

Jim Harris

Utah DEQ DERR

168 N. 1950 W., 1st Floor

**AMERICAN** 

Salt Lake City, Ut 84116

WEST

TEL: (801) 536-4145

ANALYTICAL LABORATORIES

RE: Red Butte

Dear Jim Harris:

Lab Set ID: 1011529

463 West 3600 South

Salt Lake City, UT

American West Analytical Laboratories received 6 sample(s) on 11/30/2010 for the analyses presented in the following report.

84115

All analyses were performed in accordance to The NELAC Institute protocols unless noted otherwise. American West Analytical Laboratories is certified by The NELAC Institute in Utah and Texas; and is state certified in Colorado and Idaho. Certification document is available upon request. If you have any questions or concerns regarding this report please feel free to call.

(801) 263-8686

Toll Free (888) 263-8686

Fax (801)263-8687

email: awal@awal-labs.com

The abbreviation "Surr" found in organic reports indicates a surrogate compound that is intentionally added by the laboratory to determine sample injection, extraction, and/or purging efficiency. The "Reporting Limit" found on the report is equivalent to the practical quantitaion limit (PQL). This is the minimum concentration that can be reported by the method referenced and the sample matrix. The reporting limit must not be confused with any regulatory limit. Analytical results are reported to three significant figures for quality control and calculation purposes.

Kyle F. Gross Laboratory Director

Jose Rocha

Thank You,

QA Officer

Approved by: \_\_\_\_\_ Laboratory Director or designee

Report Date: 12/6/2010 Page 1 of 162



**WEST** 

# **INORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001

Client Sample ID: 4992095 Red Butte AB Garden

**Collection Date:** 11/30/2010 0950h **Received Date:** 11/30/2010 1538h

ANALYTICAL LABORATORIES

463 West 3600 South Salt Lake City, UT 84115

Analytical Results	Units	Date Prepared	Date Analyzed	Method Used	Reporting Limit	Analytical Result	Qual
Chemical Oxygen Demand	mg/L		12/2/2010 1200	h HACH 8000	10.0	< 10.0	
Total Recoverable Petroleum Hydrocarbons	mg/L		12/1/2010 1354	h E1664A-SGT	3.00	< 3.00	

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Kyle F. Gross Laboratory Director

> Jose Rocha QA Officer

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**WEST** 

# **INORGANIC ANALYTICAL REPORT**

**Client:** Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte 1011529-002 Lab Sample ID:

Client Sample ID: 4992085 Red Butte BL Garden

**Collection Date:** 11/30/2010 1030h **Received Date:** 11/30/2010 1538h

**ANALYTICAL** 

LABORATORIES

LABORATORIES	Analytical Results	Units	Date Prepared	Date Analyzed	Method Used	Reporting Limit	Analytical Result	Qual
463 West 3600 South	Chemical Oxygen Demand	mg/L		12/2/2010 1200	h HACH 8000	10.0	< 10.0	
463 West 3600 South Salt Lake City, UT 84115	Total Recoverable Petroleum Hydrocarbons	mg/L		12/1/2010 1354	h E1664A-SGT	3.00	< 3.00	

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> Jose Rocha **QA** Officer

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**WEST** 

# **INORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

**Collection Date:** 11/30/2010 1215h **Received Date:** 11/30/2010 1538h

ANALYTICAL LABORATORIES

463 West 3600 South Salt Lake City, UT 84115

<b>Analytical Results</b>	Units	Date Prepared	Date Analyzed	Method l Used	Reporting Limit	Analytical Result	Qual
Chemical Oxygen Demand	mg/L		12/2/2010 120	00h HACH 8000	10.0	< 10.0	
Total Recoverable Petroleum Hydrocarbons	mg/L		12/1/2010 13:	54h E1664A-SGT	3.00	4.70	

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**WEST** 

# **INORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004

Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

**Collection Date:** 11/30/2010 1405h **Received Date:** 11/30/2010 1538h

ANALYTICAL LABORATORIES

463 West 3600 South Salt Lake City, UT 84115

<b>Analytical Results</b>	Units	Date Prepared	Date Analyzed	Method d Used	Reporting Limit	Analytical Result	Qual
Chemical Oxygen Demand	mg/L		12/2/2010 120	00h HACH 8000	10.0	10.0	
Total Recoverable Petroleum Hydrocarbons	mg/L		12/1/2010 13:	54h E1664A-SGT	3.00	< 3.00	

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# **INORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005

**Client Sample ID:** 4992290 J.R. @ 1700 S. **Collection Date:** 11/30/2010 1430h **Received Date:** 11/30/2010 1538h

ANALYTICAL

LABORATORIES

463 West 3600 South Salt Lake City, UT

Analytical Results	Units	Date Prepared	Date Analyze	ed	Method Used	Reporting Limit	Analytical Result	Qual
Chemical Oxygen Demand	mg/L		12/2/2010 1:	1200h	HACH 8000	10.0	13.0	
Total Recoverable Petroleum Hydrocarbons	mg/L		12/1/2010 1:	1354h I	E1664A-SGT	3.00	< 3.00	

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# **INORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

**Collection Date:** 11/30/2010 1315h **Received Date:** 11/30/2010 1538h

ANALYTICAL LABORATORIES

463 West 3600 South Salt Lake City, UT

Analytical Results	Units	Date Prepared	Date Analyzed	Method Used	Reporting Limit	Analytical Result	Qual
Chemical Oxygen Demand	mg/L		12/2/2010 1200	h HACH 8000	10.0	< 10.0	
Total Recoverable Petroleum Hydrocarbons	mg/L		12/1/2010 1354	h E1664A-SGT	3.00	< 3.00	

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**WEST** 

### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001D

AMERICAN Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/6/2010 0057h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/3/2010 0843h

Method Used: SW8015D

Surr: C36

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-ORO by GC/FID Method 8015D/3510C

Units: mg/L
Dilution Factor: 1
CAS Reporting Analytical
Compound Number Limit Result Qual

Oil Range Organics (ORO) (C28-C36)
0.0515 < 0.0515

10-200

123

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**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002D

Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/6/2010 0115h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/3/2010 0843h

Method Used: SW8015D

LABORATORIES

463 West 3600 South

Salt Lake City, UT

ANALYTICAL

Analytical Results TPH-ORO by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Limit Compound Number Result Qual Oil Range Organics (ORO) (C28-C36) 0.0500 < 0.0500 Surr: C36 10-200 117

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**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003D

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

 Collection Date:
 11/30/2010 1215h
 Analyzed: 12/6/2010 0134h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/3/2010 0843h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-ORO by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Limit Compound Number Result Qual Oil Range Organics (ORO) (C28-C36) 0.0500 < 0.0500 Surr: C36 10-200 121

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**WEST** 

### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004D

Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/6/2010 0152h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/3/2010 0843h

Method Used: SW8015D

Surr: C36

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

ANALYTICAL

Analytical Results TPH-ORO by GC/FID Method 8015D/3510C

Units: mg/L
Dilution Factor: 1
CAS Reporting Analytical
Compound Number Limit Result Qual

Oil Range Organics (ORO) (C28-C36)
0.0500 < 0.0500

10-200

124

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**WEST** 

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# ORGANIC ANALYTICAL REPORT

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005D

**Client Sample ID:** 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/6/2010 0210h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/3/2010 0843h

Method Used: SW8015D

**Analytical Results** 

**LABORATORIES** 

463 West 3600 South Salt Lake City, UT

ANALYTICAL

TPH-ORO by GC/FID Method 8015D/3510C

Units: mg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
Oil Range Organics (ORO) (C28-C36)		0.0500	< 0.0500	
Surr: C36		10-200	122	

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**WEST** 

### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006D

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/6/2010 0228h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/3/2010 0843h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-ORO by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Limit Compound Number Result Qual Oil Range Organics (ORO) (C28-C36) 0.0500 < 0.0500 Surr: C36 10-200 124

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ANALYTICAL

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### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001E

Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/2/2010 1335h

 Received Date:
 11/30/2010 1538h
 Extracted: 11/30/2010 1740h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

Analytical Results TPH-DRO (C10-C28) by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Compound Number Limit Result Qual Diesel Range Organics C10-C28 68476-34-6 0.526 < 0.526 Surr: 4-Bromofluorobenzene 460-00-4 10-190 **77.0** 

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> Jose Rocha QA Officer

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**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002E

AMERICAN Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/2/2010 1439h

 Received Date:
 11/30/2010 1538h
 Extracted: 11/30/2010 1740h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-DRO (C10-C28) by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Compound Number Limit Result Qual Diesel Range Organics C10-C28 68476-34-6 0.500 < 0.500 Surr: 4-Bromofluorobenzene 460-00-4 10-190 68.0

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Kyle F. Gross Laboratory Director

Jose Rocha

QA Officer

Report Date: 12/6/2010 Page 15 of 162



**WEST** 

### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003E

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

 Collection Date:
 11/30/2010 1215h
 Analyzed: 12/2/2010 1500h

 Received Date:
 11/30/2010 1538h
 Extracted: 11/30/2010 1740h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-DRO (C10-C28) by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Compound Number Limit Result Qual Diesel Range Organics C10-C28 68476-34-6 0.500 < 0.500 Surr: 4-Bromofluorobenzene 460-00-4 10-190 54.0

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> Jose Rocha QA Officer

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**WEST** 

### ORGANIC ANALYTICAL REPORT

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004E

Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/2/2010 1521h

 Received Date:
 11/30/2010 1538h
 Extracted: 11/30/2010 1740h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-DRO (C10-C28) by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Compound Number Limit Result Qual Diesel Range Organics C10-C28 68476-34-6 0.510 < 0.510 Surr: 4-Bromofluorobenzene 460-00-4 10-190 48.5

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Kyle F. Gross Laboratory Director

> Jose Rocha QA Officer

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**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005E

**Client Sample ID:** 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/2/2010 1542h

 Received Date:
 11/30/2010 1538h
 Extracted: 11/30/2010 1740h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-DRO (C10-C28) by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Compound Number Limit Result Qual Diesel Range Organics C10-C28 68476-34-6 0.500 < 0.500 Surr: 4-Bromofluorobenzene 460-00-4 10-190 41.8

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> Jose Rocha QA Officer

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**WEST** 

### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006E

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/2/2010 1603h

 Received Date:
 11/30/2010 1538h
 Extracted: 11/30/2010 1740h

Method Used: SW8015D

**LABORATORIES** 

463 West 3600 South

Salt Lake City, UT

84115

ANALYTICAL

Analytical Results TPH-DRO (C10-C28) by GC/FID Method 8015D/3510C

Units: mg/L **Dilution Factor: 1 CAS** Reporting **Analytical** Compound Number Limit Result Qual Diesel Range Organics C10-C28 68476-34-6 0.500 < 0.500 Surr: 4-Bromofluorobenzene 460-00-4 10-190 46.9

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Kyle F. Gross Laboratory Director

> Jose Rocha QA Officer

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Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001F

AMERICAN Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/2/2010 2232h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: SW8270D

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**ANALYTICAL** 

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Acenaphthene	83-32-9	10.5	< 10.5	
	Acenaphthylene	208-96-8	10.5	< 10.5	
	Anthracene	120-12-7	10.5	< 10.5	
	Benz(a)anthracene	56-55-3	10.5	< 10.5	
(801) 263-8686	Benzo(a)pyrene	50-32-8	10.5	< 10.5	
Toll Free (888) 263-8686	Benzo(b)fluoranthene	205-99-2	10.5	< 10.5	
Fax (801)263-8687	Benzo(g,h,i)perylene	191-24-2	10.5	< 10.5	
email: awal@awal-labs.com	Benzo(k)fluoranthene	207-08-9	10.5	< 10.5	
	Chrysene	218-01-9	10.5	< 10.5	
Kyle F. Gross	Dibenz(a,h)anthracene	53-70-3	10.5	< 10.5	
Laboratory Director	Fluoranthene	206-44-0	10.5	< 10.5	
Laboratory Director	Fluorene	86-73-7	10.5	< 10.5	
Jose Rocha	Indeno(1,2,3-cd)pyrene	193-39-5	10.5	< 10.5	
	Phenanthrene	85-01-8	10.5	< 10.5	
QA Officer	Pyrene	129-00-0	10.5	< 10.5	
	C11-C12 Aliphatic hydrocarbons		10.5	< 10.5	
	C13-C16 Aliphatic hydrocarbons		10.5	< 10.5	
	C17-C21 Aliphatic hydrocarbons		10.5	< 10.5	
	C22-C35 Aliphatic hydrocarbons		10.5	< 10.5	
	C11-C13 Alkyl Naphthalenes		10.5	< 10.5	
	Total C12-C22 PAH**		10.5	< 10.5	
	Surr: Terphenyl-d14	1718-51-0	10-199	106	

<sup>\*\* -</sup> This value is a summation of the PAH compounds listed above.



Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002F

AMERICAN Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/2/2010 2300h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

ANALYTICAL Method Used: SW8270D

LABORATORIES

**WEST** 

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
(801) 263-8686	Benzo(a)pyrene	50-32-8	10.0	< 10.0	
Toll Free (888) 263-8686	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
Fax (801)263-8687	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
email: awal@awal-labs.com	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Chrysene	218-01-9	10.0	< 10.0	
Kyle F. Gross	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
Laboratory Director	Fluoranthene	206-44-0	10.0	< 10.0	
Laboratory Director	Fluorene	86-73-7	10.0	< 10.0	
Iona Danka	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Jose Rocha	Phenanthrene	85-01-8	10.0	< 10.0	
QA Officer	Pyrene	129-00-0	10.0	< 10.0	
	C11-C12 Aliphatic hydrocarbons		10.0	< 10.0	
	C13-C16 Aliphatic hydrocarbons		10.0	< 10.0	
	C17-C21 Aliphatic hydrocarbons		10.0	< 10.0	
	C22-C35 Aliphatic hydrocarbons		10.0	< 10.0	
	C11-C13 Alkyl Naphthalenes		10.0	< 10.0	
	Total C12-C22 PAH**		10.0	< 10.0	
	Surr: Terphenyl-d14	1718-51-0	10-199	108	

<sup>\*\*</sup> - This value is a summation of the PAH compounds listed above.



Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003F

AMERICAN Client Sample ID: 4992084 Red Butte @ 1500 E.

 Collection Date:
 11/30/2010 1215h
 Analyzed: 12/2/2010 2327h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

ANALYTICAL Method Used: SW8270D

LABORATORIES
Analytical Results

**WEST** 

SVOA Fractionation by GC/MS Method 8270D/3510C

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
(801) 263-8686	Benzo(a)pyrene	50-32-8	10.0	< 10.0	
Toll Free (888) 263-8686	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
Fax (801)263-8687	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
email: awal@awal-labs.com	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Chrysene	218-01-9	10.0	< 10.0	
Kyle F. Gross	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
Laboratory Director	Fluoranthene	206-44-0	10.0	< 10.0	
Laboratory Director	Fluorene	86-73-7	10.0	< 10.0	
r D 1	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Jose Rocha	Phenanthrene	85-01-8	10.0	< 10.0	
QA Officer	Pyrene	129-00-0	10.0	< 10.0	
	C11-C12 Aliphatic hydrocarbons		10.0	< 10.0	
	C13-C16 Aliphatic hydrocarbons		10.0	< 10.0	
	C17-C21 Aliphatic hydrocarbons		10.0	< 10.0	
	C22-C35 Aliphatic hydrocarbons		10.0	< 10.0	
	C11-C13 Alkyl Naphthalenes		10.0	< 10.0	
	Total C12-C22 PAH**		10.0	< 10.0	
	Surr: Terphenyl-d14	1718-51-0	10-199	110	

<sup>\*\* -</sup> This value is a summation of the PAH compounds listed above.



**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004F

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/2/2010 2355h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: SW8270D

**LABORATORIES** 

**ANALYTICAL** 

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
(801) 263-8686	Benzo(a)pyrene	50-32-8	10.0	< 10.0	
Toll Free (888) 263-8686	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
Fax (801)263-8687	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
email: awal@awal-labs.com	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Chrysene	218-01-9	10.0	< 10.0	
Kyle F. Gross	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
Laboratory Director	Fluoranthene	206-44-0	10.0	< 10.0	
Laboratory Director	Fluorene	86-73-7	10.0	< 10.0	
r D 1	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Jose Rocha	Phenanthrene	85-01-8	10.0	< 10.0	
QA Officer	Pyrene	129-00-0	10.0	< 10.0	
	C11-C12 Aliphatic hydrocarbons		10.0	< 10.0	
	C13-C16 Aliphatic hydrocarbons		10.0	< 10.0	
	C17-C21 Aliphatic hydrocarbons		10.0	< 10.0	
	C22-C35 Aliphatic hydrocarbons		10.0	< 10.0	
	C11-C13 Alkyl Naphthalenes		10.0	< 10.0	
	Total C12-C22 PAH**		10.0	< 10.0	
	Surr: Terphenyl-d14	1718-51-0	10-199	107	

<sup>\*\* -</sup> This value is a summation of the PAH compounds listed above.



Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005F

AMERICAN Client Sample ID: 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/3/2010 0022h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: SW8270D

LABORATORIES

**WEST** 

**ANALYTICAL** 

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
(801) 263-8686	Benzo(a)pyrene	50-32-8	10.0	< 10.0	
Toll Free (888) 263-8686	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
Fax (801)263-8687	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
email: awal@awal-labs.com	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Chrysene	218-01-9	10.0	< 10.0	
Kyle F. Gross	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
Laboratory Director	Fluoranthene	206-44-0	10.0	< 10.0	
Laboratory Director	Fluorene	86-73-7	10.0	< 10.0	
r D 1	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Jose Rocha	Phenanthrene	85-01-8	10.0	< 10.0	
QA Officer	Pyrene	129-00-0	10.0	< 10.0	
	C11-C12 Aliphatic hydrocarbons		10.0	< 10.0	
	C13-C16 Aliphatic hydrocarbons		10.0	< 10.0	
	C17-C21 Aliphatic hydrocarbons		10.0	< 10.0	
	C22-C35 Aliphatic hydrocarbons		10.0	< 10.0	
	C11-C13 Alkyl Naphthalenes		10.0	< 10.0	
	Total C12-C22 PAH**		10.0	< 10.0	
	Surr: Terphenyl-d14	1718-51-0	10-199	113	

<sup>\*\* -</sup> This value is a summation of the PAH compounds listed above.



**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006F

AMERICAN Client Sample ID: 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/3/2010 0049h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: SW8270D

**LABORATORIES** 

**ANALYTICAL** 

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
(801) 263-8686	Benzo(a)pyrene	50-32-8	10.0	< 10.0	
Toll Free (888) 263-8686	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
Fax (801)263-8687	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
email: awal@awal-labs.com	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Chrysene	218-01-9	10.0	< 10.0	
Kyle F. Gross	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
Laboratory Director	Fluoranthene	206-44-0	10.0	< 10.0	
Laboratory Director	Fluorene	86-73-7	10.0	< 10.0	
Iona Danka	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Jose Rocha	Phenanthrene	85-01-8	10.0	< 10.0	
QA Officer	Pyrene	129-00-0	10.0	< 10.0	
	C11-C12 Aliphatic hydrocarbons		10.0	< 10.0	
	C13-C16 Aliphatic hydrocarbons		10.0	< 10.0	
	C17-C21 Aliphatic hydrocarbons		10.0	< 10.0	
	C22-C35 Aliphatic hydrocarbons		10.0	< 10.0	
	C11-C13 Alkyl Naphthalenes		10.0	< 10.0	
	Total C12-C22 PAH**		10.0	< 10.0	
	Surr: Terphenyl-d14	1718-51-0	10-199	107	

<sup>\*\*</sup> - This value is a summation of the PAH compounds listed above.



