID: ZR94MB

QC Report No: ZR94-Geoengineers

LIMS ID: 15-235

Project: Gas Works Park-Play Area Investigat

Matrix: Filter

0186-846-01

Data Release Authorized: 

Date Sampled: NA

Reported: 01/14/15

Date Received: NA

Percent Total Solids: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	µg/Sample	Q
6010C	01/12/15	6010C	01/13/15	7440-38-2	Arsenic	2	2	U
6010C	01/12/15	6010C	01/13/15	7439-89-6	Iron	2	47	

U-Analyte undetected at given RL

RL-Reporting Limit

# Calibration Verification



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZR94

UNITS: ug/L

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	ICP	IP011371	2000.0	2033.26	101.7	2000.0	2007.09	100.4	2059.02	103.0	2070.00	103.5	2066.19	103.3	2088.81	104.4
Iron	FE	ICP	IP011371	2000.0	2042.80	102.1	2000.0	2047.96	102.4	2043.03	102.2	2060.58	103.0	2073.09	103.7	2077.70	103.9

Control Limits: Mercury 80-120; Other Metals 90-110



# Calibration Verification



CLIENT: Geoenineers

PROJECT: Gas Works Park-Play

SDG: ZR94

UNITS: ug/L

ANALYTE	EL	M	RUN	CCVTV	CCV6	%R	CCV7	%R	CCV8	%R	CCV9	%R	CCV10	%R	CCV11	%R
Arsenic	AS	ICP	IP011371	2000.0	2087.44	104.4										
Iron	FE	ICP	IP011371	2000.0	2085.69	104.3										

Control Limits: Mercury 80-120; Other Metals 90-110

**CRDL Standard**

CLIENT: Geoenineers

PROJECT: Gas Works Park-Play

SDG: ZR94



UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Arsenic	AS	ICP	IF011371	50.0	50.12	100.2	53.24	106.5								
Iron	FE	ICP	IF011371	50.0	52.49	105.0	53.14	106.3								

ZR94 : 00018

Control Limits: no control limits have been established by the EPA at this time.

# Calibration Blanks

CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZR94



UNITS:ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Arsenic	AS	ICP	IP011371	10.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U
Iron	FE	ICP	IP011371	100.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U



# Calibration Blanks



CLIENT: Geengineers

PROJECT: Gas Works Park-Play

SDG: ZR94

UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	CCB6	CCB7	CCB8	CCB9	CCB10	CCB11	C
Arsenic	AS	ICP	IP011371	10.0	50.0	50.0						U
Iron	FE	ICP	IP011371	100.0	50.0	50.0						U

# ICP Interference Check Sample



CLIENT: Geoenigneers

ICS SOURCE: I.V.

PROJECT: Gas Works Park-Play

RUNID: IP011371

SDG: ZR94

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSA1 %R	ICSA2	ICSA2 %R	ICSA3	ICSA3 %R	ICSA4	ICSA4 %R
Aluminum	200000	200000	200194.3	200788.3	203582.5	204001.6	102.0	204001.6	102.0	204001.6
Antimony	1000	1000	15.2	1050.1	16.3	1089.0	108.9	1089.0	108.9	1089.0
Arsenic	1000	1000	22.7	1043.1	22.5	1078.8	107.9	1078.8	107.9	1078.8
Barium	1000	1000	-3.2	1026.6	-2.8	1056.8	105.7	1056.8	105.7	1056.8
Beryllium	1000	1000	0.1	1008.2	0.1	1015.6	101.6	1015.6	101.6	1015.6
Boron			-6.6	-4.1	-6.2	-5.6				
Cadmium	1000	1000	0.6	1053.2	1.1	1068.1	106.8	1068.1	106.8	1068.1
Calcium	100000	100000	101014.4	100804.5	102372.0	102295.0	102.3	102295.0	102.3	102295.0
Chromium	1000	1000	-1.9	1041.6	-2.3	1068.2	106.8	1068.2	106.8	1068.2
Cobalt	1000	1000	2.1	990.9	2.0	1018.7	101.9	1018.7	101.9	1018.7
Copper	1000	1000	0.5	1089.0	0.5	1102.7	110.3	1102.7	110.3	1102.7
Iron	200000	200000	196967.1	197129.7	198859.7	200589.4	100.3	200589.4	100.3	200589.4
Lead	1000	1000	-10.0	991.9	-9.3	1023.3	102.3	1023.3	102.3	1023.3
Magnesium	100000	100000	103068.5	99639.1	105581.2	101480.2	101.5	101480.2	101.5	101480.2
Manganese	1000	1000	0.4	979.2	0.6	996.7	99.7	996.7	99.7	996.7
Molybdenum			2.2	2.4	2.1	2.2				
Nickel	1000	1000	1.1	984.5	0.1	1012.2	101.2	1012.2	101.2	1012.2
Potassium			-12.6	-2.7	0.4	-1.4				
Selenium	1000	1000	29.4	1064.2	37.4	1095.0	109.5	1095.0	109.5	1095.0
Silicon	1000	1000	12.7	6.5	10.5	6.7				
Silver	1000	1000	-0.8	1084.5	-0.8	1102.4	110.2	1102.4	110.2	1102.4
Sodium			6.3	4.5	8.5	10.1				
Strontium			3.0	2.9	3.0	3.0				
Thallium	1000	1000	-3.2	944.5	-0.8	974.2	97.4	974.2	97.4	974.2
Tin			-20.3	-21.9	-21.1	-21.6				
Titanium			0.7	0.5	1.1	0.6				
Vanadium	1000	1000	-3.6	1005.4	-3.3	1022.0	102.2	1022.0	102.2	1022.0
Zinc	1000	1000	3.0	971.0	3.2	994.2	99.4	994.2	99.4	994.2