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001F

Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/2/2010 2232h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1´-Biphenyl	92-52-4	10.5	< 10.5	
01113	1,2,4,5-Tetrachlorobenzene	95-94-3	10.5	< 10.5	
	1,2,4-Trichlorobenzene	120-82-1	10.5	< 10.5	
	1,2-Dichlorobenzene	95-50-1	10.5	< 10.5	
(801) 263-8686	1,3,5-Trinitrobenzene	99-35-4	10.5	< 10.5	
Toll Free (888) 263-8686	1,4-Naphthoquinone	130-15-4	10.5	< 10.5	
Fax (801)263-8687	1,3-Dichlorobenzene	541-73-1	10.5	< 10.5	
email: awal@awal-labs.com	1,3-Dinitrobenzene	99-65-0	10.5	< 10.5	
	1,4-Dichlorobenzene	106-46-7	10.5	< 10.5	
Kyle F. Gross	1,4-Phenylenediamine	106-50-3	10.5	< 10.5	
•	1-Chloronaphthalene	90-13-1	10.5	< 10.5	
Laboratory Director	1-Methylnaphthalene	90-12-0	10.5	< 10.5	
r D 1	1-Naphthylamine	134-32-7	10.5	< 10.5	
Jose Rocha	2,3,4,6-Tetrachlorophenol	58-90-2	10.5	< 10.5	
QA Officer	2,4,5-Trichlorophenol	95-95-4	10.5	< 10.5	
	2,4,6-Trichlorophenol	88-06-2	10.5	< 10.5	
	2,4-Dichlorophenol	120-83-2	10.5	< 10.5	
	2,4-Dimethylphenol	105-67-9	10.5	< 10.5	
	2,4-Dinitrophenol	51-28-5	21.1	< 21.1	
	2,4-Dinitrotoluene	121-14-2	10.5	< 10.5	
	2,6-Dichlorophenol	87-65-0	10.5	< 10.5	
	2,6-Dinitrotoluene	606-20-2	10.5	< 10.5	
	2-Acetylaminofluorene	53-96-3	10.5	< 10.5	
	2-Chloronaphthalene	91-58-7	10.5	< 10.5	
	2-Chlorophenol	95-57-8	10.5	< 10.5	
	2-Methylnaphthalene	91-57-6	10.5	< 10.5	
	2-Methylphenol	95-48-7	10.5	< 10.5	
	2-Naphthylamine	91-59-8	10.5	< 10.5	
	2-Nitroaniline	88-74-4	10.5	< 10.5	

Report Date: 12/6/2010 Page 26 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001F

Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/2/2010 2232h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	2-Nitrophenol	88-75-5	10.5	< 10.5	
	2-Picoline	109-06-8	10.5	< 10.5	
	3&4-Methylphenol		10.5	< 10.5	
	3,3'-Dichlorobenzidine	91-94-1	10.5	< 10.5	
(801) 263-8686	3,3'-Dimethylbenzidine	119-93-7	10.5	< 10.5	
Toll Free (888) 263-8686	3-Methylcholanthrene	56-49-5	10.5	< 10.5	
Fax (801)263-8687	3-Nitroaniline	99-09-2	10.5	< 10.5	
email: awal@awal-labs.com	4,6-Dinitro-2-methylphenol	534-52-1	10.5	< 10.5	
	4-Aminobiphenyl	92-67-1	10.5	< 10.5	
Kyle F. Gross	4-Bromophenyl phenyl ether	101-55-3	10.5	< 10.5	
•	4-Chloro-3-methylphenol	59-50-7	10.5	< 10.5	
Laboratory Director	4-Chloroaniline	106-47-8	10.5	< 10.5	
r D 1	4-Chlorophenyl phenyl ether	7005-72-3	10.5	< 10.5	
Jose Rocha	4-Nitroaniline	100-01-6	10.5	< 10.5	
QA Officer	4-Nitrophenol	100-02-7	10.5	< 10.5	
	5-Nitro-o-toluidine	99-55-8	10.5	< 10.5	
	7,12-Dimethylbenz(a)anthracene	57-97-6	10.5	< 10.5	
	a,a-Dimethylphenethylamine	122-09-8	10.5	< 10.5	
	Acenaphthene	83-32-9	10.5	< 10.5	
	Acenaphthylene	208-96-8	10.5	< 10.5	
	Acetophenone	98-86-2	10.5	< 10.5	
	alpha-Terpineol	98-55-5	10.5	< 10.5	
	Aniline	62-53-3	10.5	< 10.5	
	Anthracene	120-12-7	10.5	< 10.5	
	Aramite	140-57-8	10.5	< 10.5	
	Azobenzene	103-33-3	10.5	< 10.5	
	Benz(a)anthracene	56-55-3	10.5	< 10.5	
	Benzidine	92-87-5	10.5	< 10.5	
	Benzo(a)pyrene	50-32-8	10.5	< 10.5	

Report Date: 12/6/2010 Page 27 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001F

Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/2/2010 2232h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Benzo(b)fluoranthene	205-99-2	10.5	< 10.5	
01113	Benzo(g,h,i)perylene	191-24-2	10.5	< 10.5	
	Benzo(k)fluoranthene	207-08-9	10.5	< 10.5	
	Benzoic acid	65-85-0	21.1	< 21.1	
(801) 263-8686	Benzyl alcohol	100-51-6	10.5	< 10.5	
Toll Free (888) 263-8686	Bis(2-chloroethoxy)methane	111-91-1	10.5	< 10.5	
Fax (801)263-8687	Bis(2-chloroethyl) ether	111-44-4	10.5	< 10.5	
email: awal@awal-labs.com	Bis(2-chloroisopropyl) ether	108-60-1	10.5	< 10.5	
	Bis(2-ethylhexyl) phthalate	117-81-7	10.5	< 10.5	
Kyle F. Gross	bis(2-ethylhexyl)adipate	103-23-1	10.5	< 10.5	
·	Butyl benzyl phthalate	85-68-7	10.5	< 10.5	
Laboratory Director	Carbazole	86-74-8	10.5	< 10.5	
, p. 1	Chlorobenzilate	510-15-6	10.5	< 10.5	
Jose Rocha	Chrysene	218-01-9	10.5	< 10.5	
QA Officer	Di-n-butyl phthalate	84-74-2	10.5	< 10.5	
	Di-n-octyl phthalate	117-84-0	10.5	< 10.5	
	Diallate (cis or trans)	2303-16-4	10.5	< 10.5	
	Dibenz(a,h)anthracene	53-70-3	10.5	< 10.5	
	Dibenzofuran	132-64-9	10.5	< 10.5	
	Diethyl phthalate	84-66-2	10.5	< 10.5	
	Dimethoate	60-51-5	10.5	< 10.5	
	Dimethyl phthalate	131-11-3	10.5	< 10.5	
	Dimethylaminoazobenzene	60-11-7	10.5	< 10.5	
	Dinoseb	88-85-7	10.5	< 10.5	
	Diphenylamine	122-39-4	10.5	< 10.5	
	Disulfoton	298-04-4	10.5	< 10.5	
	Ethyl methanesulfonate	62-50-0	10.5	< 10.5	
	Famphur	52-85-7	10.5	< 10.5	
	Fluoranthene	206-44-0	10.5	< 10.5	

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**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001F

Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/2/2010 2232h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Fluorene	86-73-7	10.5	< 10.5	
0.110	Hexachlorobenzene	118-74-1	10.5	< 10.5	
	Hexachlorobutadiene	87-68-3	10.5	< 10.5	
	Hexachlorocyclopentadiene	77-47-4	10.5	< 10.5	
(801) 263-8686	Hexachloroethane	67-72-1	10.5	< 10.5	
Toll Free (888) 263-8686	Hexachlorophene	70-30-4	10.5	< 10.5	
Fax (801)263-8687	Hexachloropropene	1888-71-7	10.5	< 10.5	
email: awal@awal-labs.com	Indene	95-13-6	10.5	< 10.5	
	Indeno(1,2,3-cd)pyrene	193-39-5	10.5	< 10.5	
Kyle F. Gross	Isodrin	465-73-6	10.5	< 10.5	
•	Isophorone	78-59-1	10.5	< 10.5	
Laboratory Director	Isosafrole	120-58-1	10.5	< 10.5	
r D 1	Kepone	143-50-0	10.5	< 10.5	
Jose Rocha	Methapyrilene	91-80-5	10.5	< 10.5	
QA Officer	Methyl methanesulfonate	66-27-3	10.5	< 10.5	
	n-Decane	124-18-5	10.5	< 10.5	
	N-Nitrosodi-n-butylamine	924-16-3	10.5	< 10.5	
	N-Nitrosodiethylamine	55-18-5	10.5	< 10.5	
	N-Nitrosodimethylamine	62-75-9	10.5	< 10.5	
	N-Nitrosodiphenylamine	86-30-6	10.5	< 10.5	
	N-Nitrosodi-n-propylamine	621-64-7	10.5	< 10.5	
	N-Nitrosomethylethylamine	10595-95-6	10.5	< 10.5	
	N-Nitrosomorpholine	59-89-2	10.5	< 10.5	
	N-Nitrosopiperidine	100-75-4	10.5	< 10.5	
	N-Nitrosopyrrolidine	930-55-2	10.5	< 10.5	
	n-Octadecane	593-45-3	10.5	< 10.5	
	Naphthalene	91-20-3	10.5	< 10.5	
	Nitrobenzene	98-95-3	10.5	< 10.5	
	Nitroquinoline-1-oxide	56-57-5	10.5	< 10.5	

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**ANALYTICAL** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001F

Client Sample ID: 4992095 Red Butte AB Garden

 Collection Date:
 11/30/2010 0950h
 Analyzed: 12/2/2010 2232h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

LABORATORIES Analytical Results

SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	O,O,O-Triethyl phosphorothioate	126-68-1	10.5	< 10.5	
	o-Toluidine	95-53-4	10.5	< 10.5	
	Parathion	56-38-2	10.5	< 10.5	
	Methyl parathion	298-00-0	10.5	< 10.5	
(801) 263-8686	Pentachlorobenzene	608-93-5	10.5	< 10.5	
Toll Free (888) 263-8686	Pentachloronitrobenzene	82-68-8	10.5	< 10.5	
Fax (801)263-8687	Pentachlorophenol	87-86-5	10.5	< 10.5	
email: awal@awal-labs.com	Phenacetin	62-44-2	10.5	< 10.5	
	Phenanthrene	85-01-8	10.5	< 10.5	
Kyle F. Gross	Phenol	108-95-2	10.5	< 10.5	
•	Phorate	298-02-2	10.5	< 10.5	
Laboratory Director	Pronamide	23950-58-5	10.5	< 10.5	
I D 1	Pyrene	129-00-0	10.5	< 10.5	
Jose Rocha	Pyridine	110-86-1	10.5	< 10.5	
QA Officer	Quinoline	91-22-5	10.5	< 10.5	
	Safrole	94-59-7	10.5	< 10.5	
	Tetraethyl dithiopyrophosphate	3689-24-5	10.5	< 10.5	
	Thionazin	297-97-2	10.5	< 10.5	
	Surr: 2,4,6-Tribromophenol	118-79-6	10-159	60.4	
	Surr: 2-Fluorobiphenyl	321-60-8	10-124	57.0	
	Surr: 2-Fluorophenol	367-12-4	14-106	36.1	
	Surr: Nitrobenzene-d5	4165-60-0	10-180	46.9	
	Surr: Phenol-d6	13127-88-3	10-122	18.1	
	Surr: Terphenyl-d14	1718-51-0	10-199	106	

 ${\it The sample was analyzed for TICs and no unknown peaks were detected.}$ 



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002F

Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/2/2010 2300h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1´-Biphenyl	92-52-4	10.0	< 10.0	
	1,2,4,5-Tetrachlorobenzene	95-94-3	10.0	< 10.0	
	1,2,4-Trichlorobenzene	120-82-1	10.0	< 10.0	
	1,2-Dichlorobenzene	95-50-1	10.0	< 10.0	
(801) 263-8686	1,3,5-Trinitrobenzene	99-35-4	10.0	< 10.0	
Toll Free (888) 263-8686	1,4-Naphthoquinone	130-15-4	10.0	< 10.0	
Fax (801)263-8687	1,3-Dichlorobenzene	541-73-1	10.0	< 10.0	
email: awal@awal-labs.com	1,3-Dinitrobenzene	99-65-0	10.0	< 10.0	
	1,4-Dichlorobenzene	106-46-7	10.0	< 10.0	
Kyle F. Gross	1,4-Phenylenediamine	106-50-3	10.0	< 10.0	
Laboratory Director	1-Chloronaphthalene	90-13-1	10.0	< 10.0	
Laboratory Director	1-Methylnaphthalene	90-12-0	10.0	< 10.0	
I D 1	1-Naphthylamine	134-32-7	10.0	< 10.0	
Jose Rocha	2,3,4,6-Tetrachlorophenol	58-90-2	10.0	< 10.0	
QA Officer	2,4,5-Trichlorophenol	95-95-4	10.0	< 10.0	
	2,4,6-Trichlorophenol	88-06-2	10.0	< 10.0	
	2,4-Dichlorophenol	120-83-2	10.0	< 10.0	
	2,4-Dimethylphenol	105-67-9	10.0	< 10.0	
	2,4-Dinitrophenol	51-28-5	20.0	< 20.0	
	2,4-Dinitrotoluene	121-14-2	10.0	< 10.0	
	2,6-Dichlorophenol	87-65-0	10.0	< 10.0	
	2,6-Dinitrotoluene	606-20-2	10.0	< 10.0	
	2-Acetylaminofluorene	53-96-3	10.0	< 10.0	
	2-Chloronaphthalene	91-58-7	10.0	< 10.0	
	2-Chlorophenol	95-57-8	10.0	< 10.0	
	2-Methylnaphthalene	91-57-6	10.0	< 10.0	
	2-Methylphenol	95-48-7	10.0	< 10.0	
	2-Naphthylamine	91-59-8	10.0	< 10.0	
	2-Nitroaniline	88-74-4	10.0	< 10.0	

Report Date: 12/6/2010 Page 31 of 162



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**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002F

Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/2/2010 2300h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	2-Nitrophenol	88-75-5	10.0	< 10.0	
	2-Picoline	109-06-8	10.0	< 10.0	
	3&4-Methylphenol		10.0	< 10.0	
	3,3'-Dichlorobenzidine	91-94-1	10.0	< 10.0	
(801) 263-8686	3,3'-Dimethylbenzidine	119-93-7	10.0	< 10.0	
Toll Free (888) 263-8686	3-Methylcholanthrene	56-49-5	10.0	< 10.0	
Fax (801)263-8687	3-Nitroaniline	99-09-2	10.0	< 10.0	
email: awal@awal-labs.com	4,6-Dinitro-2-methylphenol	534-52-1	10.0	< 10.0	
	4-Aminobiphenyl	92-67-1	10.0	< 10.0	
Kyle F. Gross	4-Bromophenyl phenyl ether	101-55-3	10.0	< 10.0	
•	4-Chloro-3-methylphenol	59-50-7	10.0	< 10.0	
Laboratory Director	4-Chloroaniline	106-47-8	10.0	< 10.0	
r n i	4-Chlorophenyl phenyl ether	7005-72-3	10.0	< 10.0	
Jose Rocha	4-Nitroaniline	100-01-6	10.0	< 10.0	
QA Officer	4-Nitrophenol	100-02-7	10.0	< 10.0	
	5-Nitro-o-toluidine	99-55-8	10.0	< 10.0	
	7,12-Dimethylbenz(a)anthracene	57-97-6	10.0	< 10.0	
	a,a-Dimethylphenethylamine	122-09-8	10.0	< 10.0	
	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Acetophenone	98-86-2	10.0	< 10.0	
	alpha-Terpineol	98-55-5	10.0	< 10.0	
	Aniline	62-53-3	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Aramite	140-57-8	10.0	< 10.0	
	Azobenzene	103-33-3	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
	Benzidine	92-87-5	10.0	< 10.0	
	Benzo(a)pyrene	50-32-8	10.0	< 10.0	

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**LABORATORIES** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002F

AMERICAN Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/2/2010 2300h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
0.110	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Benzoic acid	65-85-0	20.0	< 20.0	
(801) 263-8686	Benzyl alcohol	100-51-6	10.0	< 10.0	
Toll Free (888) 263-8686	Bis(2-chloroethoxy)methane	111-91-1	10.0	< 10.0	
Fax (801)263-8687	Bis(2-chloroethyl) ether	111-44-4	10.0	< 10.0	
email: awal@awal-labs.com	Bis(2-chloroisopropyl) ether	108-60-1	10.0	< 10.0	
	Bis(2-ethylhexyl) phthalate	117-81-7	10.0	< 10.0	
Kyle F. Gross	bis(2-ethylhexyl)adipate	103-23-1	10.0	< 10.0	
•	Butyl benzyl phthalate	85-68-7	10.0	< 10.0	
Laboratory Director	Carbazole	86-74-8	10.0	< 10.0	
. D. 1	Chlorobenzilate	510-15-6	10.0	< 10.0	
Jose Rocha	Chrysene	218-01-9	10.0	< 10.0	
QA Officer	Di-n-butyl phthalate	84-74-2	10.0	< 10.0	
	Di-n-octyl phthalate	117-84-0	10.0	< 10.0	
	Diallate (cis or trans)	2303-16-4	10.0	< 10.0	
	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
	Dibenzofuran	132-64-9	10.0	< 10.0	
	Diethyl phthalate	84-66-2	10.0	< 10.0	
	Dimethoate	60-51-5	10.0	< 10.0	
	Dimethyl phthalate	131-11-3	10.0	< 10.0	
	Dimethylaminoazobenzene	60-11-7	10.0	< 10.0	
	Dinoseb	88-85-7	10.0	< 10.0	
	Diphenylamine	122-39-4	10.0	< 10.0	
	Disulfoton	298-04-4	10.0	< 10.0	
	Ethyl methanesulfonate	62-50-0	10.0	< 10.0	
	Famphur	52-85-7	10.0	< 10.0	
	Fluoranthene	206-44-0	10.0	< 10.0	

Report Date: 12/6/2010 Page 33 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002F

Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/2/2010 2300h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
Salt Lake City, UT	Elyana	86-73-7	10.0	< 10.0	
84115	Fluorene Hexachlorobenzene	118-74-1	10.0		
	Hexachlorobutadiene	87-68-3	10.0	< 10.0 < 10.0	
		77-47-4	10.0	< 10.0	
(801) 263-8686	Hexachlorocyclopentadiene Hexachloroethane	67-72-1	10.0	< 10.0	
Toll Free (888) 263-8686					
	Hexachlorophene	70-30-4	10.0	< 10.0	
Fax (801)263-8687	Hexachloropropene	1888-71-7	10.0	< 10.0	
email: awal@awal-labs.com	Indene	95-13-6	10.0	< 10.0	
	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Kyle F. Gross	Isodrin	465-73-6	10.0	< 10.0	
Laboratory Director	Isophorone	78-59-1	10.0	< 10.0	
	Isosafrole	120-58-1	10.0	< 10.0	
Jose Rocha	Kepone	143-50-0	10.0	< 10.0	
QA Officer	Methapyrilene	91-80-5	10.0	< 10.0	
Q11 Officer	Methyl methanesulfonate	66-27-3	10.0	< 10.0	
	n-Decane	124-18-5	10.0	< 10.0	
	N-Nitrosodi-n-butylamine	924-16-3	10.0	< 10.0	
	N-Nitrosodiethylamine	55-18-5	10.0	< 10.0	
	N-Nitrosodimethylamine	62-75-9	10.0	< 10.0	
	N-Nitrosodiphenylamine	86-30-6	10.0	< 10.0	
	N-Nitrosodi-n-propylamine	621-64-7	10.0	< 10.0	
	N-Nitrosomethylethylamine	10595-95-6	10.0	< 10.0	
	N-Nitrosomorpholine	59-89-2	10.0	< 10.0	
	N-Nitrosopiperidine	100-75-4	10.0	< 10.0	
	N-Nitrosopyrrolidine	930-55-2	10.0	< 10.0	
	n-Octadecane	593-45-3	10.0	< 10.0	
	Naphthalene	91-20-3	10.0	< 10.0	
	Nitrobenzene	98-95-3	10.0	< 10.0	
	Nitroquinoline-1-oxide	56-57-5	10.0	< 10.0	

Report Date: 12/6/2010 Page 34 of 162



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**ANALYTICAL** 

**LABORATORIES** 

### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002F

AMERICAN Client Sample ID: 4992085 Red Butte BL Garden

 Collection Date:
 11/30/2010 1030h
 Analyzed: 12/2/2010 2300h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	O,O,O-Triethyl phosphorothioate	126-68-1	10.0	< 10.0	
	o-Toluidine	95-53-4	10.0	< 10.0	
	Parathion	56-38-2	10.0	< 10.0	
	Methyl parathion	298-00-0	10.0	< 10.0	
(801) 263-8686	Pentachlorobenzene	608-93-5	10.0	< 10.0	
Toll Free (888) 263-8686	Pentachloronitrobenzene	82-68-8	10.0	< 10.0	
Fax (801)263-8687	Pentachlorophenol	87-86-5	10.0	< 10.0	
email: awal@awal-labs.com	Phenacetin	62-44-2	10.0	< 10.0	
	Phenanthrene	85-01-8	10.0	< 10.0	
Kyle F. Gross	Phenol	108-95-2	10.0	< 10.0	
Laboratory Director	Phorate	298-02-2	10.0	< 10.0	
Laboratory Director	Pronamide	23950-58-5	10.0	< 10.0	
r D 1	Pyrene	129-00-0	10.0	< 10.0	
Jose Rocha	Pyridine	110-86-1	10.0	< 10.0	
QA Officer	Quinoline	91-22-5	10.0	< 10.0	
	Safrole	94-59-7	10.0	< 10.0	
	Tetraethyl dithiopyrophosphate	3689-24-5	10.0	< 10.0	
	Thionazin	297-97-2	10.0	< 10.0	
	Surr: 2,4,6-Tribromophenol	118-79-6	10-159	55.4	
	Surr: 2-Fluorobiphenyl	321-60-8	10-124	33.4	
	Surr: 2-Fluorophenol	367-12-4	14-106	29.8	
	Surr: Nitrobenzene-d5	4165-60-0	10-180	20.7	
	Surr: Phenol-d6	13127-88-3	10-122	25.5	
	Surr: Terphenyl-d14	1718-51-0	10-199	108	

The sample was analyzed for TICs and no unknown peaks were detected.



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003F

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

 Collection Date:
 11/30/2010 1215h
 Analyzed: 12/2/2010 2327h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1´-Biphenyl	92-52-4	10.0	< 10.0	_
	1,2,4,5-Tetrachlorobenzene	95-94-3	10.0	< 10.0	
	1,2,4-Trichlorobenzene	120-82-1	10.0	< 10.0	
	1,2-Dichlorobenzene	95-50-1	10.0	< 10.0	
(801) 263-8686	1,3,5-Trinitrobenzene	99-35-4	10.0	< 10.0	
Toll Free (888) 263-8686	1,4-Naphthoquinone	130-15-4	10.0	< 10.0	
Fax (801)263-8687	1,3-Dichlorobenzene	541-73-1	10.0	< 10.0	
email: awal@awal-labs.com	1,3-Dinitrobenzene	99-65-0	10.0	< 10.0	
	1,4-Dichlorobenzene	106-46-7	10.0	< 10.0	
Kyle F. Gross	1,4-Phenylenediamine	106-50-3	10.0	< 10.0	
Laboratory Director	1-Chloronaphthalene	90-13-1	10.0	< 10.0	
Laboratory Director	1-Methylnaphthalene	90-12-0	10.0	< 10.0	
I D 1	1-Naphthylamine	134-32-7	10.0	< 10.0	
Jose Rocha	2,3,4,6-Tetrachlorophenol	58-90-2	10.0	< 10.0	
QA Officer	2,4,5-Trichlorophenol	95-95-4	10.0	< 10.0	
	2,4,6-Trichlorophenol	88-06-2	10.0	< 10.0	
	2,4-Dichlorophenol	120-83-2	10.0	< 10.0	
	2,4-Dimethylphenol	105-67-9	10.0	< 10.0	
	2,4-Dinitrophenol	51-28-5	20.0	< 20.0	
	2,4-Dinitrotoluene	121-14-2	10.0	< 10.0	
	2,6-Dichlorophenol	87-65-0	10.0	< 10.0	
	2,6-Dinitrotoluene	606-20-2	10.0	< 10.0	
	2-Acetylaminofluorene	53-96-3	10.0	< 10.0	
	2-Chloronaphthalene	91-58-7	10.0	< 10.0	
	2-Chlorophenol	95-57-8	10.0	< 10.0	
	2-Methylnaphthalene	91-57-6	10.0	< 10.0	
	2-Methylphenol	95-48-7	10.0	< 10.0	
	2-Naphthylamine	91-59-8	10.0	< 10.0	
	2-Nitroaniline	88-74-4	10.0	< 10.0	

Report Date: 12/6/2010 Page 36 of 162



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**LABORATORIES** 

**WEST** 

#### ORGANIC ANALYTICAL REPORT

**Client:** Utah DEQ DERR **Contact:** Jim Harris

**Project:** Red Butte 1011529-003F Lab Sample ID:

Client Sample ID: 4992084 Red Butte @ 1500 E.

**Collection Date:** 11/30/2010 1215h Analyzed: 12/2/2010 2327h **Received Date:** 11/30/2010 1538h Extracted: 12/2/2010 0923h

**Method Used: EPA625** 

SVOA by GC/MS Method 625/3510 **Analytical Results** 

Units: µg/L **Dilution Factor: 1 CAS** Reporting 463 West 3600 South **Analytical** Number Compound Limit Result Qual Salt Lake City, UT 88-75-5 10.0 < 10.0 2-Nitrophenol 84115 109-06-8 10.0 2-Picoline < 10.0 3&4-Methylphenol 10.0 < 10.0 91-94-1 3,3'-Dichlorobenzidine 10.0 < 10.0 (801) 263-8686 3,3'-Dimethylbenzidine 119-93-7 10.0 < 10.0 Toll Free (888) 263-8686 3-Methylcholanthrene 56-49-5 10.0 < 10.0 3-Nitroaniline 99-09-2 10.0 < 10.0 Fax (801)263-8687 4,6-Dinitro-2-methylphenol 534-52-1 10.0 < 10.0 email: awal@awal-labs.com 4-Aminobiphenyl 92-67-1 10.0 < 10.0 4-Bromophenyl phenyl ether 101-55-3 10.0 < 10.0 Kyle F. Gross 4-Chloro-3-methylphenol 59-50-7 10.0 < 10.0 Laboratory Director 4-Chloroaniline 106-47-8 10.0 < 10.0 4-Chlorophenyl phenyl ether 7005-72-3 10.0 < 10.0 Jose Rocha 4-Nitroaniline 100-01-6 10.0 < 10.0 **QA** Officer 4-Nitrophenol 10.0 < 10.0 100-02-7 5-Nitro-o-toluidine 99-55-8 10.0 < 10.0 57-97-6 10.0 7,12-Dimethylbenz(a)anthracene < 10.0 a,a-Dimethylphenethylamine 122-09-8 10.0 < 10.0 < 10.0 Acenaphthene 83-32-9 10.0 208-96-8 10.0 < 10.0 Acenaphthylene 98-86-2 Acetophenone 10.0 < 10.0 98-55-5 alpha-Terpineol 10.0 < 10.0 Aniline 62-53-3 10.0 < 10.0 Anthracene 120-12-7 10.0 < 10.0 10.0 Aramite 140-57-8 < 10.0 Azobenzene 103-33-3 10.0 < 10.0 10.0 Benz(a)anthracene 56-55-3 < 10.0 Benzidine 92-87-5 10.0 < 10.0 50-32-8 10.0 < 10.0 Benzo(a)pyrene

Report Date: 12/6/2010 Page 37 of 162



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**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003F

AMERICAN Client Sample ID: 4992084 Red Butte @ 1500 E.

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 11/30/2010 1215h
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Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Benzoic acid	65-85-0	20.0	< 20.0	
(801) 263-8686	Benzyl alcohol	100-51-6	10.0	< 10.0	
Toll Free (888) 263-8686	Bis(2-chloroethoxy)methane	111-91-1	10.0	< 10.0	
Fax (801)263-8687	Bis(2-chloroethyl) ether	111-44-4	10.0	< 10.0	
email: awal@awal-labs.com	Bis(2-chloroisopropyl) ether	108-60-1	10.0	< 10.0	
	Bis(2-ethylhexyl) phthalate	117-81-7	10.0	< 10.0	
Kyle F. Gross	bis(2-ethylhexyl)adipate	103-23-1	10.0	< 10.0	
ř	Butyl benzyl phthalate	85-68-7	10.0	< 10.0	
Laboratory Director	Carbazole	86-74-8	10.0	< 10.0	
T D 1	Chlorobenzilate	510-15-6	10.0	< 10.0	
Jose Rocha	Chrysene	218-01-9	10.0	< 10.0	
QA Officer	Di-n-butyl phthalate	84-74-2	10.0	< 10.0	
	Di-n-octyl phthalate	117-84-0	10.0	< 10.0	
	Diallate (cis or trans)	2303-16-4	10.0	< 10.0	
	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
	Dibenzofuran	132-64-9	10.0	< 10.0	
	Diethyl phthalate	84-66-2	10.0	< 10.0	
	Dimethoate	60-51-5	10.0	< 10.0	
	Dimethyl phthalate	131-11-3	10.0	< 10.0	
	Dimethylaminoazobenzene	60-11-7	10.0	< 10.0	
	Dinoseb	88-85-7	10.0	< 10.0	
	Diphenylamine	122-39-4	10.0	< 10.0	
	Disulfoton	298-04-4	10.0	< 10.0	
	Ethyl methanesulfonate	62-50-0	10.0	< 10.0	
	Famphur	52-85-7	10.0	< 10.0	
	Fluoranthene	206-44-0	10.0	< 10.0	

Report Date: 12/6/2010 Page 38 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003F

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

 Collection Date:
 11/30/2010 1215h
 Analyzed: 12/2/2010 2327h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Fluorene	86-73-7	10.0	< 10.0	
	Hexachlorobenzene	118-74-1	10.0	< 10.0	
	Hexachlorobutadiene	87-68-3	10.0	< 10.0	
	Hexachlorocyclopentadiene	77-47-4	10.0	< 10.0	
(801) 263-8686	Hexachloroethane	67-72-1	10.0	< 10.0	
Toll Free (888) 263-8686	Hexachlorophene	70-30-4	10.0	< 10.0	
Fax (801)263-8687	Hexachloropropene	1888-71-7	10.0	< 10.0	
email: awal@awal-labs.com	Indene	95-13-6	10.0	< 10.0	
	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Kyle F. Gross	Isodrin	465-73-6	10.0	< 10.0	
•	Isophorone	78-59-1	10.0	< 10.0	
Laboratory Director	Isosafrole	120-58-1	10.0	< 10.0	
I D 1	Kepone	143-50-0	10.0	< 10.0	
Jose Rocha	Methapyrilene	91-80-5	10.0	< 10.0	
QA Officer	Methyl methanesulfonate	66-27-3	10.0	< 10.0	
	n-Decane	124-18-5	10.0	< 10.0	
	N-Nitrosodi-n-butylamine	924-16-3	10.0	< 10.0	
	N-Nitrosodiethylamine	55-18-5	10.0	< 10.0	
	N-Nitrosodimethylamine	62-75-9	10.0	< 10.0	
	N-Nitrosodiphenylamine	86-30-6	10.0	< 10.0	
	N-Nitrosodi-n-propylamine	621-64-7	10.0	< 10.0	
	N-Nitrosomethylethylamine	10595-95-6	10.0	< 10.0	
	N-Nitrosomorpholine	59-89-2	10.0	< 10.0	
	N-Nitrosopiperidine	100-75-4	10.0	< 10.0	
	N-Nitrosopyrrolidine	930-55-2	10.0	< 10.0	
	n-Octadecane	593-45-3	10.0	< 10.0	
	Naphthalene	91-20-3	10.0	< 10.0	
	Nitrobenzene	98-95-3	10.0	< 10.0	
	Nitroquinoline-1-oxide	56-57-5	10.0	< 10.0	

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**ANALYTICAL** 

**LABORATORIES** 

# ORGANIC ANALYTICAL REPORT

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003F

AMERICAN Client Sample ID: 4992084 Red Butte @ 1500 E.

 Collection Date:
 11/30/2010 1215h
 Analyzed: 12/2/2010 2327h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	O,O,O-Triethyl phosphorothioate	126-68-1	10.0	< 10.0	
	o-Toluidine	95-53-4	10.0	< 10.0	
	Parathion	56-38-2	10.0	< 10.0	
	Methyl parathion	298-00-0	10.0	< 10.0	
(801) 263-8686	Pentachlorobenzene	608-93-5	10.0	< 10.0	
Toll Free (888) 263-8686	Pentachloronitrobenzene	82-68-8	10.0	< 10.0	
Fax (801)263-8687	Pentachlorophenol	87-86-5	10.0	< 10.0	
email: awal@awal-labs.com	Phenacetin	62-44-2	10.0	< 10.0	
	Phenanthrene	85-01-8	10.0	< 10.0	
Kyle F. Gross	Phenol	108-95-2	10.0	< 10.0	
Laboratory Director	Phorate	298-02-2	10.0	< 10.0	
Laboratory Director	Pronamide	23950-58-5	10.0	< 10.0	
Jose Rocha	Pyrene	129-00-0	10.0	< 10.0	
	Pyridine	110-86-1	10.0	< 10.0	
QA Officer	Quinoline	91-22-5	10.0	< 10.0	
	Safrole	94-59-7	10.0	< 10.0	
	Tetraethyl dithiopyrophosphate	3689-24-5	10.0	< 10.0	
	Thionazin	297-97-2	10.0	< 10.0	
	Surr: 2,4,6-Tribromophenol	118-79-6	10-159	62.0	
	Surr: 2-Fluorobiphenyl	321-60-8	10-124	60.7	
	Surr: 2-Fluorophenol	367-12-4	14-106	41.3	
	Surr: Nitrobenzene-d5	4165-60-0	10-180	55.4	
	Surr: Phenol-d6	13127-88-3	10-122	19.8	
	Surr: Terphenyl-d14	1718-51-0	10-199	110	

The sample was analyzed for TICs and no unknown peaks were detected.



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004F

Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/2/2010 2355h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1´-Biphenyl	92-52-4	10.0	< 10.0	_
	1,2,4,5-Tetrachlorobenzene	95-94-3	10.0	< 10.0	
	1,2,4-Trichlorobenzene	120-82-1	10.0	< 10.0	
	1,2-Dichlorobenzene	95-50-1	10.0	< 10.0	
(801) 263-8686	1,3,5-Trinitrobenzene	99-35-4	10.0	< 10.0	
Toll Free (888) 263-8686	1,4-Naphthoquinone	130-15-4	10.0	< 10.0	
Fax (801)263-8687	1,3-Dichlorobenzene	541-73-1	10.0	< 10.0	
email: awal@awal-labs.com	1,3-Dinitrobenzene	99-65-0	10.0	< 10.0	
	1,4-Dichlorobenzene	106-46-7	10.0	< 10.0	
Kyle F. Gross	1,4-Phenylenediamine	106-50-3	10.0	< 10.0	
•	1-Chloronaphthalene	90-13-1	10.0	< 10.0	
Laboratory Director	1-Methylnaphthalene	90-12-0	10.0	< 10.0	
I D 1	1-Naphthylamine	134-32-7	10.0	< 10.0	
Jose Rocha	2,3,4,6-Tetrachlorophenol	58-90-2	10.0	< 10.0	
QA Officer	2,4,5-Trichlorophenol	95-95-4	10.0	< 10.0	
	2,4,6-Trichlorophenol	88-06-2	10.0	< 10.0	
	2,4-Dichlorophenol	120-83-2	10.0	< 10.0	
	2,4-Dimethylphenol	105-67-9	10.0	< 10.0	
	2,4-Dinitrophenol	51-28-5	20.0	< 20.0	
	2,4-Dinitrotoluene	121-14-2	10.0	< 10.0	
	2,6-Dichlorophenol	87-65-0	10.0	< 10.0	
	2,6-Dinitrotoluene	606-20-2	10.0	< 10.0	
	2-Acetylaminofluorene	53-96-3	10.0	< 10.0	
	2-Chloronaphthalene	91-58-7	10.0	< 10.0	
	2-Chlorophenol	95-57-8	10.0	< 10.0	
	2-Methylnaphthalene	91-57-6	10.0	< 10.0	
	2-Methylphenol	95-48-7	10.0	< 10.0	
	2-Naphthylamine	91-59-8	10.0	< 10.0	
	2-Nitroaniline	88-74-4	10.0	< 10.0	

Report Date: 12/6/2010 Page 41 of 162



**ANALYTICAL** 

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004F

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/2/2010 2355h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	2-Nitrophenol	88-75-5	10.0	< 10.0	
	2-Picoline	109-06-8	10.0	< 10.0	
	3&4-Methylphenol		10.0	< 10.0	
	3,3'-Dichlorobenzidine	91-94-1	10.0	< 10.0	
(801) 263-8686	3,3'-Dimethylbenzidine	119-93-7	10.0	< 10.0	
Toll Free (888) 263-8686	3-Methylcholanthrene	56-49-5	10.0	< 10.0	
Fax (801)263-8687	3-Nitroaniline	99-09-2	10.0	< 10.0	
email: awal@awal-labs.com	4,6-Dinitro-2-methylphenol	534-52-1	10.0	< 10.0	
	4-Aminobiphenyl	92-67-1	10.0	< 10.0	
Kyle F. Gross	4-Bromophenyl phenyl ether	101-55-3	10.0	< 10.0	
•	4-Chloro-3-methylphenol	59-50-7	10.0	< 10.0	
Laboratory Director	4-Chloroaniline	106-47-8	10.0	< 10.0	
	4-Chlorophenyl phenyl ether	7005-72-3	10.0	< 10.0	
Jose Rocha	4-Nitroaniline	100-01-6	10.0	< 10.0	
QA Officer	4-Nitrophenol	100-02-7	10.0	< 10.0	
	5-Nitro-o-toluidine	99-55-8	10.0	< 10.0	
	7,12-Dimethylbenz(a)anthracene	57-97-6	10.0	< 10.0	
	a,a-Dimethylphenethylamine	122-09-8	10.0	< 10.0	
	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Acetophenone	98-86-2	10.0	< 10.0	
	alpha-Terpineol	98-55-5	10.0	< 10.0	
	Aniline	62-53-3	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Aramite	140-57-8	10.0	< 10.0	
	Azobenzene	103-33-3	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
	Benzidine	92-87-5	10.0	< 10.0	
	Benzo(a)pyrene	50-32-8	10.0	< 10.0	

Report Date: 12/6/2010 Page 42 of 162



**ANALYTICAL** 

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004F

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/2/2010 2355h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	_
	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Benzoic acid	65-85-0	20.0	< 20.0	
(801) 263-8686	Benzyl alcohol	100-51-6	10.0	< 10.0	
Toll Free (888) 263-8686	Bis(2-chloroethoxy)methane	111-91-1	10.0	< 10.0	
Fax (801)263-8687	Bis(2-chloroethyl) ether	111-44-4	10.0	< 10.0	
email: awal@awal-labs.com	Bis(2-chloroisopropyl) ether	108-60-1	10.0	< 10.0	
	Bis(2-ethylhexyl) phthalate	117-81-7	10.0	< 10.0	
Kyle F. Gross	bis(2-ethylhexyl)adipate	103-23-1	10.0	< 10.0	
•	Butyl benzyl phthalate	85-68-7	10.0	< 10.0	
Laboratory Director	Carbazole	86-74-8	10.0	< 10.0	
I D 1	Chlorobenzilate	510-15-6	10.0	< 10.0	
Jose Rocha	Chrysene	218-01-9	10.0	< 10.0	
QA Officer	Di-n-butyl phthalate	84-74-2	10.0	< 10.0	
	Di-n-octyl phthalate	117-84-0	10.0	< 10.0	
	Diallate (cis or trans)	2303-16-4	10.0	< 10.0	
	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
	Dibenzofuran	132-64-9	10.0	< 10.0	
	Diethyl phthalate	84-66-2	10.0	< 10.0	
	Dimethoate	60-51-5	10.0	< 10.0	
	Dimethyl phthalate	131-11-3	10.0	< 10.0	
	Dimethylaminoazobenzene	60-11-7	10.0	< 10.0	
	Dinoseb	88-85-7	10.0	< 10.0	
	Diphenylamine	122-39-4	10.0	< 10.0	
	Disulfoton	298-04-4	10.0	< 10.0	
	Ethyl methanesulfonate	62-50-0	10.0	< 10.0	
	Famphur	52-85-7	10.0	< 10.0	
	Fluoranthene	206-44-0	10.0	< 10.0	

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**ANALYTICAL** 

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004F

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/2/2010 2355h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Fluorene	86-73-7	10.0	< 10.0	
04113	Hexachlorobenzene	118-74-1	10.0	< 10.0	
	Hexachlorobutadiene	87-68-3	10.0	< 10.0	
	Hexachlorocyclopentadiene	77-47-4	10.0	< 10.0	
(801) 263-8686	Hexachloroethane	67-72-1	10.0	< 10.0	
Toll Free (888) 263-8686	Hexachlorophene	70-30-4	10.0	< 10.0	
Fax (801)263-8687	Hexachloropropene	1888-71-7	10.0	< 10.0	
email: awal@awal-labs.com	Indene	95-13-6	10.0	< 10.0	
	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Kyle F. Gross	Isodrin	465-73-6	10.0	< 10.0	
·	Isophorone	78-59-1	10.0	< 10.0	
Laboratory Director	Isosafrole	120-58-1	10.0	< 10.0	
	Kepone	143-50-0	10.0	< 10.0	
Jose Rocha	Methapyrilene	91-80-5	10.0	< 10.0	
QA Officer	Methyl methanesulfonate	66-27-3	10.0	< 10.0	
	n-Decane	124-18-5	10.0	< 10.0	
	N-Nitrosodi-n-butylamine	924-16-3	10.0	< 10.0	
	N-Nitrosodiethylamine	55-18-5	10.0	< 10.0	
	N-Nitrosodimethylamine	62-75-9	10.0	< 10.0	
	N-Nitrosodiphenylamine	86-30-6	10.0	< 10.0	
	N-Nitrosodi-n-propylamine	621-64-7	10.0	< 10.0	
	N-Nitrosomethylethylamine	10595-95-6	10.0	< 10.0	
	N-Nitrosomorpholine	59-89-2	10.0	< 10.0	
	N-Nitrosopiperidine	100-75-4	10.0	< 10.0	
	N-Nitrosopyrrolidine	930-55-2	10.0	< 10.0	
	n-Octadecane	593-45-3	10.0	< 10.0	
	Naphthalene	91-20-3	10.0	< 10.0	
	Nitrobenzene	98-95-3	10.0	< 10.0	
	Nitroquinoline-1-oxide	56-57-5	10.0	< 10.0	

Report Date: 12/6/2010 Page 44 of 162



**ANALYTICAL** 

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004F

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

 Collection Date:
 11/30/2010 1405h
 Analyzed: 12/2/2010 2355h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual		
84115	O,O,O-Triethyl phosphorothioate	126-68-1	10.0	< 10.0			
	o-Toluidine	95-53-4	10.0	< 10.0			
	Parathion	56-38-2	10.0	< 10.0			
	Methyl parathion	298-00-0	10.0	< 10.0			
(801) 263-8686	Pentachlorobenzene	608-93-5	10.0	< 10.0			
Toll Free (888) 263-8686	Pentachloronitrobenzene	82-68-8	10.0	< 10.0			
Fax (801)263-8687	Pentachlorophenol	87-86-5	10.0	< 10.0			
email: awal@awal-labs.com	Phenacetin	62-44-2	10.0	< 10.0			
	Phenanthrene	85-01-8	10.0	< 10.0			
Kyle F. Gross	Phenol	108-95-2	10.0	< 10.0			
•	Phorate	298-02-2	10.0	< 10.0			
Laboratory Director	Pronamide	23950-58-5	10.0	< 10.0			
. D. 1	Pyrene	129-00-0	10.0	< 10.0			
Jose Rocha	Pyridine	110-86-1	10.0	< 10.0			
QA Officer	Quinoline	91-22-5	10.0	< 10.0			
	Safrole	94-59-7	10.0	< 10.0			
	Tetraethyl dithiopyrophosphate	3689-24-5	10.0	< 10.0			
	Thionazin	297-97-2	10.0	< 10.0			
	Surr: 2,4,6-Tribromophenol	118-79-6	10-159	61.3			
	Surr: 2-Fluorobiphenyl	321-60-8	10-124	46.0			
	Surr: 2-Fluorophenol	367-12-4	14-106	28.1			
	Surr: Nitrobenzene-d5	4165-60-0	10-180	37.0			
	Surr: Phenol-d6	13127-88-3	10-122	28.7			
	Surr: Terphenyl-d14	1718-51-0	10-199	107			

The sample was analyzed for TICs and no unknown peaks were detected.