ZR94 : 00021

# IDLs and ICP Linear Ranges



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

SDG: ZR94

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
Arsenic	AS	ICP	OPTIMA ICP 2	197.20		10	50.0	4/1/2012	30000.0	6/10/2014
Iron	FE	ICP	OPTIMA ICP 2	259.94		100	50.0	4/1/2012	250000.0	6/10/2014



# ICP Interelement Correction Factors



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

IEC DATE: 10/27/2014

SDG: ZR94

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	AL	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	14.0198100	0.000000	0.000000
Arsenic	188.98	0.000000	0.000000	0.000000	0.000000	0.0764100	0.000000	-1.0860250	1.6016350	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.1911190	0.000000	0.000000	0.1532030
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0098370	0.000000	0.000000
Boron	249.67	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	2.1785010	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	5.7686950	0.000000	0.000000	0.000000	0.000000	0.1134910	0.000000	0.000000	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.000000	0.000000	0.000000	0.000000	0.0105610	0.000000	0.000000	0.000000	0.000000	0.000000
Cobalt	228.62	0.000000	0.000000	0.0853470	0.000000	0.000000	0.000000	0.000000	-0.0421260	0.000000	0.000000
Copper	324.75	0.000000	0.000000	0.000000	0.000000	0.0021980	0.000000	-0.1622490	-0.0152760	0.000000	-0.0447290
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.9508650	0.000000	0.000000
Lead	220.35	-0.2096680	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-1.7516130	1.5325290	0.0499030
Magnesium	279.08	0.000000	0.000000	0.000000	0.000000	0.1119060	0.000000	-1.6746270	-1.1710960	0.000000	0.6610200
Manganese	257.61	0.0060150	0.000000	0.000000	0.000000	0.0041750	0.000000	0.0137770	0.000000	0.000000	-0.0046620
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.0156160	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.5290750	0.000000	0.000000	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	-3.5634780	0.000000	-0.5863590	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Thallium	190.80	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	6.1573050	0.000000	0.000000	0.000000
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.1475120	0.000000	0.000000	0.000000	0.000000	0.000000
Titanium	334.90	0.000000	0.000000	0.000000	0.000000	0.0797880	0.000000	0.000000	0.1965610	0.000000	0.000000
Vanadium	292.40	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-3.7864850	0.000000	0.0538250
Zinc	206.20	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000



# ICP Interelement Correction Factors



CLIENT: Geoengineers

PROJECT: Gas Works Park-Play

IEC DATE: 10/27/2014

SDG: ZR94

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PB	SB	TI	TL	V	ZN
Aluminum	308.22	0.000000	0.000000	15.2105350	0.000000	0.000000	0.000000	1.9181250	0.000000	14.9692830	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.6574750	0.000000	-4.0497020	0.000000
Arsenic	188.98	0.000000	0.000000	3.6569730	0.000000	0.000000	0.000000	-26.2017890	0.000000	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.1112510	0.000000	0.000000	0.000000	0.000000	0.2137080	0.000000
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0122940	0.000000	0.2849470	0.000000
Boron	249.67	0.000000	0.000000	-1.1347080	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	0.000000	0.000000	-0.9151470	0.000000	0.000000	0.000000	0.000000	0.0642140	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.1073910	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.3457620	0.000000
Cobalt	228.62	0.000000	0.000000	-0.1331780	0.1621790	0.000000	0.000000	1.7359160	0.000000	0.000000	0.000000
Copper	324.75	0.0055170	0.000000	0.3194440	0.000000	0.000000	0.000000	0.1761040	0.000000	0.000000	0.000000
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	7.5009230	0.000000
Lead	220.35	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Magnesium	279.08	0.000000	0.000000	-5.2575470	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Manganese	257.61	0.000000	0.000000	0.000000	0.000000	-0.1944900	0.000000	0.000000	0.000000	-0.0205160	0.000000
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	-0.5357680	0.000000	0.4509940	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.6217450	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.2780320	0.000000
Thallium	190.80	0.000000	0.000000	-1.4387970	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	81.4212140
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.0427690	-0.5371860	-0.2280960	0.000000	0.000000	0.000000
Titanium	334.90	0.000000	0.000000	0.9643590	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Vanadium	292.40	0.000000	-0.1516390	-0.4437580	0.000000	0.000000	0.000000	0.5348410	0.000000	0.000000	0.000000
Zinc	206.20	0.000000	0.000000	0.2610130	0.000000	-0.0494060	0.000000	0.000000	0.000000	0.000000	0.000000

ZR94 : 00024

# Preparation Log



CLIENT: Geoengineers  
PROJECT: Gas Works Park-Play  
SDG: ZR94

ANALYSIS METHOD: ICP  
ARI PREP CODE: PHN  
PREPDATE: 1/12/2015

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
MW-36D-141215	ZR94B	1.000	0.0	50.0
MW-36D-141215D	ZR94BDUP	1.000	0.0	50.0
PBS	ZR94MB1	1.000	0.0	50.0