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005F

**Client Sample ID:** 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/3/2010 0022h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1'-Biphenyl	92-52-4	10.0	< 10.0	
01113	1,2,4,5-Tetrachlorobenzene	95-94-3	10.0	< 10.0	
	1,2,4-Trichlorobenzene	120-82-1	10.0	< 10.0	
	1,2-Dichlorobenzene	95-50-1	10.0	< 10.0	
(801) 263-8686	1,3,5-Trinitrobenzene	99-35-4	10.0	< 10.0	
Toll Free (888) 263-8686	1,4-Naphthoquinone	130-15-4	10.0	< 10.0	
Fax (801)263-8687	1,3-Dichlorobenzene	541-73-1	10.0	< 10.0	
email: awal@awal-labs.com	1,3-Dinitrobenzene	99-65-0	10.0	< 10.0	
	1,4-Dichlorobenzene	106-46-7	10.0	< 10.0	
Kyle F. Gross	1,4-Phenylenediamine	106-50-3	10.0	< 10.0	
•	1-Chloronaphthalene	90-13-1	10.0	< 10.0	
Laboratory Director	1-Methylnaphthalene	90-12-0	10.0	< 10.0	
	1-Naphthylamine	134-32-7	10.0	< 10.0	
Jose Rocha	2,3,4,6-Tetrachlorophenol	58-90-2	10.0	< 10.0	
QA Officer	2,4,5-Trichlorophenol	95-95-4	10.0	< 10.0	
	2,4,6-Trichlorophenol	88-06-2	10.0	< 10.0	
	2,4-Dichlorophenol	120-83-2	10.0	< 10.0	
	2,4-Dimethylphenol	105-67-9	10.0	< 10.0	
	2,4-Dinitrophenol	51-28-5	20.0	< 20.0	
	2,4-Dinitrotoluene	121-14-2	10.0	< 10.0	
	2,6-Dichlorophenol	87-65-0	10.0	< 10.0	
	2,6-Dinitrotoluene	606-20-2	10.0	< 10.0	
	2-Acetylaminofluorene	53-96-3	10.0	< 10.0	
	2-Chloronaphthalene	91-58-7	10.0	< 10.0	
	2-Chlorophenol	95-57-8	10.0	< 10.0	
	2-Methylnaphthalene	91-57-6	10.0	< 10.0	
	2-Methylphenol	95-48-7	10.0	< 10.0	
	2-Naphthylamine	91-59-8	10.0	< 10.0	
	2-Nitroaniline	88-74-4	10.0	< 10.0	

Report Date: 12/6/2010 Page 46 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005F

**Client Sample ID:** 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/3/2010 0022h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	2-Nitrophenol	88-75-5	10.0	< 10.0	
	2-Picoline	109-06-8	10.0	< 10.0	
	3&4-Methylphenol		10.0	< 10.0	
	3,3'-Dichlorobenzidine	91-94-1	10.0	< 10.0	
(801) 263-8686	3,3'-Dimethylbenzidine	119-93-7	10.0	< 10.0	
Toll Free (888) 263-8686	3-Methylcholanthrene	56-49-5	10.0	< 10.0	
Fax (801)263-8687	3-Nitroaniline	99-09-2	10.0	< 10.0	
email: awal@awal-labs.com	4,6-Dinitro-2-methylphenol	534-52-1	10.0	< 10.0	
	4-Aminobiphenyl	92-67-1	10.0	< 10.0	
Kyle F. Gross	4-Bromophenyl phenyl ether	101-55-3	10.0	< 10.0	
•	4-Chloro-3-methylphenol	59-50-7	10.0	< 10.0	
Laboratory Director	4-Chloroaniline	106-47-8	10.0	< 10.0	
T D 1	4-Chlorophenyl phenyl ether	7005-72-3	10.0	< 10.0	
Jose Rocha	4-Nitroaniline	100-01-6	10.0	< 10.0	
QA Officer	4-Nitrophenol	100-02-7	10.0	< 10.0	
	5-Nitro-o-toluidine	99-55-8	10.0	< 10.0	
	7,12-Dimethylbenz(a)anthracene	57-97-6	10.0	< 10.0	
	a,a-Dimethylphenethylamine	122-09-8	10.0	< 10.0	
	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Acetophenone	98-86-2	10.0	< 10.0	
	alpha-Terpineol	98-55-5	10.0	< 10.0	
	Aniline	62-53-3	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Aramite	140-57-8	10.0	< 10.0	
	Azobenzene	103-33-3	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
	Benzidine	92-87-5	10.0	< 10.0	
	Benzo(a)pyrene	50-32-8	10.0	< 10.0	

Report Date: 12/6/2010 Page 47 of 162



**ANALYTICAL** 

**LABORATORIES** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005F

AMERICAN Client Sample ID: 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/3/2010 0022h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Benzoic acid	65-85-0	20.0	< 20.0	
(801) 263-8686	Benzyl alcohol	100-51-6	10.0	< 10.0	
Toll Free (888) 263-8686	Bis(2-chloroethoxy)methane	111-91-1	10.0	< 10.0	
Fax (801)263-8687	Bis(2-chloroethyl) ether	111-44-4	10.0	< 10.0	
email: awal@awal-labs.com	Bis(2-chloroisopropyl) ether	108-60-1	10.0	< 10.0	
	Bis(2-ethylhexyl) phthalate	117-81-7	10.0	< 10.0	
Kyle F. Gross	bis(2-ethylhexyl)adipate	103-23-1	10.0	< 10.0	
ř	Butyl benzyl phthalate	85-68-7	10.0	< 10.0	
Laboratory Director	Carbazole	86-74-8	10.0	< 10.0	
T D 1	Chlorobenzilate	510-15-6	10.0	< 10.0	
Jose Rocha	Chrysene	218-01-9	10.0	< 10.0	
QA Officer	Di-n-butyl phthalate	84-74-2	10.0	< 10.0	
	Di-n-octyl phthalate	117-84-0	10.0	< 10.0	
	Diallate (cis or trans)	2303-16-4	10.0	< 10.0	
	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
	Dibenzofuran	132-64-9	10.0	< 10.0	
	Diethyl phthalate	84-66-2	10.0	< 10.0	
	Dimethoate	60-51-5	10.0	< 10.0	
	Dimethyl phthalate	131-11-3	10.0	< 10.0	
	Dimethylaminoazobenzene	60-11-7	10.0	< 10.0	
	Dinoseb	88-85-7	10.0	< 10.0	
	Diphenylamine	122-39-4	10.0	< 10.0	
	Disulfoton	298-04-4	10.0	< 10.0	
	Ethyl methanesulfonate	62-50-0	10.0	< 10.0	
	Famphur	52-85-7	10.0	< 10.0	
	Fluoranthene	206-44-0	10.0	< 10.0	

Report Date: 12/6/2010 Page 48 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005F

**Client Sample ID:** 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/3/2010 0022h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Fluorene	86-73-7	10.0	< 10.0	
01113	Hexachlorobenzene	118-74-1	10.0	< 10.0	
	Hexachlorobutadiene	87-68-3	10.0	< 10.0	
	Hexachlorocyclopentadiene	77-47-4	10.0	< 10.0	
(801) 263-8686	Hexachloroethane	67-72-1	10.0	< 10.0	
Toll Free (888) 263-8686	Hexachlorophene	70-30-4	10.0	< 10.0	
Fax (801)263-8687	Hexachloropropene	1888-71-7	10.0	< 10.0	
email: awal@awal-labs.com	Indene	95-13-6	10.0	< 10.0	
	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Kyle F. Gross	Isodrin	465-73-6	10.0	< 10.0	
Laboratory Director	Isophorone	78-59-1	10.0	< 10.0	
Laboratory Director	Isosafrole	120-58-1	10.0	< 10.0	
r D 1	Kepone	143-50-0	10.0	< 10.0	
Jose Rocha	Methapyrilene	91-80-5	10.0	< 10.0	
QA Officer	Methyl methanesulfonate	66-27-3	10.0	< 10.0	
	n-Decane	124-18-5	10.0	< 10.0	
	N-Nitrosodi-n-butylamine	924-16-3	10.0	< 10.0	
	N-Nitrosodiethylamine	55-18-5	10.0	< 10.0	
	N-Nitrosodimethylamine	62-75-9	10.0	< 10.0	
	N-Nitrosodiphenylamine	86-30-6	10.0	< 10.0	
	N-Nitrosodi-n-propylamine	621-64-7	10.0	< 10.0	
	N-Nitrosomethylethylamine	10595-95-6	10.0	< 10.0	
	N-Nitrosomorpholine	59-89-2	10.0	< 10.0	
	N-Nitrosopiperidine	100-75-4	10.0	< 10.0	
	N-Nitrosopyrrolidine	930-55-2	10.0	< 10.0	
	n-Octadecane	593-45-3	10.0	< 10.0	
	Naphthalene	91-20-3	10.0	< 10.0	
	Nitrobenzene	98-95-3	10.0	< 10.0	
	Nitroquinoline-1-oxide	56-57-5	10.0	< 10.0	

Report Date: 12/6/2010 Page 49 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005F

**Client Sample ID:** 4992290 J.R. @ 1700 S.

 Collection Date:
 11/30/2010 1430h
 Analyzed: 12/3/2010 0022h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	O,O,O-Triethyl phosphorothioate	126-68-1	10.0	< 10.0	
	o-Toluidine	95-53-4	10.0	< 10.0	
	Parathion	56-38-2	10.0	< 10.0	
	Methyl parathion	298-00-0	10.0	< 10.0	
(801) 263-8686	Pentachlorobenzene	608-93-5	10.0	< 10.0	
Toll Free (888) 263-8686	Pentachloronitrobenzene	82-68-8	10.0	< 10.0	
Fax (801)263-8687	Pentachlorophenol	87-86-5	10.0	< 10.0	
email: awal@awal-labs.com	Phenacetin	62-44-2	10.0	< 10.0	
	Phenanthrene	85-01-8	10.0	< 10.0	
Kyle F. Gross	Phenol	108-95-2	10.0	< 10.0	
Laboratory Director	Phorate	298-02-2	10.0	< 10.0	
Laboratory Director	Pronamide	23950-58-5	10.0	< 10.0	
I Dl	Pyrene	129-00-0	10.0	< 10.0	
Jose Rocha	Pyridine	110-86-1	10.0	< 10.0	
QA Officer	Quinoline	91-22-5	10.0	< 10.0	
	Safrole	94-59-7	10.0	< 10.0	
	Tetraethyl dithiopyrophosphate	3689-24-5	10.0	< 10.0	
	Thionazin	297-97-2	10.0	< 10.0	
	Surr: 2,4,6-Tribromophenol	118-79-6	10-159	57.8	
	Surr: 2-Fluorobiphenyl	321-60-8	10-124	23.0	
	Surr: 2-Fluorophenol	367-12-4	14-106	20.3	
	Surr: Nitrobenzene-d5	4165-60-0	10-180	12.0	
	Surr: Phenol-d6	13127-88-3	10-122	17.1	
	Surr: Terphenyl-d14	1718-51-0	10-199	113	

 $\label{thm:continuous} \textit{The sample was analyzed for TICs and no unknown peaks were detected.}$ 



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006F

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/3/2010 0049h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1´-Biphenyl	92-52-4	10.0	< 10.0	
	1,2,4,5-Tetrachlorobenzene	95-94-3	10.0	< 10.0	
	1,2,4-Trichlorobenzene	120-82-1	10.0	< 10.0	
	1,2-Dichlorobenzene	95-50-1	10.0	< 10.0	
(801) 263-8686	1,3,5-Trinitrobenzene	99-35-4	10.0	< 10.0	1
Toll Free (888) 263-8686	1,4-Naphthoquinone	130-15-4	10.0	< 10.0	
Fax (801)263-8687	1,3-Dichlorobenzene	541-73-1	10.0	< 10.0	
email: awal@awal-labs.com	1,3-Dinitrobenzene	99-65-0	10.0	< 10.0	
	1,4-Dichlorobenzene	106-46-7	10.0	< 10.0	
Kyle F. Gross	1,4-Phenylenediamine	106-50-3	10.0	< 10.0	
ř	1-Chloronaphthalene	90-13-1	10.0	< 10.0	
Laboratory Director	1-Methylnaphthalene	90-12-0	10.0	< 10.0	
	1-Naphthylamine	134-32-7	10.0	< 10.0	
Jose Rocha	2,3,4,6-Tetrachlorophenol	58-90-2	10.0	< 10.0	
QA Officer	2,4,5-Trichlorophenol	95-95-4	10.0	< 10.0	
	2,4,6-Trichlorophenol	88-06-2	10.0	< 10.0	
	2,4-Dichlorophenol	120-83-2	10.0	< 10.0	
	2,4-Dimethylphenol	105-67-9	10.0	< 10.0	
	2,4-Dinitrophenol	51-28-5	20.0	< 20.0	
	2,4-Dinitrotoluene	121-14-2	10.0	< 10.0	
	2,6-Dichlorophenol	87-65-0	10.0	< 10.0	
	2,6-Dinitrotoluene	606-20-2	10.0	< 10.0	
	2-Acetylaminofluorene	53-96-3	10.0	< 10.0	1
	2-Chloronaphthalene	91-58-7	10.0	< 10.0	
	2-Chlorophenol	95-57-8	10.0	< 10.0	
	2-Methylnaphthalene	91-57-6	10.0	< 10.0	
	2-Methylphenol	95-48-7	10.0	< 10.0	
	2-Naphthylamine	91-59-8	10.0	< 10.0	
	2-Nitroaniline	88-74-4	10.0	< 10.0	

Report Date: 12/6/2010 Page 51 of 162



**ANALYTICAL** 

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006F

AMERICAN Client Sample ID: 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/3/2010 0049h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	2-Nitrophenol	88-75-5	10.0	< 10.0	
	2-Picoline	109-06-8	10.0	< 10.0	
	3&4-Methylphenol		10.0	< 10.0	
	3,3'-Dichlorobenzidine	91-94-1	10.0	< 10.0	
(801) 263-8686	3,3'-Dimethylbenzidine	119-93-7	10.0	< 10.0	
Toll Free (888) 263-8686	3-Methylcholanthrene	56-49-5	10.0	< 10.0	
Fax (801)263-8687	3-Nitroaniline	99-09-2	10.0	< 10.0	
email: awal@awal-labs.com	4,6-Dinitro-2-methylphenol	534-52-1	10.0	< 10.0	
	4-Aminobiphenyl	92-67-1	10.0	< 10.0	
Kyle F. Gross	4-Bromophenyl phenyl ether	101-55-3	10.0	< 10.0	
Laboratory Director	4-Chloro-3-methylphenol	59-50-7	10.0	< 10.0	
Laboratory Director	4-Chloroaniline	106-47-8	10.0	< 10.0	
T D 1	4-Chlorophenyl phenyl ether	7005-72-3	10.0	< 10.0	
Jose Rocha	4-Nitroaniline	100-01-6	10.0	< 10.0	
QA Officer	4-Nitrophenol	100-02-7	10.0	< 10.0	1
	5-Nitro-o-toluidine	99-55-8	10.0	< 10.0	
	7,12-Dimethylbenz(a)anthracene	57-97-6	10.0	< 10.0	
	a,a-Dimethylphenethylamine	122-09-8	10.0	< 10.0	
	Acenaphthene	83-32-9	10.0	< 10.0	
	Acenaphthylene	208-96-8	10.0	< 10.0	
	Acetophenone	98-86-2	10.0	< 10.0	
	alpha-Terpineol	98-55-5	10.0	< 10.0	
	Aniline	62-53-3	10.0	< 10.0	
	Anthracene	120-12-7	10.0	< 10.0	
	Aramite	140-57-8	10.0	< 10.0	
	Azobenzene	103-33-3	10.0	< 10.0	
	Benz(a)anthracene	56-55-3	10.0	< 10.0	
	Benzidine	92-87-5	10.0	< 10.0	
	Benzo(a)pyrene	50-32-8	10.0	< 10.0	

Report Date: 12/6/2010 Page 52 of 162



**ANALYTICAL** 

**LABORATORIES** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006F

AMERICAN Client Sample ID: 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/3/2010 0049h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

Method Used: EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Benzo(b)fluoranthene	205-99-2	10.0	< 10.0	
01113	Benzo(g,h,i)perylene	191-24-2	10.0	< 10.0	
	Benzo(k)fluoranthene	207-08-9	10.0	< 10.0	
	Benzoic acid	65-85-0	20.0	< 20.0	
(801) 263-8686	Benzyl alcohol	100-51-6	10.0	< 10.0	
Toll Free (888) 263-8686	Bis(2-chloroethoxy)methane	111-91-1	10.0	< 10.0	
Fax (801)263-8687	Bis(2-chloroethyl) ether	111-44-4	10.0	< 10.0	
email: awal@awal-labs.com	Bis(2-chloroisopropyl) ether	108-60-1	10.0	< 10.0	
	Bis(2-ethylhexyl) phthalate	117-81-7	10.0	< 10.0	1@
Kyle F. Gross	bis(2-ethylhexyl)adipate	103-23-1	10.0	< 10.0	
•	Butyl benzyl phthalate	85-68-7	10.0	< 10.0	
Laboratory Director	Carbazole	86-74-8	10.0	< 10.0	
. D. 1	Chlorobenzilate	510-15-6	10.0	< 10.0	
Jose Rocha	Chrysene	218-01-9	10.0	< 10.0	
QA Officer	Di-n-butyl phthalate	84-74-2	10.0	< 10.0	
	Di-n-octyl phthalate	117-84-0	10.0	< 10.0	1
	Diallate (cis or trans)	2303-16-4	10.0	< 10.0	
	Dibenz(a,h)anthracene	53-70-3	10.0	< 10.0	
	Dibenzofuran	132-64-9	10.0	< 10.0	
	Diethyl phthalate	84-66-2	10.0	< 10.0	
	Dimethoate	60-51-5	10.0	< 10.0	
	Dimethyl phthalate	131-11-3	10.0	< 10.0	
	Dimethylaminoazobenzene	60-11-7	10.0	< 10.0	
	Dinoseb	88-85-7	10.0	< 10.0	
	Diphenylamine	122-39-4	10.0	< 10.0	
	Disulfoton	298-04-4	10.0	< 10.0	
	Ethyl methanesulfonate	62-50-0	10.0	< 10.0	
	Famphur	52-85-7	10.0	< 10.0	1
	Fluoranthene	206-44-0	10.0	< 10.0	

Report Date: 12/6/2010 Page 53 of 162



**ANALYTICAL** 

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006F

AMERICAN Client Sample ID: 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/3/2010 0049h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Fluorene	86-73-7	10.0	< 10.0	
01113	Hexachlorobenzene	118-74-1	10.0	< 10.0	
	Hexachlorobutadiene	87-68-3	10.0	< 10.0	
	Hexachlorocyclopentadiene	77-47-4	10.0	< 10.0	
(801) 263-8686	Hexachloroethane	67-72-1	10.0	< 10.0	
Toll Free (888) 263-8686	Hexachlorophene	70-30-4	10.0	< 10.0	
Fax (801)263-8687	Hexachloropropene	1888-71-7	10.0	< 10.0	
email: awal@awal-labs.com	Indene	95-13-6	10.0	< 10.0	1
	Indeno(1,2,3-cd)pyrene	193-39-5	10.0	< 10.0	
Kyle F. Gross	Isodrin	465-73-6	10.0	< 10.0	
·	Isophorone	78-59-1	10.0	< 10.0	
Laboratory Director	Isosafrole	120-58-1	10.0	< 10.0	
r D 1	Kepone	143-50-0	10.0	< 10.0	1
Jose Rocha	Methapyrilene	91-80-5	10.0	< 10.0	
QA Officer	Methyl methanesulfonate	66-27-3	10.0	< 10.0	
	n-Decane	124-18-5	10.0	< 10.0	
	N-Nitrosodi-n-butylamine	924-16-3	10.0	< 10.0	
	N-Nitrosodiethylamine	55-18-5	10.0	< 10.0	
	N-Nitrosodimethylamine	62-75-9	10.0	< 10.0	
	N-Nitrosodiphenylamine	86-30-6	10.0	< 10.0	
	N-Nitrosodi-n-propylamine	621-64-7	10.0	< 10.0	
	N-Nitrosomethylethylamine	10595-95-6	10.0	< 10.0	
	N-Nitrosomorpholine	59-89-2	10.0	< 10.0	
	N-Nitrosopiperidine	100-75-4	10.0	< 10.0	
	N-Nitrosopyrrolidine	930-55-2	10.0	< 10.0	
	n-Octadecane	593-45-3	10.0	< 10.0	
	Naphthalene	91-20-3	10.0	< 10.0	
	Nitrobenzene	98-95-3	10.0	< 10.0	
	Nitroquinoline-1-oxide	56-57-5	10.0	< 10.0	

Report Date: 12/6/2010 Page 54 of 162



ANALYTICAL

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006F

AMERICAN Client Sample ID: 4992083 Red Butte @ 1100 E.

 Collection Date:
 11/30/2010 1315h
 Analyzed: 12/3/2010 0049h

 Received Date:
 11/30/2010 1538h
 Extracted: 12/2/2010 0923h

**Method Used:** EPA625

Analytical Results SVOA by GC/MS Method 625/3510

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	O,O,O-Triethyl phosphorothioate	126-68-1	10.0	< 10.0	
	o-Toluidine	95-53-4	10.0	< 10.0	1
	Parathion	56-38-2	10.0	< 10.0	
	Methyl parathion	298-00-0	10.0	< 10.0	
(801) 263-8686	Pentachlorobenzene	608-93-5	10.0	< 10.0	
Toll Free (888) 263-8686	Pentachloronitrobenzene	82-68-8	10.0	< 10.0	
Fax (801)263-8687	Pentachlorophenol	87-86-5	10.0	< 10.0	
email: awal@awal-labs.com	Phenacetin	62-44-2	10.0	< 10.0	
	Phenanthrene	85-01-8	10.0	< 10.0	
Kyle F. Gross	Phenol	108-95-2	10.0	< 10.0	
Laboratory Director	Phorate	298-02-2	10.0	< 10.0	
Laboratory Director	Pronamide	23950-58-5	10.0	< 10.0	1
Jose Rocha	Pyrene	129-00-0	10.0	< 10.0	
	Pyridine	110-86-1	10.0	< 10.0	
QA Officer	Quinoline	91-22-5	10.0	< 10.0	1
	Safrole	94-59-7	10.0	< 10.0	
	Tetraethyl dithiopyrophosphate	3689-24-5	10.0	< 10.0	
	Thionazin	297-97-2	10.0	< 10.0	
	Surr: 2,4,6-Tribromophenol	118-79-6	10-159	52.3	
	Surr: 2-Fluorobiphenyl	321-60-8	10-124	42.8	
	Surr: 2-Fluorophenol	367-12-4	14-106	28.4	
	Surr: Nitrobenzene-d5	4165-60-0	10-180	35.9	
	Surr: Phenol-d6	13127-88-3	10-122	25.7	
	Surr: Terphenyl-d14	1718-51-0	10-199	107	

<sup>&</sup>lt;sup>1</sup> - Matrix spike recovery indicates matrix interference. The method is in control as indicated by the LCS.

The sample was analyzed for TICs and no unknown peaks were detected.

<sup>@ -</sup> High RPD due to suspected sample non-homogeneity or matrix interference.



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001B

Client Sample ID: 4992095 Red Butte AB Garden

**Collection Date:** 11/30/2010 0950h **Analyzed:** 12/2/2010 0827h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1,1,2-Tetrachloroethane	630-20-6	2.00	< 2.00	
	1,1,1-Trichloroethane	71-55-6	2.00	< 2.00	
	1,1,2,2-Tetrachloroethane	79-34-5	2.00	< 2.00	
	1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	2.00	< 2.00	
(801) 263-8686	1,1,2-Trichloroethane	79-00-5	2.00	< 2.00	
Toll Free (888) 263-8686	1,1-Dichloropropene	563-58-6	2.00	< 2.00	
Fax (801)263-8687	1,1-Dichloroethane	75-34-3	2.00	< 2.00	
email: awal@awal-labs.com	1,1-Dichloroethene	75-35-4	2.00	< 2.00	
	1,2,3-Trichlorobenzene	87-61-6	2.00	< 2.00	
Kyle F. Gross	1,2,3-Trichloropropane	96-18-4	2.00	< 2.00	
Laboratory Director	1,2,3-Trimethylbenzene	526-73-8	2.00	< 2.00	
Laboratory Director	1,2,4-Trichlorobenzene	120-82-1	2.00	< 2.00	
T D 1	1,2,4-Trimethylbenzene	95-63-6	2.00	< 2.00	
Jose Rocha	1,2-Dibromo-3-chloropropane	96-12-8	2.00	< 2.00	
QA Officer	1,2-Dibromoethane	106-93-4	2.00	< 2.00	
	1,2-Dichlorobenzene	95-50-1	2.00	< 2.00	
	1,2-Dichloroethane	107-06-2	2.00	< 2.00	
	1,2-Dichloropropane	78-87-5	2.00	< 2.00	
	1,3,5-Trimethylbenzene	108-67-8	2.00	< 2.00	
	1,3-Dichlorobenzene	541-73-1	2.00	< 2.00	
	1,3-Dichloropropane	142-28-9	2.00	< 2.00	
	1,4-Dichlorobenzene	106-46-7	2.00	< 2.00	
	1,4-Dioxane	123-91-1	40.0	< 40.0	
	2,2-Dichloropropane	594-20-7	2.00	< 2.00	
	2-Butanone	78-93-3	10.0	< 10.0	
	2-Chloroethyl vinyl ether	110-75-8	5.00	< 5.00	
	2-Chlorotoluene	95-49-8	2.00	< 2.00	
	2-Hexanone	591-78-6	5.00	< 5.00	
	2-Nitropropane	79-46-9	2.00	< 2.00	

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**ANALYTICAL** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001B

Client Sample ID: 4992095 Red Butte AB Garden

**Collection Date:** 11/30/2010 0950h **Analyzed:** 12/2/2010 0827h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	4-Chlorotoluene	106-43-4	2.00	< 2.00	
	4-Isopropyltoluene	99-87-6	2.00	< 2.00	
	4-Methyl-2-pentanone	108-10-1	5.00	< 5.00	
	Acetone	67-64-1	10.0	< 10.0	
(801) 263-8686	Acetonitrile	75-05-8	5.00	< 5.00	
Toll Free (888) 263-8686	Acrolein	107-02-8	5.00	< 5.00	
Fax (801)263-8687	Acrylonitrile	107-13-1	10.0	< 10.0	
email: awal@awal-labs.com	Allyl chloride	107-05-1	5.00	< 5.00	
	Benzene	71-43-2	2.00	< 2.00	
Kyle F. Gross	Benzyl chloride	100-44-7	5.00	< 5.00	
•	Bis(2-chloroisopropyl) ether	108-60-1	5.00	< 5.00	
Laboratory Director	Bromobenzene	108-86-1	2.00	< 2.00	
T D 1	Bromochloromethane	74-97-5	2.00	< 2.00	
Jose Rocha	Bromodichloromethane	75-27-4	2.00	< 2.00	
QA Officer	Bromoform	75-25-2	2.00	< 2.00	
	Bromomethane	74-83-9	5.00	< 5.00	
	Butyl acetate	123-86-4	5.00	< 5.00	
	Carbon disulfide	75-15-0	2.00	< 2.00	
	Carbon tetrachloride	56-23-5	2.00	< 2.00	
	Chlorobenzene	108-90-7	2.00	< 2.00	
	Chloroethane	75-00-3	2.00	< 2.00	
	Chloroform	67-66-3	2.00	< 2.00	
	Chloromethane	74-87-3	5.00	< 5.00	
	Chloroprene	126-99-8	2.00	< 2.00	
	cis-1,2-Dichloroethene	156-59-2	2.00	< 2.00	
	cis-1,3-Dichloropropene	10061-01-5	2.00	< 2.00	
	Cyclohexane	110-82-7	2.00	< 2.00	
	Cyclohexanone	108-94-1	50.0	< 50.0	
	Dibromochloromethane	124-48-1	2.00	< 2.00	

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#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001B

Client Sample ID: 4992095 Red Butte AB Garden

**Collection Date:** 11/30/2010 0950h **Analyzed:** 12/2/2010 0827h

**Received Date:** 11/30/2010 1538h

ANALYTICAL Method Used: EPA624

LABORATORIES
Analytical Results

**WEST** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Dibromomethane	74-95-3	2.00	< 2.00	
0.1110	Dichlorodifluoromethane	75-71-8	2.00	< 2.00	
	Ethyl acetate	141-78-6	10.0	< 10.0	
	Ethyl ether	60-29-7	10.0	< 10.0	
(801) 263-8686	Ethyl methacrylate	97-63-2	2.00	< 2.00	
Toll Free (888) 263-8686	Ethylbenzene	100-41-4	2.00	< 2.00	
Fax (801)263-8687	Hexachlorobutadiene	87-68-3	2.00	< 2.00	
email: awal@awal-labs.com	Iodomethane	74-88-4	5.00	< 5.00	
	Isobutyl alcohol	78-83-1	100	< 100	
Kyle F. Gross	Isopropyl acetate	108-21-4	2.00	< 2.00	
•	Isopropyl alcohol	67-63-0	25.0	< 25.0	
Laboratory Director	Isopropylbenzene	98-82-8	2.00	< 2.00	
	Isopropyltoluene	99-87-6	2.00	< 2.00	
Jose Rocha	m,p-Xylene	179601-23-1	2.00	< 2.00	
QA Officer	Methacrylonitrile	126-98-7	5.00	< 5.00	
	Methyl Acetate	79-20-9	5.00	< 5.00	
	Methyl methacrylate	80-62-6	5.00	< 5.00	
	Methyl tert-butyl ether	1634-04-4	2.00	< 2.00	
	Methylcyclohexane	108-87-2	2.00	< 2.00	
	Methylene chloride	75-09-2	2.00	< 2.00	
	n-Amyl acetate	628-63-7	2.00	< 2.00	
	n-Butyl alcohol	71-36-3	25.0	< 25.0	
	n-Butylbenzene	104-51-8	2.00	< 2.00	
	n-Hexane	110-54-3	2.00	< 2.00	
	n-Octane	111-65-9	2.00	< 2.00	
	n-Propylbenzene	103-65-1	2.00	< 2.00	
	Naphthalene	91-20-3	2.00	< 2.00	
	o-Xylene	95-47-6	2.00	< 2.00	
	Pentachloroethane	76-01-7	2.00	< 2.00	

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**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-001B

Client Sample ID: 4992095 Red Butte AB Garden

**Collection Date:** 11/30/2010 0950h **Analyzed:** 12/2/2010 0827h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

**LABORATORIES** 

**ANALYTICAL** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Propionitrile	107-12-0	25.0	< 25.0	
	Propyl acetate	109-60-4	2.00	< 2.00	
	sec-Butylbenzene	135-98-8	2.00	< 2.00	
	Styrene	100-42-5	2.00	< 2.00	
(801) 263-8686	tert-Butyl alcohol	76-65-0	20.0	< 20.0	
Toll Free (888) 263-8686	tert-Butylbenzene	98-06-6	2.00	< 2.00	
Fax (801)263-8687	Tetrachloroethene	127-18-4	2.00	< 2.00	
email: awal@awal-labs.com	Tetrahydrofuran	109-99-9	2.00	< 2.00	
	Toluene	108-88-3	2.00	< 2.00	
Kyle F. Gross	trans-1,2-Dichloroethene	156-60-5	2.00	< 2.00	
Laboratory Director	trans-1,3-Dichloropropene	10061-02-6	2.00	< 2.00	
Laboratory Director	trans-1,4-Dichloro-2-butene	110-57-6	2.00	< 2.00	
r D 1	Trichloroethene	79-01-6	2.00	< 2.00	
Jose Rocha	Trichlorofluoromethane	75-69-4	2.00	< 2.00	
QA Officer	Vinyl acetate	108-05-4	5.00	< 5.00	
	Vinyl chloride	75-01-4	1.00	< 1.00	
	Xylenes, Total	1330-20-7	2.00	< 2.00	
	TPH C11-C15 (DRO)		20.0	< 20.0	
	TPH C6-C10 (GRO)		20.0	< 20.0	
	Surr: 1,2-Dichloroethane-d4	17060-07-0	77-144	102	
	Surr: 4-Bromofluorobenzene	460-00-4	80-123	104	
	Surr: Dibromofluoromethane	1868-53-7	80-124	99.2	
	Surr: Toluene-d8	2037-26-5	80-125	98.9	



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002B

Client Sample ID: 4992085 Red Butte BL Garden

**Collection Date:** 11/30/2010 1030h **Analyzed:** 12/2/2010 0846h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1,1,2-Tetrachloroethane	630-20-6	2.00	< 2.00	
	1,1,1-Trichloroethane	71-55-6	2.00	< 2.00	
	1,1,2,2-Tetrachloroethane	79-34-5	2.00	< 2.00	
	1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	2.00	< 2.00	
(801) 263-8686	1,1,2-Trichloroethane	79-00-5	2.00	< 2.00	
Toll Free (888) 263-8686	1,1-Dichloropropene	563-58-6	2.00	< 2.00	
Fax (801)263-8687	1,1-Dichloroethane	75-34-3	2.00	< 2.00	
email: awal@awal-labs.com	1,1-Dichloroethene	75-35-4	2.00	< 2.00	
	1,2,3-Trichlorobenzene	87-61-6	2.00	< 2.00	
Kyle F. Gross	1,2,3-Trichloropropane	96-18-4	2.00	< 2.00	
Laboratory Director	1,2,3-Trimethylbenzene	526-73-8	2.00	< 2.00	
Laboratory Director	1,2,4-Trichlorobenzene	120-82-1	2.00	< 2.00	
T D 1	1,2,4-Trimethylbenzene	95-63-6	2.00	< 2.00	
Jose Rocha	1,2-Dibromo-3-chloropropane	96-12-8	2.00	< 2.00	
QA Officer	1,2-Dibromoethane	106-93-4	2.00	< 2.00	
	1,2-Dichlorobenzene	95-50-1	2.00	< 2.00	
	1,2-Dichloroethane	107-06-2	2.00	< 2.00	
	1,2-Dichloropropane	78-87-5	2.00	< 2.00	
	1,3,5-Trimethylbenzene	108-67-8	2.00	< 2.00	
	1,3-Dichlorobenzene	541-73-1	2.00	< 2.00	
	1,3-Dichloropropane	142-28-9	2.00	< 2.00	
	1,4-Dichlorobenzene	106-46-7	2.00	< 2.00	
	1,4-Dioxane	123-91-1	40.0	< 40.0	
	2,2-Dichloropropane	594-20-7	2.00	< 2.00	
	2-Butanone	78-93-3	10.0	< 10.0	
	2-Chloroethyl vinyl ether	110-75-8	5.00	< 5.00	
	2-Chlorotoluene	95-49-8	2.00	< 2.00	
	2-Hexanone	591-78-6	5.00	< 5.00	
	2-Nitropropane	79-46-9	2.00	< 2.00	

Report Date: 12/6/2010 Page 60 of 162



**ANALYTICAL** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002B

Client Sample ID: 4992085 Red Butte BL Garden

**Collection Date:** 11/30/2010 1030h **Analyzed:** 12/2/2010 0846h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES

Analytical Results VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	4-Chlorotoluene	106-43-4	2.00	< 2.00	
	4-Isopropyltoluene	99-87-6	2.00	< 2.00	
	4-Methyl-2-pentanone	108-10-1	5.00	< 5.00	
	Acetone	67-64-1	10.0	< 10.0	
(801) 263-8686	Acetonitrile	75-05-8	5.00	< 5.00	
Toll Free (888) 263-8686	Acrolein	107-02-8	5.00	< 5.00	
Fax (801)263-8687	Acrylonitrile	107-13-1	10.0	< 10.0	
email: awal@awal-labs.com	Allyl chloride	107-05-1	5.00	< 5.00	
	Benzene	71-43-2	2.00	< 2.00	
Kyle F. Gross	Benzyl chloride	100-44-7	5.00	< 5.00	
Laboratory Director	Bis(2-chloroisopropyl) ether	108-60-1	5.00	< 5.00	
Laboratory Director	Bromobenzene	108-86-1	2.00	< 2.00	
T D 1	Bromochloromethane	74-97-5	2.00	< 2.00	
Jose Rocha	Bromodichloromethane	75-27-4	2.00	< 2.00	
QA Officer	Bromoform	75-25-2	2.00	< 2.00	
	Bromomethane	74-83-9	5.00	< 5.00	
	Butyl acetate	123-86-4	5.00	< 5.00	
	Carbon disulfide	75-15-0	2.00	< 2.00	
	Carbon tetrachloride	56-23-5	2.00	< 2.00	
	Chlorobenzene	108-90-7	2.00	< 2.00	
	Chloroethane	75-00-3	2.00	< 2.00	
	Chloroform	67-66-3	2.00	< 2.00	
	Chloromethane	74-87-3	5.00	< 5.00	
	Chloroprene	126-99-8	2.00	< 2.00	
	cis-1,2-Dichloroethene	156-59-2	2.00	< 2.00	
	cis-1,3-Dichloropropene	10061-01-5	2.00	< 2.00	
	Cyclohexane	110-82-7	2.00	< 2.00	
	Cyclohexanone	108-94-1	50.0	< 50.0	
	Dibromochloromethane	124-48-1	2.00	< 2.00	

Report Date: 12/6/2010 Page 61 of 162



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002B

Client Sample ID: 4992085 Red Butte BL Garden

**Collection Date:** 11/30/2010 1030h **Analyzed:** 12/2/2010 0846h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Dibromomethane	74-95-3	2.00	< 2.00	
04113	Dichlorodifluoromethane	75-71-8	2.00	< 2.00	
	Ethyl acetate	141-78-6	10.0	< 10.0	
	Ethyl ether	60-29-7	10.0	< 10.0	
(801) 263-8686	Ethyl methacrylate	97-63-2	2.00	< 2.00	
Toll Free (888) 263-8686	Ethylbenzene	100-41-4	2.00	< 2.00	
Fax (801)263-8687	Hexachlorobutadiene	87-68-3	2.00	< 2.00	
email: awal@awal-labs.com	Iodomethane	74-88-4	5.00	< 5.00	
	Isobutyl alcohol	78-83-1	100	< 100	
Wala E. Carre	Isopropyl acetate	108-21-4	2.00	< 2.00	
Kyle F. Gross	Isopropyl alcohol	67-63-0	25.0	< 25.0	
Laboratory Director	Isopropylbenzene	98-82-8	2.00	< 2.00	
	Isopropyltoluene	99-87-6	2.00	< 2.00	
Jose Rocha	m,p-Xylene	179601-23-1	2.00	< 2.00	
QA Officer	Methacrylonitrile	126-98-7	5.00	< 5.00	
	Methyl Acetate	79-20-9	5.00	< 5.00	
	Methyl methacrylate	80-62-6	5.00	< 5.00	
	Methyl tert-butyl ether	1634-04-4	2.00	< 2.00	
	Methylcyclohexane	108-87-2	2.00	< 2.00	
	Methylene chloride	75-09-2	2.00	< 2.00	
	n-Amyl acetate	628-63-7	2.00	< 2.00	
	n-Butyl alcohol	71-36-3	25.0	< 25.0	
	n-Butylbenzene	104-51-8	2.00	< 2.00	
	n-Hexane	110-54-3	2.00	< 2.00	
	n-Octane	111-65-9	2.00	< 2.00	
	n-Propylbenzene	103-65-1	2.00	< 2.00	
	Naphthalene	91-20-3	2.00	< 2.00	
	o-Xylene	95-47-6	2.00	< 2.00	
	Pentachloroethane	76-01-7	2.00	< 2.00	

Report Date: 12/6/2010 Page 62 of 162



#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-002B

AMERICAN Client Sample ID: 4992085 Red Butte BL Garden

**Collection Date:** 11/30/2010 1030h **Analyzed:** 12/2/2010 0846h

**Received Date:** 11/30/2010 1538h

**Method Used:** EPA624

LABORATORIES
Analytical Results

**ANALYTICAL** 

**WEST** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Propionitrile	107-12-0	25.0	< 25.0	
	Propyl acetate	109-60-4	2.00	< 2.00	
	sec-Butylbenzene	135-98-8	2.00	< 2.00	
	Styrene	100-42-5	2.00	< 2.00	
(801) 263-8686	tert-Butyl alcohol	76-65-0	20.0	< 20.0	
Toll Free (888) 263-8686	tert-Butylbenzene	98-06-6	2.00	< 2.00	
Fax (801)263-8687	Tetrachloroethene	127-18-4	2.00	< 2.00	
email: awal@awal-labs.com	Tetrahydrofuran	109-99-9	2.00	< 2.00	
	Toluene	108-88-3	2.00	< 2.00	
Kyle F. Gross	trans-1,2-Dichloroethene	156-60-5	2.00	< 2.00	
Laboratory Director	trans-1,3-Dichloropropene	10061-02-6	2.00	< 2.00	
Laboratory Director	trans-1,4-Dichloro-2-butene	110-57-6	2.00	< 2.00	
I Dl	Trichloroethene	79-01-6	2.00	< 2.00	
Jose Rocha	Trichlorofluoromethane	75-69-4	2.00	< 2.00	
QA Officer	Vinyl acetate	108-05-4	5.00	< 5.00	
	Vinyl chloride	75-01-4	1.00	< 1.00	
	Xylenes, Total	1330-20-7	2.00	< 2.00	
	TPH C11-C15 (DRO)		20.0	< 20.0	
	TPH C6-C10 (GRO)		20.0	< 20.0	
	Surr: 1,2-Dichloroethane-d4	17060-07-0	77-144	103	
	Surr: 4-Bromofluorobenzene	460-00-4	80-123	106	
	Surr: Dibromofluoromethane	1868-53-7	80-124	100	
	Surr: Toluene-d8	2037-26-5	80-125	99.4	



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003B

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

**Collection Date:** 11/30/2010 1215h **Analyzed:** 12/2/2010 0905h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1,1,2-Tetrachloroethane	630-20-6	2.00	< 2.00	
01113	1,1,1-Trichloroethane	71-55-6	2.00	< 2.00	
	1,1,2,2-Tetrachloroethane	79-34-5	2.00	< 2.00	
	1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	2.00	< 2.00	
(801) 263-8686	1,1,2-Trichloroethane	79-00-5	2.00	< 2.00	
Toll Free (888) 263-8686	1,1-Dichloropropene	563-58-6	2.00	< 2.00	
Fax (801)263-8687	1,1-Dichloroethane	75-34-3	2.00	< 2.00	
email: awal@awal-labs.com	1,1-Dichloroethene	75-35-4	2.00	< 2.00	
	1,2,3-Trichlorobenzene	87-61-6	2.00	< 2.00	
Kyle F. Gross	1,2,3-Trichloropropane	96-18-4	2.00	< 2.00	
•	1,2,3-Trimethylbenzene	526-73-8	2.00	< 2.00	
Laboratory Director	1,2,4-Trichlorobenzene	120-82-1	2.00	< 2.00	
. D. 1	1,2,4-Trimethylbenzene	95-63-6	2.00	< 2.00	
Jose Rocha	1,2-Dibromo-3-chloropropane	96-12-8	2.00	< 2.00	
QA Officer	1,2-Dibromoethane	106-93-4	2.00	< 2.00	
	1,2-Dichlorobenzene	95-50-1	2.00	< 2.00	
	1,2-Dichloroethane	107-06-2	2.00	< 2.00	
	1,2-Dichloropropane	78-87-5	2.00	< 2.00	
	1,3,5-Trimethylbenzene	108-67-8	2.00	< 2.00	
	1,3-Dichlorobenzene	541-73-1	2.00	< 2.00	
	1,3-Dichloropropane	142-28-9	2.00	< 2.00	
	1,4-Dichlorobenzene	106-46-7	2.00	< 2.00	
	1,4-Dioxane	123-91-1	40.0	< 40.0	
	2,2-Dichloropropane	594-20-7	2.00	< 2.00	
	2-Butanone	78-93-3	10.0	< 10.0	
	2-Chloroethyl vinyl ether	110-75-8	5.00	< 5.00	
	2-Chlorotoluene	95-49-8	2.00	< 2.00	
	2-Hexanone	591-78-6	5.00	< 5.00	
	2-Nitropropane	79-46-9	2.00	< 2.00	

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**ANALYTICAL** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003B

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

**Collection Date:** 11/30/2010 1215h **Analyzed:** 12/2/2010 0905h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	4-Chlorotoluene	106-43-4	2.00	< 2.00	
	4-Isopropyltoluene	99-87-6	2.00	< 2.00	
	4-Methyl-2-pentanone	108-10-1	5.00	< 5.00	
	Acetone	67-64-1	10.0	< 10.0	
(801) 263-8686	Acetonitrile	75-05-8	5.00	< 5.00	
Toll Free (888) 263-8686	Acrolein	107-02-8	5.00	< 5.00	
Fax (801)263-8687	Acrylonitrile	107-13-1	10.0	< 10.0	
email: awal@awal-labs.com	Allyl chloride	107-05-1	5.00	< 5.00	
	Benzene	71-43-2	2.00	< 2.00	
Kyle F. Gross	Benzyl chloride	100-44-7	5.00	< 5.00	
Laboratory Director	Bis(2-chloroisopropyl) ether	108-60-1	5.00	< 5.00	
Laboratory Director	Bromobenzene	108-86-1	2.00	< 2.00	
T D 1	Bromochloromethane	74-97-5	2.00	< 2.00	
Jose Rocha	Bromodichloromethane	75-27-4	2.00	< 2.00	
QA Officer	Bromoform	75-25-2	2.00	< 2.00	
	Bromomethane	74-83-9	5.00	< 5.00	
	Butyl acetate	123-86-4	5.00	< 5.00	
	Carbon disulfide	75-15-0	2.00	< 2.00	
	Carbon tetrachloride	56-23-5	2.00	< 2.00	
	Chlorobenzene	108-90-7	2.00	< 2.00	
	Chloroethane	75-00-3	2.00	< 2.00	
	Chloroform	67-66-3	2.00	< 2.00	
	Chloromethane	74-87-3	5.00	< 5.00	
	Chloroprene	126-99-8	2.00	< 2.00	
	cis-1,2-Dichloroethene	156-59-2	2.00	< 2.00	
	cis-1,3-Dichloropropene	10061-01-5	2.00	< 2.00	
	Cyclohexane	110-82-7	2.00	< 2.00	
	Cyclohexanone	108-94-1	50.0	< 50.0	
	Dibromochloromethane	124-48-1	2.00	< 2.00	

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**ANALYTICAL** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003B

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

**Collection Date:** 11/30/2010 1215h **Analyzed:** 12/2/2010 0905h

**Received Date:** 11/30/2010 1538h

**Method Used:** EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Dibromomethane	74-95-3	2.00	< 2.00	
	Dichlorodifluoromethane	75-71-8	2.00	< 2.00	
	Ethyl acetate	141-78-6	10.0	< 10.0	
	Ethyl ether	60-29-7	10.0	< 10.0	
(801) 263-8686	Ethyl methacrylate	97-63-2	2.00	< 2.00	
Toll Free (888) 263-8686	Ethylbenzene	100-41-4	2.00	< 2.00	
Fax (801)263-8687	Hexachlorobutadiene	87-68-3	2.00	< 2.00	
email: awal@awal-labs.com	Iodomethane	74-88-4	5.00	< 5.00	
	Isobutyl alcohol	78-83-1	100	< 100	
Kyle F. Gross	Isopropyl acetate	108-21-4	2.00	< 2.00	
•	Isopropyl alcohol	67-63-0	25.0	< 25.0	
Laboratory Director	Isopropylbenzene	98-82-8	2.00	< 2.00	
T D 1	Isopropyltoluene	99-87-6	2.00	< 2.00	
Jose Rocha	m,p-Xylene	179601-23-1	2.00	< 2.00	
QA Officer	Methacrylonitrile	126-98-7	5.00	< 5.00	
	Methyl Acetate	79-20-9	5.00	< 5.00	
	Methyl methacrylate	80-62-6	5.00	< 5.00	
	Methyl tert-butyl ether	1634-04-4	2.00	< 2.00	
	Methylcyclohexane	108-87-2	2.00	< 2.00	
	Methylene chloride	75-09-2	2.00	< 2.00	
	n-Amyl acetate	628-63-7	2.00	< 2.00	
	n-Butyl alcohol	71-36-3	25.0	< 25.0	
	n-Butylbenzene	104-51-8	2.00	< 2.00	
	n-Hexane	110-54-3	2.00	< 2.00	
	n-Octane	111-65-9	2.00	< 2.00	
	n-Propylbenzene	103-65-1	2.00	< 2.00	
	Naphthalene	91-20-3	2.00	< 2.00	
	o-Xylene	95-47-6	2.00	< 2.00	
	Pentachloroethane	76-01-7	2.00	< 2.00	

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**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-003B

**Client Sample ID:** 4992084 Red Butte @ 1500 E.

**Collection Date:** 11/30/2010 1215h **Analyzed:** 12/2/2010 0905h

**Received Date:** 11/30/2010 1538h

**Method Used:** EPA624

**Analytical Results** 

LABORATORIES

**ANALYTICAL** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Propionitrile	107-12-0	25.0	< 25.0	
	Propyl acetate	109-60-4	2.00	< 2.00	
	sec-Butylbenzene	135-98-8	2.00	< 2.00	
	Styrene	100-42-5	2.00	< 2.00	
(801) 263-8686	tert-Butyl alcohol	76-65-0	20.0	< 20.0	
Toll Free (888) 263-8686	tert-Butylbenzene	98-06-6	2.00	< 2.00	
Fax (801)263-8687	Tetrachloroethene	127-18-4	2.00	< 2.00	
email: awal@awal-labs.com	Tetrahydrofuran	109-99-9	2.00	< 2.00	
	Toluene	108-88-3	2.00	< 2.00	
Kyle F. Gross	trans-1,2-Dichloroethene	156-60-5	2.00	< 2.00	
Laboratory Director	trans-1,3-Dichloropropene	10061-02-6	2.00	< 2.00	
Laboratory Director	trans-1,4-Dichloro-2-butene	110-57-6	2.00	< 2.00	
r D 1	Trichloroethene	79-01-6	2.00	< 2.00	
Jose Rocha	Trichlorofluoromethane	75-69-4	2.00	< 2.00	
QA Officer	Vinyl acetate	108-05-4	5.00	< 5.00	
	Vinyl chloride	75-01-4	1.00	< 1.00	
	Xylenes, Total	1330-20-7	2.00	< 2.00	
	TPH C11-C15 (DRO)		20.0	< 20.0	
	TPH C6-C10 (GRO)		20.0	< 20.0	
	Surr: 1,2-Dichloroethane-d4	17060-07-0	77-144	105	
	Surr: 4-Bromofluorobenzene	460-00-4	80-123	104	
	Surr: Dibromofluoromethane	1868-53-7	80-124	101	
	Surr: Toluene-d8	2037-26-5	80-125	98.6	



**ANALYTICAL** 

**LABORATORIES** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004B

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

**Collection Date:** 11/30/2010 1405h **Analyzed:** 12/2/2010 0924h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1,1,2-Tetrachloroethane	630-20-6	2.00	< 2.00	
0.220	1,1,1-Trichloroethane	71-55-6	2.00	< 2.00	
	1,1,2,2-Tetrachloroethane	79-34-5	2.00	< 2.00	
	1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	2.00	< 2.00	
(801) 263-8686	1,1,2-Trichloroethane	79-00-5	2.00	< 2.00	
Toll Free (888) 263-8686	1,1-Dichloropropene	563-58-6	2.00	< 2.00	
Fax (801)263-8687	1,1-Dichloroethane	75-34-3	2.00	< 2.00	
email: awal@awal-labs.com	1,1-Dichloroethene	75-35-4	2.00	< 2.00	
	1,2,3-Trichlorobenzene	87-61-6	2.00	< 2.00	
Kyle F. Gross	1,2,3-Trichloropropane	96-18-4	2.00	< 2.00	
Laboratory Director	1,2,3-Trimethylbenzene	526-73-8	2.00	< 2.00	
Laboratory Director	1,2,4-Trichlorobenzene	120-82-1	2.00	< 2.00	
T D 1	1,2,4-Trimethylbenzene	95-63-6	2.00	< 2.00	
Jose Rocha	1,2-Dibromo-3-chloropropane	96-12-8	2.00	< 2.00	
QA Officer	1,2-Dibromoethane	106-93-4	2.00	< 2.00	
	1,2-Dichlorobenzene	95-50-1	2.00	< 2.00	
	1,2-Dichloroethane	107-06-2	2.00	< 2.00	
	1,2-Dichloropropane	78-87-5	2.00	< 2.00	
	1,3,5-Trimethylbenzene	108-67-8	2.00	< 2.00	
	1,3-Dichlorobenzene	541-73-1	2.00	< 2.00	
	1,3-Dichloropropane	142-28-9	2.00	< 2.00	
	1,4-Dichlorobenzene	106-46-7	2.00	< 2.00	
	1,4-Dioxane	123-91-1	40.0	< 40.0	
	2,2-Dichloropropane	594-20-7	2.00	< 2.00	
	2-Butanone	78-93-3	10.0	< 10.0	
	2-Chloroethyl vinyl ether	110-75-8	5.00	< 5.00	
	2-Chlorotoluene	95-49-8	2.00	< 2.00	
	2-Hexanone	591-78-6	5.00	< 5.00	
	2-Nitropropane	79-46-9	2.00	< 2.00	

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# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004B

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

**Collection Date:** 11/30/2010 1405h **Analyzed:** 12/2/2010 0924h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

**WEST** 

**ANALYTICAL** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	4-Chlorotoluene	106-43-4	2.00	< 2.00	
0.220	4-Isopropyltoluene	99-87-6	2.00	< 2.00	
	4-Methyl-2-pentanone	108-10-1	5.00	< 5.00	
	Acetone	67-64-1	10.0	< 10.0	
(801) 263-8686	Acetonitrile	75-05-8	5.00	< 5.00	
Toll Free (888) 263-8686	Acrolein	107-02-8	5.00	< 5.00	
Fax (801)263-8687	Acrylonitrile	107-13-1	10.0	< 10.0	
email: awal@awal-labs.com	Allyl chloride	107-05-1	5.00	< 5.00	
	Benzene	71-43-2	2.00	< 2.00	
Kyle F. Gross	Benzyl chloride	100-44-7	5.00	< 5.00	
•	Bis(2-chloroisopropyl) ether	108-60-1	5.00	< 5.00	
Laboratory Director	Bromobenzene	108-86-1	2.00	< 2.00	
T D 1	Bromochloromethane	74-97-5	2.00	< 2.00	
Jose Rocha	Bromodichloromethane	75-27-4	2.00	< 2.00	
QA Officer	Bromoform	75-25-2	2.00	< 2.00	
	Bromomethane	74-83-9	5.00	< 5.00	
	Butyl acetate	123-86-4	5.00	< 5.00	
	Carbon disulfide	75-15-0	2.00	< 2.00	
	Carbon tetrachloride	56-23-5	2.00	< 2.00	
	Chlorobenzene	108-90-7	2.00	< 2.00	
	Chloroethane	75-00-3	2.00	< 2.00	
	Chloroform	67-66-3	2.00	< 2.00	
	Chloromethane	74-87-3	5.00	< 5.00	
	Chloroprene	126-99-8	2.00	< 2.00	
	cis-1,2-Dichloroethene	156-59-2	2.00	< 2.00	
	cis-1,3-Dichloropropene	10061-01-5	2.00	< 2.00	
	Cyclohexane	110-82-7	2.00	< 2.00	
	Cyclohexanone	108-94-1	50.0	< 50.0	
	Dibromochloromethane	124-48-1	2.00	< 2.00	

Report Date: 12/6/2010 Page 69 of 162



#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004B

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

**Collection Date:** 11/30/2010 1405h **Analyzed:** 12/2/2010 0924h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

**WEST** 

**ANALYTICAL** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
Salt Lake City, UT	Dibromomethane	74-95-3	2.00	< 2.00	
84115	Dichlorodifluoromethane	75-71-8	2.00	< 2.00	
		73-71-8 141-78-6	10.0	< 10.0	
	Ethyl acetate				
(801) 263-8686	Ethyl methodrilate	60-29-7 97-63-2	10.0	< 10.0	
Toll Free (888) 263-8686	Ethyl methacrylate	100-41-4	2.00	< 2.00	
	Ethylbenzene		2.00	< 2.00	
Fax (801)263-8687	Hexachlorobutadiene	87-68-3	2.00	< 2.00	
email: awal@awal-labs.com	Iodomethane	74-88-4	5.00	< 5.00	
	Isobutyl alcohol	78-83-1	100	< 100	
Kyle F. Gross	Isopropyl acetate	108-21-4	2.00	< 2.00	
Laboratory Director	Isopropyl alcohol	67-63-0	25.0	< 25.0	
	Isopropylbenzene	98-82-8	2.00	< 2.00	
Jose Rocha	Isopropyltoluene	99-87-6	2.00	< 2.00	
QA Officer	m,p-Xylene	179601-23-1	2.00	< 2.00	
Q11 Officer	Methacrylonitrile	126-98-7	5.00	< 5.00	
	Methyl Acetate	79-20-9	5.00	< 5.00	
	Methyl methacrylate	80-62-6	5.00	< 5.00	
	Methyl tert-butyl ether	1634-04-4	2.00	< 2.00	
	Methylcyclohexane	108-87-2	2.00	< 2.00	
	Methylene chloride	75-09-2	2.00	< 2.00	
	n-Amyl acetate	628-63-7	2.00	< 2.00	
	n-Butyl alcohol	71-36-3	25.0	< 25.0	
	n-Butylbenzene	104-51-8	2.00	< 2.00	
	n-Hexane	110-54-3	2.00	< 2.00	
	n-Octane	111-65-9	2.00	< 2.00	
	n-Propylbenzene	103-65-1	2.00	< 2.00	
	Naphthalene	91-20-3	2.00	< 2.00	
	o-Xylene	95-47-6	2.00	< 2.00	
	Pentachloroethane	76-01-7	2.00	< 2.00	

Report Date: 12/6/2010 Page 70 of 162



# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-004B

AMERICAN Client Sample ID: 4992050 J.R. Below 900 S. Storm Sewer

**Collection Date:** 11/30/2010 1405h **Analyzed:** 12/2/2010 0924h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

**LABORATORIES** 

**ANALYTICAL** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Propionitrile	107-12-0	25.0	< 25.0	
	Propyl acetate	109-60-4	2.00	< 2.00	
	sec-Butylbenzene	135-98-8	2.00	< 2.00	
	Styrene	100-42-5	2.00	< 2.00	
(801) 263-8686	tert-Butyl alcohol	76-65-0	20.0	< 20.0	
Toll Free (888) 263-8686	tert-Butylbenzene	98-06-6	2.00	< 2.00	
Fax (801)263-8687	Tetrachloroethene	127-18-4	2.00	< 2.00	
email: awal@awal-labs.com	Tetrahydrofuran	109-99-9	2.00	< 2.00	
	Toluene	108-88-3	2.00	< 2.00	
Kyle F. Gross	trans-1,2-Dichloroethene	156-60-5	2.00	< 2.00	
Laboratory Director	trans-1,3-Dichloropropene	10061-02-6	2.00	< 2.00	
Laboratory Director	trans-1,4-Dichloro-2-butene	110-57-6	2.00	< 2.00	
r D 1	Trichloroethene	79-01-6	2.00	< 2.00	
Jose Rocha	Trichlorofluoromethane	75-69-4	2.00	< 2.00	
QA Officer	Vinyl acetate	108-05-4	5.00	< 5.00	
	Vinyl chloride	75-01-4	1.00	< 1.00	
	Xylenes, Total	1330-20-7	2.00	< 2.00	
	TPH C11-C15 (DRO)		20.0	< 20.0	
	TPH C6-C10 (GRO)		20.0	< 20.0	
	Surr: 1,2-Dichloroethane-d4	17060-07-0	77-144	108	
	Surr: 4-Bromofluorobenzene	460-00-4	80-123	103	
	Surr: Dibromofluoromethane	1868-53-7	80-124	102	
	Surr: Toluene-d8	2037-26-5	80-125	97.2	



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

#### **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005B

**Client Sample ID:** 4992290 J.R. @ 1700 S.

**Collection Date:** 11/30/2010 1430h **Analyzed:** 12/2/2010 0944h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1,1,2-Tetrachloroethane	630-20-6	2.00	< 2.00	
01113	1,1,1-Trichloroethane	71-55-6	2.00	< 2.00	
	1,1,2,2-Tetrachloroethane	79-34-5	2.00	< 2.00	
	1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	2.00	< 2.00	
(801) 263-8686	1,1,2-Trichloroethane	79-00-5	2.00	< 2.00	
Toll Free (888) 263-8686	1,1-Dichloropropene	563-58-6	2.00	< 2.00	
Fax (801)263-8687	1,1-Dichloroethane	75-34-3	2.00	< 2.00	
email: awal@awal-labs.com	1,1-Dichloroethene	75-35-4	2.00	< 2.00	
	1,2,3-Trichlorobenzene	87-61-6	2.00	< 2.00	
Kyle F. Gross	1,2,3-Trichloropropane	96-18-4	2.00	< 2.00	
•	1,2,3-Trimethylbenzene	526-73-8	2.00	< 2.00	
Laboratory Director	1,2,4-Trichlorobenzene	120-82-1	2.00	< 2.00	
	1,2,4-Trimethylbenzene	95-63-6	2.00	< 2.00	
Jose Rocha	1,2-Dibromo-3-chloropropane	96-12-8	2.00	< 2.00	
QA Officer	1,2-Dibromoethane	106-93-4	2.00	< 2.00	
	1,2-Dichlorobenzene	95-50-1	2.00	< 2.00	
	1,2-Dichloroethane	107-06-2	2.00	< 2.00	
	1,2-Dichloropropane	78-87-5	2.00	< 2.00	
	1,3,5-Trimethylbenzene	108-67-8	2.00	< 2.00	
	1,3-Dichlorobenzene	541-73-1	2.00	< 2.00	
	1,3-Dichloropropane	142-28-9	2.00	< 2.00	
	1,4-Dichlorobenzene	106-46-7	2.00	< 2.00	
	1,4-Dioxane	123-91-1	40.0	< 40.0	
	2,2-Dichloropropane	594-20-7	2.00	< 2.00	
	2-Butanone	78-93-3	10.0	< 10.0	
	2-Chloroethyl vinyl ether	110-75-8	5.00	< 5.00	
	2-Chlorotoluene	95-49-8	2.00	< 2.00	
	2-Hexanone	591-78-6	5.00	< 5.00	
	2-Nitropropane	79-46-9	2.00	< 2.00	

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**ANALYTICAL** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005B

**Client Sample ID:** 4992290 J.R. @ 1700 S.

**Collection Date:** 11/30/2010 1430h **Analyzed:** 12/2/2010 0944h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	4-Chlorotoluene	106-43-4	2.00	< 2.00	
0.122	4-Isopropyltoluene	99-87-6	2.00	< 2.00	
	4-Methyl-2-pentanone	108-10-1	5.00	< 5.00	
	Acetone	67-64-1	10.0	< 10.0	
(801) 263-8686	Acetonitrile	75-05-8	5.00	< 5.00	
Toll Free (888) 263-8686	Acrolein	107-02-8	5.00	< 5.00	
Fax (801)263-8687	Acrylonitrile	107-13-1	10.0	< 10.0	
email: awal@awal-labs.com	Allyl chloride	107-05-1	5.00	< 5.00	
	Benzene	71-43-2	2.00	< 2.00	
Kyle F. Gross	Benzyl chloride	100-44-7	5.00	< 5.00	
·	Bis(2-chloroisopropyl) ether	108-60-1	5.00	< 5.00	
Laboratory Director	Bromobenzene	108-86-1	2.00	< 2.00	
	Bromochloromethane	74-97-5	2.00	< 2.00	
Jose Rocha	Bromodichloromethane	75-27-4	2.00	< 2.00	
QA Officer	Bromoform	75-25-2	2.00	< 2.00	
	Bromomethane	74-83-9	5.00	< 5.00	
	Butyl acetate	123-86-4	5.00	< 5.00	
	Carbon disulfide	75-15-0	2.00	< 2.00	
	Carbon tetrachloride	56-23-5	2.00	< 2.00	
	Chlorobenzene	108-90-7	2.00	< 2.00	
	Chloroethane	75-00-3	2.00	< 2.00	
	Chloroform	67-66-3	2.00	< 2.00	
	Chloromethane	74-87-3	5.00	< 5.00	
	Chloroprene	126-99-8	2.00	< 2.00	
	cis-1,2-Dichloroethene	156-59-2	2.00	< 2.00	
	cis-1,3-Dichloropropene	10061-01-5	2.00	< 2.00	
	Cyclohexane	110-82-7	2.00	< 2.00	
	Cyclohexanone	108-94-1	50.0	< 50.0	
	Dibromochloromethane	124-48-1	2.00	< 2.00	

Report Date: 12/6/2010 Page 73 of 162



**ANALYTICAL** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005B

Client Sample ID: 4992290 J.R. @ 1700 S.

**Collection Date:** 11/30/2010 1430h **Analyzed:** 12/2/2010 0944h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Dibromomethane	74-95-3	2.00	< 2.00	
0.112	Dichlorodifluoromethane	75-71-8	2.00	< 2.00	
	Ethyl acetate	141-78-6	10.0	< 10.0	
	Ethyl ether	60-29-7	10.0	< 10.0	
(801) 263-8686	Ethyl methacrylate	97-63-2	2.00	< 2.00	
Toll Free (888) 263-8686	Ethylbenzene	100-41-4	2.00	< 2.00	
Fax (801)263-8687	Hexachlorobutadiene	87-68-3	2.00	< 2.00	
email: awal@awal-labs.com	Iodomethane	74-88-4	5.00	< 5.00	
	Isobutyl alcohol	78-83-1	100	< 100	
Kyle F. Gross	Isopropyl acetate	108-21-4	2.00	< 2.00	
•	Isopropyl alcohol	67-63-0	25.0	< 25.0	
Laboratory Director	Isopropylbenzene	98-82-8	2.00	< 2.00	
	Isopropyltoluene	99-87-6	2.00	< 2.00	
Jose Rocha	m,p-Xylene	179601-23-1	2.00	< 2.00	
QA Officer	Methacrylonitrile	126-98-7	5.00	< 5.00	
	Methyl Acetate	79-20-9	5.00	< 5.00	
	Methyl methacrylate	80-62-6	5.00	< 5.00	
	Methyl tert-butyl ether	1634-04-4	2.00	< 2.00	
	Methylcyclohexane	108-87-2	2.00	< 2.00	
	Methylene chloride	75-09-2	2.00	< 2.00	
	n-Amyl acetate	628-63-7	2.00	< 2.00	
	n-Butyl alcohol	71-36-3	25.0	< 25.0	
	n-Butylbenzene	104-51-8	2.00	< 2.00	
	n-Hexane	110-54-3	2.00	< 2.00	
	n-Octane	111-65-9	2.00	< 2.00	
	n-Propylbenzene	103-65-1	2.00	< 2.00	
	Naphthalene	91-20-3	2.00	< 2.00	
	o-Xylene	95-47-6	2.00	< 2.00	
	Pentachloroethane	76-01-7	2.00	< 2.00	

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**ANALYTICAL** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-005B

**Client Sample ID:** 4992290 J.R. @ 1700 S.

**Collection Date:** 11/30/2010 1430h **Analyzed:** 12/2/2010 0944h

**Received Date:** 11/30/2010 1538h

**Method Used:** EPA624

LABORATORIES

Analytical Results VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Propionitrile	107-12-0	25.0	< 25.0	
	Propyl acetate	109-60-4	2.00	< 2.00	
	sec-Butylbenzene	135-98-8	2.00	< 2.00	
	Styrene	100-42-5	2.00	< 2.00	
(801) 263-8686	tert-Butyl alcohol	76-65-0	20.0	< 20.0	
Toll Free (888) 263-8686	tert-Butylbenzene	98-06-6	2.00	< 2.00	
Fax (801)263-8687	Tetrachloroethene	127-18-4	2.00	< 2.00	
email: awal@awal-labs.com	Tetrahydrofuran	109-99-9	2.00	< 2.00	
	Toluene	108-88-3	2.00	< 2.00	
Kyle F. Gross	trans-1,2-Dichloroethene	156-60-5	2.00	< 2.00	
Laboratory Director	trans-1,3-Dichloropropene	10061-02-6	2.00	< 2.00	
Laboratory Director	trans-1,4-Dichloro-2-butene	110-57-6	2.00	< 2.00	
Iona Danka	Trichloroethene	79-01-6	2.00	< 2.00	
Jose Rocha	Trichlorofluoromethane	75-69-4	2.00	< 2.00	
QA Officer	Vinyl acetate	108-05-4	5.00	< 5.00	
	Vinyl chloride	75-01-4	1.00	< 1.00	
	Xylenes, Total	1330-20-7	2.00	< 2.00	
	TPH C11-C15 (DRO)		20.0	< 20.0	
	TPH C6-C10 (GRO)		20.0	< 20.0	
	Surr: 1,2-Dichloroethane-d4	17060-07-0	77-144	111	
	Surr: 4-Bromofluorobenzene	460-00-4	80-123	105	
	Surr: Dibromofluoromethane	1868-53-7	80-124	103	
	Surr: Toluene-d8	2037-26-5	80-125	97.9	



**ANALYTICAL** 

**LABORATORIES** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006B

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

**Collection Date:** 11/30/2010 1315h **Analyzed:** 12/2/2010 0809h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

**Analytical Results** 

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	1,1,1,2-Tetrachloroethane	630-20-6	2.00	< 2.00	
	1,1,1-Trichloroethane	71-55-6	2.00	< 2.00	
	1,1,2,2-Tetrachloroethane	79-34-5	2.00	< 2.00	
	1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	2.00	< 2.00	
(801) 263-8686	1,1,2-Trichloroethane	79-00-5	2.00	< 2.00	
Toll Free (888) 263-8686	1,1-Dichloropropene	563-58-6	2.00	< 2.00	
Fax (801)263-8687	1,1-Dichloroethane	75-34-3	2.00	< 2.00	
email: awal@awal-labs.com	1,1-Dichloroethene	75-35-4	2.00	< 2.00	
	1,2,3-Trichlorobenzene	87-61-6	2.00	< 2.00	
Kyle F. Gross	1,2,3-Trichloropropane	96-18-4	2.00	< 2.00	
Laboratory Director	1,2,3-Trimethylbenzene	526-73-8	2.00	< 2.00	
Laboratory Director	1,2,4-Trichlorobenzene	120-82-1	2.00	< 2.00	
I D 1	1,2,4-Trimethylbenzene	95-63-6	2.00	< 2.00	
Jose Rocha	1,2-Dibromo-3-chloropropane	96-12-8	2.00	< 2.00	
QA Officer	1,2-Dibromoethane	106-93-4	2.00	< 2.00	
	1,2-Dichlorobenzene	95-50-1	2.00	< 2.00	
	1,2-Dichloroethane	107-06-2	2.00	< 2.00	
	1,2-Dichloropropane	78-87-5	2.00	< 2.00	
	1,3,5-Trimethylbenzene	108-67-8	2.00	< 2.00	
	1,3-Dichlorobenzene	541-73-1	2.00	< 2.00	
	1,3-Dichloropropane	142-28-9	2.00	< 2.00	
	1,4-Dichlorobenzene	106-46-7	2.00	< 2.00	
	1,4-Dioxane	123-91-1	40.0	< 40.0	
	2,2-Dichloropropane	594-20-7	2.00	< 2.00	
	2-Butanone	78-93-3	10.0	< 10.0	
	2-Chloroethyl vinyl ether	110-75-8	5.00	< 5.00	1
	2-Chlorotoluene	95-49-8	2.00	< 2.00	
	2-Hexanone	591-78-6	5.00	< 5.00	
	2-Nitropropane	79-46-9	2.00	< 2.00	

Report Date: 12/6/2010 Page 76 of 162



**ANALYTICAL** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006B

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

**Collection Date:** 11/30/2010 1315h **Analyzed:** 12/2/2010 0809h

**Received Date:** 11/30/2010 1538h

**Method Used:** EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	4-Chlorotoluene	106-43-4	2.00	< 2.00	
0.220	4-Isopropyltoluene	99-87-6	2.00	< 2.00	
	4-Methyl-2-pentanone	108-10-1	5.00	< 5.00	
	Acetone	67-64-1	10.0	< 10.0	
(801) 263-8686	Acetonitrile	75-05-8	5.00	< 5.00	
Toll Free (888) 263-8686	Acrolein	107-02-8	5.00	< 5.00	
Fax (801)263-8687	Acrylonitrile	107-13-1	10.0	< 10.0	
email: awal@awal-labs.com	Allyl chloride	107-05-1	5.00	< 5.00	
	Benzene	71-43-2	2.00	< 2.00	
Kyle F. Gross	Benzyl chloride	100-44-7	5.00	< 5.00	
Laboratory Director	Bis(2-chloroisopropyl) ether	108-60-1	5.00	< 5.00	
Laboratory Director	Bromobenzene	108-86-1	2.00	< 2.00	
T D 1	Bromochloromethane	74-97-5	2.00	< 2.00	
Jose Rocha	Bromodichloromethane	75-27-4	2.00	< 2.00	
QA Officer	Bromoform	75-25-2	2.00	< 2.00	
	Bromomethane	74-83-9	5.00	< 5.00	
	Butyl acetate	123-86-4	5.00	< 5.00	
	Carbon disulfide	75-15-0	2.00	< 2.00	
	Carbon tetrachloride	56-23-5	2.00	< 2.00	
	Chlorobenzene	108-90-7	2.00	< 2.00	
	Chloroethane	75-00-3	2.00	< 2.00	
	Chloroform	67-66-3	2.00	< 2.00	
	Chloromethane	74-87-3	5.00	< 5.00	
	Chloroprene	126-99-8	2.00	< 2.00	
	cis-1,2-Dichloroethene	156-59-2	2.00	< 2.00	
	cis-1,3-Dichloropropene	10061-01-5	2.00	< 2.00	
	Cyclohexane	110-82-7	2.00	< 2.00	
	Cyclohexanone	108-94-1	50.0	< 50.0	
	Dibromochloromethane	124-48-1	2.00	< 2.00	

Report Date: 12/6/2010 Page 77 of 162



**ANALYTICAL** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006B

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

**Collection Date:** 11/30/2010 1315h **Analyzed:** 12/2/2010 0809h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Dibromomethane	74-95-3	2.00	< 2.00	
	Dichlorodifluoromethane	75-71-8	2.00	< 2.00	
	Ethyl acetate	141-78-6	10.0	< 10.0	
	Ethyl ether	60-29-7	10.0	< 10.0	
(801) 263-8686	Ethyl methacrylate	97-63-2	2.00	< 2.00	
Toll Free (888) 263-8686	Ethylbenzene	100-41-4	2.00	< 2.00	
Fax (801)263-8687	Hexachlorobutadiene	87-68-3	2.00	< 2.00	
email: awal@awal-labs.com	Iodomethane	74-88-4	5.00	< 5.00	
	Isobutyl alcohol	78-83-1	100	< 100	
Kyle F. Gross	Isopropyl acetate	108-21-4	2.00	< 2.00	
Laboratory Director	Isopropyl alcohol	67-63-0	25.0	< 25.0	
Laboratory Director	Isopropylbenzene	98-82-8	2.00	< 2.00	
r n i	Isopropyltoluene	99-87-6	2.00	< 2.00	
Jose Rocha	m,p-Xylene	179601-23-1	2.00	< 2.00	
QA Officer	Methacrylonitrile	126-98-7	5.00	< 5.00	
	Methyl Acetate	79-20-9	5.00	< 5.00	
	Methyl methacrylate	80-62-6	5.00	< 5.00	
	Methyl tert-butyl ether	1634-04-4	2.00	< 2.00	
	Methylcyclohexane	108-87-2	2.00	< 2.00	
	Methylene chloride	75-09-2	2.00	< 2.00	
	n-Amyl acetate	628-63-7	2.00	< 2.00	
	n-Butyl alcohol	71-36-3	25.0	< 25.0	
	n-Butylbenzene	104-51-8	2.00	< 2.00	
	n-Hexane	110-54-3	2.00	< 2.00	@
	n-Octane	111-65-9	2.00	< 2.00	
	n-Propylbenzene	103-65-1	2.00	< 2.00	
	Naphthalene	91-20-3	2.00	< 2.00	
	o-Xylene	95-47-6	2.00	< 2.00	
	Pentachloroethane	76-01-7	2.00	< 2.00	

Report Date: 12/6/2010 Page 78 of 162



**ANALYTICAL** 

**WEST** 

# **ORGANIC ANALYTICAL REPORT**

Client: Utah DEQ DERR Contact: Jim Harris

**Project:** Red Butte **Lab Sample ID:** 1011529-006B

**Client Sample ID:** 4992083 Red Butte @ 1100 E.

**Collection Date:** 11/30/2010 1315h **Analyzed:** 12/2/2010 0809h

**Received Date:** 11/30/2010 1538h

Method Used: EPA624

LABORATORIES
Analytical Results

VOAs Custom List by GC/MS Method 624

463 West 3600 South Salt Lake City, UT	Units: µg/L Dilution Factor: 1 Compound	CAS Number	Reporting Limit	Analytical Result	Qual
84115	Propionitrile	107-12-0	25.0	< 25.0	
	Propyl acetate	109-60-4	2.00	< 2.00	
	sec-Butylbenzene	135-98-8	2.00	< 2.00	
	Styrene	100-42-5	2.00	< 2.00	
(801) 263-8686	tert-Butyl alcohol	76-65-0	20.0	< 20.0	
Toll Free (888) 263-8686	tert-Butylbenzene	98-06-6	2.00	< 2.00	
Fax (801)263-8687	Tetrachloroethene	127-18-4	2.00	< 2.00	
email: awal@awal-labs.com	Tetrahydrofuran	109-99-9	2.00	< 2.00	
	Toluene	108-88-3	2.00	< 2.00	
Kyle F. Gross	trans-1,2-Dichloroethene	156-60-5	2.00	< 2.00	
Laboratory Director	trans-1,3-Dichloropropene	10061-02-6	2.00	< 2.00	
Laboratory Director	trans-1,4-Dichloro-2-butene	110-57-6	2.00	< 2.00	
r D 1	Trichloroethene	79-01-6	2.00	< 2.00	
Jose Rocha	Trichlorofluoromethane	75-69-4	2.00	< 2.00	
QA Officer	Vinyl acetate	108-05-4	5.00	< 5.00	
	Vinyl chloride	75-01-4	1.00	< 1.00	
	Xylenes, Total	1330-20-7	2.00	< 2.00	
	TPH C11-C15 (DRO)		20.0	< 20.0	
	TPH C6-C10 (GRO)		20.0	< 20.0	
	Surr: 1,2-Dichloroethane-d4	17060-07-0	77-144	100	
	Surr: 4-Bromofluorobenzene	460-00-4	80-123	106	
	Surr: Dibromofluoromethane	1868-53-7	80-124	100	
	Surr: Toluene-d8	2037-26-5	80-125	102	

<sup>&</sup>lt;sup>1</sup> - Matrix spike recovery indicates matrix interference. The method is in control as indicated by the LCS.

 $<sup>@-</sup>High\ RPD\ due\ to\ suspected\ sample\ non-homogeneity\ or\ matrix\ interference.$ 



463 West 3600 South

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(801) 263-8686, Toll Free (888) 263-8686, Fax (801) 263-8687 e-mail: awal@awal-labs.com, web: www.awal-labs.com Kyle F. Gross

Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

**Dept:** GC

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limit	Qual	Date Analyzed
LCS-9679	Oil Range Organics (ORO) (C28-	mg/L	SW8015D	0.149	0.1600	0	93.4	10-200			12/6/2010 0039h
LCS-9679	Surr: C36	%REC	SW8015D	1.25	1.000		125	10-200			12/6/2010 0039h
LCS-9631	Diesel Range Organics C10-C28	mg/L	SW8015D	1.84	2.000	0	91.8	48-118			11/30/2010 2326h
LCS-9631	Surr: 4-Bromofluorobenzene	%REC	SW8015D	0.294	0.4000		73.5	18-95			11/30/2010 2326h

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463 West 3600 South

Salt Lake City, UT 84115

(801) 263-8686, Toll Free (888) 263-8686, Fax (801) 263-8687 e-mail: awal@awal-labs.com, web: www.awal-labs.com Kyle F. Gross

Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529

**Project:** Red Butte

**Contact:** Jim Harris

**Dept:** GC

**QC Type:** MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9679	Oil Range Organics (ORO) (C28-	mg/L	SW8015D	< 0.0500				-				12/6/2010 0021h
MB-9679	Surr: C36	%REC	SW8015D	1.23	1.000		123	10-200				12/6/2010 0021h
MB-9631	Diesel Range Organics C10-C28	mg/L	SW8015D	< 0.500				-				11/30/2010 2305h
MB-9631	Surr: 4-Bromofluorobenzene	%REC	SW8015D	0.272	0.4000		68.0	18-95				11/30/2010 2305h

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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

**Dept:** GC **QC Type:** MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limi	Qual	Date Analyzed
1011529-006DMS	Oil Range Organics (ORO) (C28-	mg/L	SW8015D	0.178	0.1600	0	111	10-200			12/6/2010 0247h
1011529-006DMS	Surr: C36	%REC	SW8015D	1.26	1.000		126	10-200			12/6/2010 0247h
1012039-001FMS	Oil Range Organics (ORO) (C28-	mg/L	SW8015D	0.174	0.1600	0	109	10-200			12/6/2010 0418h
1012039-001FMS	Surr: C36	%REC	SW8015D	1.24	1.000		124	10-200			12/6/2010 0418h
1011526-001BMS	Diesel Range Organics C10-C28	mg/L	SW8015D	4.16	4.000	0.7980	84.0	60-161			12/1/2010 0008h
1011526-001BMS	Surr: 4-Bromofluorobenzene	%REC	SW8015D	0.322	0.8000		40.3	10-190			12/1/2010 0008h
1011529-001EMS	Diesel Range Organics C10-C28	mg/L	SW8015D	4.03	4.167	0	96.7	60-161			12/2/2010 1355h
1011529-001EMS	Surr: 4-Bromofluorobenzene	%REC	SW8015D	0.688	0.8333		82.5	10-190			12/2/2010 1355h

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## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

**Dept:** GC

QC Type: MSD

					Amount	Original				RPD		
Sample ID	Analyte	Units	Method	Result	Spiked	Amount	% REC	Limits	%RPD	Limit	Qual	Date Analyzed
1011529-006DMSD	Oil Range Organics (ORO) (C28-	mg/L	SW8015D	0.169	0.1600	0	106	10-200	5.01	30		12/6/2010 0305h
1011529-006DMSD	Surr: C36	%REC	SW8015D	1.22	1.000		122	10-200				12/6/2010 0305h
1012039-001FMSD	Oil Range Organics (ORO) (C28-	mg/L	SW8015D	0.200	0.1600	0	125	10-200	13.4	30		12/6/2010 0436h
1012039-001FMSD	Surr: C36	%REC	SW8015D	1.42	1.000		142	10-200				12/6/2010 0436h
1011526-001BMSD	Diesel Range Organics C10-C28	mg/L	SW8015D	2.92	4.000	0.7980	53.0	60-161	35.2	25	3	12/1/2010 0029h
1011526-001BMSD	Surr: 4-Bromofluorobenzene	%REC	SW8015D	0.351	0.8000		43.9	10-190				12/1/2010 0029h
1011529-001EMSD	Diesel Range Organics C10-C28	mg/L	SW8015D	4.25	4.167	0	102	60-161	5.33	25		12/2/2010 1417h
1011529-001EMSD	Surr: 4-Bromofluorobenzene	%REC	SW8015D	0.696	0.8333		83.5	10-190				12/2/2010 1417h

<sup>&</sup>lt;sup>3</sup> - Matrix spike recoveries and high RPDs indicate suspected sample non-homogeneity. The method is in control as indicated by the LCS.

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Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limit	Qual	Date Analyzed
LCS-9662	1,1'-Biphenyl	μg/L	EPA625	49.0	80.00	0	61.3	22-104			12/2/2010 2205h
LCS-9662	1,2,4,5-Tetrachlorobenzene	$\mug/L$	EPA625	41.7	80.00	0	52.1	13-102			12/2/2010 2205h
LCS-9662	1,2,4-Trichlorobenzene	$\mug/L$	EPA625	25.4	80.00	0	31.8	10-104			12/2/2010 2205h
LCS-9662	1,2-Dichlorobenzene	$\mu g/L$	EPA625	17.9	80.00	0	22.3	10-68			12/2/2010 2205h
LCS-9662	1,3,5-Trinitrobenzene	$\mu g/L$	EPA625	133	80.00	0	166	18-209			12/2/2010 2205h
LCS-9662	1,3-Dichlorobenzene	μg/L	EPA625	14.3	80.00	0	17.9	10-60			12/2/2010 2205h
LCS-9662	1,3-Dinitrobenzene	$\mug/L$	EPA625	97.7	80.00	0	122	10-165			12/2/2010 2205h
LCS-9662	1,4-Dichlorobenzene	μg/L	EPA625	15.5	80.00	0	19.3	10-118			12/2/2010 2205h
LCS-9662	1,4-Naphthoquinone	μg/L	EPA625	< 10.0	80.00	0	12.4	10-187			12/2/2010 2205h
LCS-9662	1,4-Phenylenediamine	μg/L	EPA625	35.5	80.00	0	44.4	10-80			12/2/2010 2205h
LCS-9662	1-Chloronaphthalene	μg/L	EPA625	47.3	80.00	0	59.1	13-123			12/2/2010 2205h
LCS-9662	1-Methylnaphthalene	μg/L	EPA625	45.0	80.00	0	56.3	13-105			12/2/2010 2205h
LCS-9662	1-Naphthylamine	μg/L	EPA625	64.7	80.00	0	80.8	32-256			12/2/2010 2205h
LCS-9662	2,3,4,6-Tetrachlorophenol	μg/L	EPA625	85.4	80.00	0	107	44-158			12/2/2010 2205h
LCS-9662	2,4,5-Trichlorophenol	μg/L	EPA625	59.0	80.00	0	73.7	46-142			12/2/2010 2205h
LCS-9662	2,4,6-Trichlorophenol	μg/L	EPA625	49.9	80.00	0	62.3	42-113			12/2/2010 2205h
LCS-9662	2,4-Dichlorophenol	μg/L	EPA625	46.9	80.00	0	58.6	37-102			12/2/2010 2205h
LCS-9662	2,4-Dimethylphenol	μg/L	EPA625	53.7	80.00	0	67.1	37-99			12/2/2010 2205h
LCS-9662	2,4-Dinitrophenol	μg/L	EPA625	68.9	80.00	0	86.2	10-200			12/2/2010 2205h
LCS-9662	2,4-Dinitrotoluene	μg/L	EPA625	96.5	80.00	0	121	15-209			12/2/2010 2205h
LCS-9662	2,6-Dichlorophenol	μg/L	EPA625	53.8	80.00	0	67.2	44-111			12/2/2010 2205h
LCS-9662	2,6-Dinitrotoluene	μg/L	EPA625	78.4	80.00	0	98.0	13-183			12/2/2010 2205h
LCS-9662	2-Acetylaminofluorene	μg/L	EPA625	77.8	80.00	0	97.2	40-131			12/2/2010 2205h
LCS-9662	2-Chloronaphthalene	μg/L	EPA625	43.7	80.00	0	54.7	16-103			12/2/2010 2205h
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QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**QC Type:** LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limit	Qual	Date Analyzed
LCS-9662	2-Chlorophenol	μg/L	EPA625	42.7	80.00	0	53.4	21-98			12/2/2010 2205h
LCS-9662	2-Methylnaphthalene	μg/L	EPA625	38.6	80.00	0	48.2	11-92			12/2/2010 2205h
LCS-9662	2-Methylphenol	μg/L	EPA625	44.5	80.00	0	55.6	24-92			12/2/2010 2205h
LCS-9662	2-Naphthylamine	μg/L	EPA625	59.5	80.00	0	74.4	48-193			12/2/2010 2205h
LCS-9662	2-Nitroaniline	μg/L	EPA625	76.6	80.00	0	95.7	10-216			12/2/2010 2205h
LCS-9662	2-Nitrophenol	μg/L	EPA625	63.7	80.00	0	79.6	10-173			12/2/2010 2205h
LCS-9662	2-Picoline	μg/L	EPA625	19.2	80.00	0	24.1	10-66			12/2/2010 2205h
LCS-9662	3&4-Methylphenol	μg/L	EPA625	27.8	80.00	0	34.8	10-88			12/2/2010 2205h
LCS-9662	3,3´-Dichlorobenzidine	μg/L	EPA625	73.2	80.00	0	91.4	38-131			12/2/2010 2205h
LCS-9662	3,3´-Dimethylbenzidine	μg/L	EPA625	73.3	80.00	0	91.7	10-172			12/2/2010 2205h
LCS-9662	3-Methylcholanthrene	μg/L	EPA625	77.9	80.00	0	97.3	48-206			12/2/2010 2205h
LCS-9662	3-Nitroaniline	μg/L	EPA625	74.4	80.00	0	93.0	31-172			12/2/2010 2205h
LCS-9662	4,6-Dinitro-2-methylphenol	μg/L	EPA625	102	80.00	0	128	10-190			12/2/2010 2205h
LCS-9662	4-Aminobiphenyl	μg/L	EPA625	72.5	80.00	0	90.6	10-202			12/2/2010 2205h
LCS-9662	4-Bromophenyl phenyl ether	μg/L	EPA625	61.7	80.00	0	77.2	55-136			12/2/2010 2205h
LCS-9662	4-Chloro-3-methylphenol	μg/L	EPA625	59.0	80.00	0	73.7	47-113			12/2/2010 2205h
LCS-9662	4-Chloroaniline	μg/L	EPA625	46.7	80.00	0	58.4	24-124			12/2/2010 2205h
LCS-9662	4-Chlorophenyl phenyl ether	μg/L	EPA625	60.5	80.00	0	75.6	41-119			12/2/2010 2205h
LCS-9662	4-Nitroaniline	μg/L	EPA625	70.6	80.00	0	88.3	27-159			12/2/2010 2205h
LCS-9662	4-Nitrophenol	μg/L	EPA625	115	80.00	0	143	10-157			12/2/2010 2205h
LCS-9662	5-Nitro-o-toluidine	μg/L	EPA625	87.9	80.00	0	110	60-168			12/2/2010 2205h
LCS-9662	7,12-Dimethylbenz(a)anthracene	μg/L	EPA625	101	80.00	0	126	42-277			12/2/2010 2205h
LCS-9662	a,a-Dimethylphenethylamine	μg/L	EPA625	40.5	80.00	0	50.7	10-160			12/2/2010 2205h
LCS-9662	Acenaphthene	μg/L	EPA625	55.7	80.00	0	69.6	29-112			12/2/2010 2205h
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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	% RPD	RPD Limit	Qual	Date Analyzed
LCS-9662	Acenaphthylene	μg/L	EPA625	56.5	80.00	0	70.6	31-116				12/2/2010 2205h
LCS-9662	Acetophenone	$\mu g/L$	EPA625	44.3	80.00	0	55.3	10-105				12/2/2010 2205h
LCS-9662	alpha-Terpineol	$\mug/L$	EPA625	46.8	80.00	0	58.5	14-98				12/2/2010 2205h
LCS-9662	Aniline	$\mu g/L$	EPA625	29.0	80.00	0	36.3	10-94				12/2/2010 2205h
LCS-9662	Anthracene	$\mug/L$	EPA625	74.8	80.00	0	93.5	64-145				12/2/2010 2205h
LCS-9662	Aramite	$\mu g/L$	EPA625	90.4	80.00	0	113	46-162				12/2/2010 2205h
LCS-9662	Azobenzene	$\mu g/L$	EPA625	56.2	80.00	0	70.3	32-117				12/2/2010 2205h
LCS-9662	Benz(a)anthracene	$\mu g/L$	EPA625	72.0	80.00	0	90.1	50-134				12/2/2010 2205h
LCS-9662	Benzidine	$\mu g/L$	EPA625	62.4	80.00	0	78.0	10-211				12/2/2010 2205h
LCS-9662	Benzo(a)pyrene	$\mu g/L$	EPA625	79.7	80.00	0	99.7	39-152				12/2/2010 2205h
LCS-9662	Benzo(b)fluoranthene	$\mu g/L$	EPA625	71.3	80.00	0	89.2	46-256				12/2/2010 2205h
LCS-9662	Benzo(g,h,i)perylene	$\mu g/L$	EPA625	67.5	80.00	0	84.4	21-288				12/2/2010 2205h
LCS-9662	Benzo(k)fluoranthene	$\mu g/L$	EPA625	97.6	80.00	0	122	10-254				12/2/2010 2205h
LCS-9662	Benzoic acid	$\mu g/L$	EPA625	< 20.0	80.00	0	19.0	10-64				12/2/2010 2205h
LCS-9662	Benzyl alcohol	$\mu g/L$	EPA625	23.0	80.00	0	28.8	10-88				12/2/2010 2205h
LCS-9662	Bis(2-chloroethoxy)methane	$\mu g/L$	EPA625	33.1	80.00	0	41.4	10-98				12/2/2010 2205h
LCS-9662	Bis(2-chloroethyl) ether	$\mu g/L$	EPA625	31.6	80.00	0	39.4	10-99				12/2/2010 2205h
LCS-9662	Bis(2-chloroisopropyl) ether	$\mu g/L$	EPA625	30.8	80.00	0	38.6	10-92				12/2/2010 2205h
LCS-9662	Bis(2-ethylhexyl) phthalate	$\mu g/L$	EPA625	99.0	80.00	0	124	10-233				12/2/2010 2205h
LCS-9662	bis(2-ethylhexyl)adipate	$\mu g/L$	EPA625	90.0	80.00	0	112	10-200				12/2/2010 2205h
LCS-9662	Butyl benzyl phthalate	$\mu g/L$	EPA625	85.9	80.00	0	107	10-178				12/2/2010 2205h
LCS-9662	Carbazole	$\mu g/L$	EPA625	73.2	80.00	0	91.5	61-140				12/2/2010 2205h
LCS-9662	Chlorobenzilate	μg/L	EPA625	77.9	80.00	0	97.4	10-218				12/2/2010 2205h
LCS-9662	Chrysene	μg/L	EPA625	77.7	80.00	0	97.2	54-130				12/2/2010 2205h
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QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RP %RPD Lin	Date Analyzed
LCS-9662	Diallate (cis or trans)	μg/L	EPA625	60.8	80.00	0	76.1	41-132		12/2/2010 2205h
LCS-9662	Dibenz(a,h)anthracene	$\mug/L$	EPA625	70.6	80.00	0	88.2	27-199		12/2/2010 2205h
LCS-9662	Dibenzofuran	$\mug/L$	EPA625	57.9	80.00	0	72.4	38-117		12/2/2010 2205h
LCS-9662	Diethyl phthalate	$\mu g/L$	EPA625	56.8	80.00	0	71.0	20-128		12/2/2010 2205h
LCS-9662	Dimethoate	$\mu g/L$	EPA625	< 10.0	80.00	0	12.2	10-60		12/2/2010 2205h
LCS-9662	Dimethyl phthalate	$\mu g/L$	EPA625	36.3	80.00	0	45.3	10-105		12/2/2010 2205h
LCS-9662	Dimethylaminoazobenzene	$\mu g/L$	EPA625	84.0	80.00	0	105	30-151		12/2/2010 2205h
LCS-9662	Di-n-butyl phthalate	$\mu g/L$	EPA625	75.2	80.00	0	94.0	45-131		12/2/2010 2205h
LCS-9662	Di-n-octyl phthalate	μg/L	EPA625	136	80.00	0	170	10-252		12/2/2010 2205h
LCS-9662	Dinoseb	$\mug/L$	EPA625	113	80.00	0	142	10-186		12/2/2010 2205h
LCS-9662	Diphenylamine	μg/L	EPA625	67.5	80.00	0	84.4	46-128		12/2/2010 2205h
LCS-9662	Disulfoton	μg/L	EPA625	51.9	80.00	0	64.9	10-124		12/2/2010 2205h
LCS-9662	Ethyl methanesulfonate	$\mu g/L$	EPA625	39.4	80.00	0	49.2	10-105		12/2/2010 2205h
LCS-9662	Famphur	μg/L	EPA625	140	80.00	0	175	10-298		12/2/2010 2205h
LCS-9662	Fluoranthene	$\mu g/L$	EPA625	74.0	80.00	0	92.5	61-138		12/2/2010 2205h
LCS-9662	Fluorene	μg/L	EPA625	60.3	80.00	0	75.4	45-116		12/2/2010 2205h
LCS-9662	Hexachlorobenzene	μg/L	EPA625	64.4	80.00	0	80.5	55-135		12/2/2010 2205h
LCS-9662	Hexachlorobutadiene	μg/L	EPA625	18.5	80.00	0	23.2	10-79		12/2/2010 2205h
LCS-9662	Hexachlorocyclopentadiene	μg/L	EPA625	11.1	80.00	0	13.8	10-104		12/2/2010 2205h
LCS-9662	Hexachloroethane	μg/L	EPA625	13.6	80.00	0	17.0	10-58		12/2/2010 2205h
LCS-9662	Hexachlorophene	μg/L	EPA625	82.4	80.00	0	103	10-242		12/2/2010 2205h
LCS-9662	Hexachloropropene	μg/L	EPA625	14.4	80.00	0	18.0	10-79		12/2/2010 2205h
LCS-9662	Indene	μg/L	EPA625	23.9	80.00	0	29.8	10-71		12/2/2010 2205h
LCS-9662	Indeno(1,2,3-cd)pyrene	μg/L	EPA625	69.9	80.00	0	87.3	29-208		12/2/2010 2205h

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## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD Limit	Qual	Date Analyzed
LCS-9662	Isodrin	μg/L	EPA625	70.1	80.00	0	87.6	58-140			12/2/2010 2205h
LCS-9662	Isophorone	$\mu g/L$	EPA625	45.6	80.00	0	57.0	10-105			12/2/2010 2205h
LCS-9662	Isosafrole	$\mu g/L$	EPA625	44.3	80.00	0	55.4	12-108			12/2/2010 2205h
LCS-9662	Kepone	$\mu g/L$	EPA625	214	80.00	0	268	10-330			12/2/2010 2205h
LCS-9662	Methapyrilene	$\mu g/L$	EPA625	85.9	80.00	0	107	17-151			12/2/2010 2205h
LCS-9662	Methyl methanesulfonate	$\mu g/L$	EPA625	25.4	80.00	0	31.7	10-107			12/2/2010 2205h
LCS-9662	Naphthalene	$\mu g/L$	EPA625	33.2	80.00	0	41.5	10-79			12/2/2010 2205h
LCS-9662	n-Decane	$\mu g/L$	EPA625	< 10.0	80.00	0	10.5	10-45			12/2/2010 2205h
LCS-9662	Nitrobenzene	$\mu g/L$	EPA625	50.4	80.00	0	63.0	10-104			12/2/2010 2205h
LCS-9662	Nitroquinoline-1-oxide	$\mu g/L$	EPA625	29.4	80.00	0	36.8	10-165			12/2/2010 2205h
LCS-9662	N-Nitrosodiethylamine	μg/L	EPA625	36.2	80.00	0	45.2	10-96			12/2/2010 2205h
LCS-9662	N-Nitrosodimethylamine	μg/L	EPA625	16.3	80.00	0	20.4	10-55			12/2/2010 2205h
LCS-9662	N-Nitrosodi-n-butylamine	$\mu g/L$	EPA625	52.5	80.00	0	65.6	21-104			12/2/2010 2205h
LCS-9662	N-Nitrosodiphenylamine	μg/L	EPA625	66.1	80.00	0	82.6	45-126			12/2/2010 2205h
LCS-9662	N-Nitrosodi-n-propylamine	μg/L	EPA625	40.9	80.00	0	51.2	10-103			12/2/2010 2205h
LCS-9662	N-Nitrosomethylethylamine	μg/L	EPA625	28.9	80.00	0	36.2	10-84			12/2/2010 2205h
LCS-9662	N-Nitrosomorpholine	$\mu g/L$	EPA625	43.4	80.00	0	54.2	15-107			12/2/2010 2205h
LCS-9662	N-Nitrosopiperidine	μg/L	EPA625	50.5	80.00	0	63.1	10-108			12/2/2010 2205h
LCS-9662	N-Nitrosopyrrolidine	$\mu g/L$	EPA625	50.8	80.00	0	63.5	19-115			12/2/2010 2205h
LCS-9662	n-Octadecane	μg/L	EPA625	55.0	80.00	0	68.8	27-134			12/2/2010 2205h
LCS-9662	O,O,O-Triethyl phosphorothioate	μg/L	EPA625	48.4	80.00	0	60.5	10-129			12/2/2010 2205h
LCS-9662	o-Toluidine	μg/L	EPA625	84.4	80.00	0	106	21-225			12/2/2010 2205h
LCS-9662	Parathion	μg/L	EPA625	106	80.00	0	132	10-165			12/2/2010 2205h
LCS-9662	Methyl parathion	μg/L	EPA625	98.2	80.00	0	123	10-165			12/2/2010 2205h

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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
LCS-9662	Pentachlorobenzene	μg/L	EPA625	56.8	80.00	0	71.0	41-118				12/2/2010 2205h
LCS-9662	Pentachloronitrobenzene	$\mug/L$	EPA625	75.6	80.00	0	94.5	35-185				12/2/2010 2205h
LCS-9662	Pentachlorophenol	$\mug/L$	EPA625	65.4	80.00	0	81.8	14-144				12/2/2010 2205h
LCS-9662	Phenacetin	$\mug/L$	EPA625	73.0	80.00	0	91.2	42-163				12/2/2010 2205h
LCS-9662	Phenanthrene	$\mug/L$	EPA625	76.0	80.00	0	95.0	25-135				12/2/2010 2205h
LCS-9662	Phenol	$\mug/L$	EPA625	26.5	80.00	0	33.1	10-53				12/2/2010 2205h
LCS-9662	Phorate	$\mug/L$	EPA625	73.0	80.00	0	91.2	36-158				12/2/2010 2205h
LCS-9662	Pronamide	$\mug/L$	EPA625	78.2	80.00	0	97.7	33-137				12/2/2010 2205h
LCS-9662	Pyrene	$\mug/L$	EPA625	85.1	80.00	0	106	42-145				12/2/2010 2205h
LCS-9662	Pyridine	$\mug/L$	EPA625	10.3	80.00	0	12.9	10-37				12/2/2010 2205h
LCS-9662	Quinoline	$\mug/L$	EPA625	53.9	80.00	0	67.4	13-111				12/2/2010 2205h
LCS-9662	Safrole	$\mug\!/L$	EPA625	49.1	80.00	0	61.4	24-111				12/2/2010 2205h
LCS-9662	Tetraethyl dithiopyrophosphate	$\mug\!/L$	EPA625	65.9	80.00	0	82.4	52-138				12/2/2010 2205h
LCS-9662	Thionazin	$\mug\!/L$	EPA625	62.8	80.00	0	78.4	45-123				12/2/2010 2205h
LCS-9662	Surr: 2,4,6-Tribromophenol	%REC	EPA625	66.9	80.00		83.7	64-130				12/2/2010 2205h
LCS-9662	Surr: 2-Fluorobiphenyl	%REC	EPA625	47.6	40.00		119	32-128				12/2/2010 2205h
LCS-9662	Surr: 2-Fluorophenol	%REC	EPA625	25.2	80.00		31.4	10-121				12/2/2010 2205h
LCS-9662	Surr: Nitrobenzene-d5	%REC	EPA625	47.1	40.00		118	17-133				12/2/2010 2205h
LCS-9662	Surr: Phenol-d6	%REC	EPA625	18.7	80.00		23.4	10-124				12/2/2010 2205h
LCS-9662	Surr: Terphenyl-d14	%REC	EPA625	86.5	40.00		216	51-221				12/2/2010 2205h
LCS-9662	Acenaphthene	$\mug/L$	SW8270D	55.7	80.00	0	69.6	20-116				12/2/2010 2205h
LCS-9662	Benzo(a)pyrene	$\mug/L$	SW8270D	79.7	80.00	0	99.7	39-152				12/2/2010 2205h
LCS-9662	Pyrene	$\mug/L$	SW8270D	85.1	80.00	0	106	37-138				12/2/2010 2205h
LCS-9662	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	66.9	80.00		83.7	10-165				12/2/2010 2205h
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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**Dept:** MSS **QC Type:** LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	RPD Limit	Qual	Date Analyzed
LCS-9662	Surr: 2-Fluorobiphenyl	%REC	SW8270D	47.6	40.00		119	32-128			12/2/2010 2205h
LCS-9662	Surr: 2-Fluorophenol	%REC	SW8270D	25.2	80.00		31.4	10-121			12/2/2010 2205h
LCS-9662	Surr: Nitrobenzene-d5	%REC	SW8270D	47.1	40.00		118	17-133			12/2/2010 2205h
LCS-9662	Surr: Phenol-d6	%REC	SW8270D	18.7	80.00		23.4	10-124			12/2/2010 2205h
LCS-9662	Surr: Terphenyl-d14	%REC	SW8270D	86.5	40.00		216	51-221			12/2/2010 2205h

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QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9662	1,1´-Biphenyl	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,2,4,5-Tetrachlorobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,2,4-Trichlorobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,2-Dichlorobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,3,5-Trinitrobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,3-Dichlorobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,3-Dinitrobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,4-Dichlorobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,4-Naphthoquinone	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1,4-Phenylenediamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1-Chloronaphthalene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1-Methylnaphthalene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	1-Naphthylamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,3,4,6-Tetrachlorophenol	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,4,5-Trichlorophenol	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,4,6-Trichlorophenol	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,4-Dichlorophenol	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,4-Dimethylphenol	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,4-Dinitrophenol	$\mug/L$	EPA625	< 20.0				-				12/2/2010 2137h
MB-9662	2,4-Dinitrotoluene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,6-Dichlorophenol	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2,6-Dinitrotoluene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Acetylaminofluorene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Chloronaphthalene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
										Repor	t Date: 12	2/6/2010 Page 91 of 162

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Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9662	2-Chlorophenol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Methylnaphthalene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Methylphenol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Naphthylamine	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Nitroaniline	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Nitrophenol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	2-Picoline	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	3&4-Methylphenol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	3,3´-Dichlorobenzidine	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	3,3´-Dimethylbenzidine	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	3-Methylcholanthrene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	3-Nitroaniline	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4,6-Dinitro-2-methylphenol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4-Aminobiphenyl	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4-Bromophenyl phenyl ether	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4-Chloro-3-methylphenol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4-Chloroaniline	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4-Chlorophenyl phenyl ether	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4-Nitroaniline	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	4-Nitrophenol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	5-Nitro-o-toluidine	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	7,12-Dimethylbenz(a)anthracene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	a,a-Dimethylphenethylamine	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Acenaphthene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
	-									Repor	t Data: 12	2/6/2010 Page 92 of 16

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## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9662	Acenaphthylene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Acetophenone	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	alpha-Terpineol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Aniline	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Anthracene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Aramite	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Azobenzene	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Benz(a)anthracene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Benzidine	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Benzo(a)pyrene	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Benzo(b)fluoranthene	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Benzo(g,h,i)perylene	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Benzo(k)fluoranthene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Benzoic acid	μg/L	EPA625	< 20.0				-				12/2/2010 2137h
MB-9662	Benzyl alcohol	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Bis(2-chloroethoxy)methane	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Bis(2-chloroethyl) ether	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Bis(2-chloroisopropyl) ether	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Bis(2-ethylhexyl) phthalate	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	bis(2-ethylhexyl)adipate	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Butyl benzyl phthalate	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Carbazole	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Chlorobenzilate	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Chrysene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
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## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9662	Diallate (cis or trans)	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Dibenz(a,h)anthracene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Dibenzofuran	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Diethyl phthalate	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Dimethoate	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Dimethyl phthalate	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Dimethylaminoazobenzene	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Di-n-butyl phthalate	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Di-n-octyl phthalate	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Dinoseb	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Diphenylamine	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Disulfoton	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Ethyl methanesulfonate	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Famphur	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Fluoranthene	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Fluorene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Hexachlorobenzene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Hexachlorobutadiene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Hexachlorocyclopentadiene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Hexachloroethane	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Hexachlorophene	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Hexachloropropene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Indene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Indeno(1,2,3-cd)pyrene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**Dept:** MSSV **QC Type:** MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9662	Isodrin	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Isophorone	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Isosafrole	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Kepone	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Methapyrilene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Methyl methanesulfonate	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Naphthalene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	n-Decane	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Nitrobenzene	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Nitroquinoline-1-oxide	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosodiethylamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosodimethylamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosodi-n-butylamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosodiphenylamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosodi-n-propylamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosomethylethylamine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosomorpholine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosopiperidine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	N-Nitrosopyrrolidine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	n-Octadecane	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	O,O,O-Triethyl phosphorothioate	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	o-Toluidine	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Parathion	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Methyl parathion	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
										D	4 D-4 10	0/6/2010 Dogg 05 of 162

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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**QC Type:** MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9662	Pentachlorobenzene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Pentachloronitrobenzene	μg/L	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Pentachlorophenol	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Phenacetin	$\mu g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Phenanthrene	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Phenol	$\mug/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Phorate	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Pronamide	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Pyrene	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Pyridine	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Quinoline	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Safrole	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Tetraethyl dithiopyrophosphate	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Thionazin	$\mu  g/L$	EPA625	< 10.0				-				12/2/2010 2137h
MB-9662	Surr: 2,4,6-Tribromophenol	%REC	EPA625	52.3	80.00		65.4	10-165				12/2/2010 2137h
MB-9662	Surr: 2-Fluorobiphenyl	%REC	EPA625	22.3	40.00		55.7	18-113				12/2/2010 2137h
MB-9662	Surr: 2-Fluorophenol	%REC	EPA625	32.2	80.00		40.2	10-121				12/2/2010 2137h
MB-9662	Surr: Nitrobenzene-d5	%REC	EPA625	19.7	40.00		49.2	17-133				12/2/2010 2137h
MB-9662	Surr: Phenol-d6	%REC	EPA625	16.0	80.00		20.0	10-124				12/2/2010 2137h
MB-9662	Surr: Terphenyl-d14	%REC	EPA625	45.7	40.00		114	28-163				12/2/2010 2137h
MB-9662	Acenaphthene	$\mu g/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Acenaphthylene	$\mu g/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Anthracene	μg/L	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Benz(a)anthracene	μg/L	SW8270D	< 10.0				-				12/2/2010 2137h
										Papor	t Data: 1	2/6/2010 Page 96 of 162

Report Date: 12/6/2010 Page 96 of 162



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Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV
QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-9662	Benzo(a)pyrene	μg/L	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Benzo(b)fluoranthene	μg/L	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Benzo(g,h,i)perylene	$\mu g/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Benzo(k)fluoranthene	$\mu g/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	C11-C12 Aliphatic hydrocarbons	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	C11-C13 Alkyl Naphthalenes	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	C13-C16 Aliphatic hydrocarbons	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	C17-C21 Aliphatic hydrocarbons	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	C22-C35 Aliphatic hydrocarbons	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Chrysene	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Dibenz(a,h)anthracene	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Fluoranthene	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Fluorene	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Indeno(1,2,3-cd)pyrene	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Phenanthrene	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Pyrene	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Total C12-C22 PAH**	$\mug/L$	SW8270D	< 10.0				-				12/2/2010 2137h
MB-9662	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	52.3	80.00		65.4	10-165				12/2/2010 2137h
MB-9662	Surr: 2-Fluorobiphenyl	%REC	SW8270D	22.3	40.00		55.7	18-113				12/2/2010 2137h
MB-9662	Surr: 2-Fluorophenol	%REC	SW8270D	32.2	80.00		40.2	10-121				12/2/2010 2137h
MB-9662	Surr: Nitrobenzene-d5	%REC	SW8270D	19.7	40.00		49.2	17-133				12/2/2010 2137h
MB-9662	Surr: Phenol-d6	%REC	SW8270D	16.0	80.00		20.0	10-124				12/2/2010 2137h
MB-9662	Surr: Terphenyl-d14	%REC	SW8270D	45.7	40.00		114	28-163				12/2/2010 2137h

<sup>\*\* -</sup> This value is a summation of the PAH compounds listed above.

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Laboratory Director

Jose Rocha

QA Officer

### **OC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

**RPD** Amount **Original** % REC %RPD Limit Sample ID Analyte Units Method Result Spiked Amount Limits Oual Date Analyzed 1011529-006FMS 1,1'-Biphenyl μg/L EPA625 48.3 80.00 0 60.4 27-99 12/3/2010 0117h 1011529-006FMS 1,2,4,5-Tetrachlorobenzene  $\mu g/L$ EPA625 39.8 80.00 0 49.8 10-119 12/3/2010 0117h 1011529-006FMS 1,2,4-Trichlorobenzene EPA625 24.1 80.00 0 30.1 10-79 12/3/2010 0117h μg/L 1011529-006FMS 1.2-Dichlorobenzene EPA625 18.4 80.00 0 23.0 10-59 12/3/2010 0117h μg/L 1011529-006FMS 1,3,5-Trinitrobenzene  $\mu g/L$ EPA625 138 80.00 0 172 10-175 12/3/2010 0117h 1011529-006FMS EPA625 15.0 80.00 0 18.8 10-56 12/3/2010 0117h 1.3-Dichlorobenzene μg/L 1011529-006FMS 1,3-Dinitrobenzene μg/L EPA625 96.8 80.00 0 121 10-175 12/3/2010 0117h 1011529-006FMS 1,4-Dichlorobenzene EPA625 16.4 80.00 0 20.5 10-58 12/3/2010 0117h μg/L 1011529-006FMS 1,4-Naphthoquinone  $\mu g/L$ EPA625 < 10.0 80.00 0 11.6 10-177 12/3/2010 0117h 1011529-006FMS 1,4-Phenylenediamine EPA625 34.6 80.00 0 43.2 10-124 12/3/2010 0117h μg/L 1011529-006FMS 1-Chloronaphthalene EPA625 46.0 80.00 0 57.5 10-106 12/3/2010 0117h μg/L 1011529-006FMS 1-Methylnaphthalene μg/L EPA625 44.0 80.00 0 55.0 10-83 12/3/2010 0117h 1011529-006FMS 1-Naphthylamine  $\mu\,\text{g/L}$ EPA625 63.4 80.00 0 79.2 10-122 12/3/2010 0117h 1011529-006FMS 2,3,4,6-Tetrachlorophenol μg/L EPA625 76.8 80.00 0 96.0 10-157 12/3/2010 0117h 1011529-006FMS 2,4,5-Trichlorophenol  $\mu g/L$ EPA625 60.2 80.00 0 75.2 10-148 12/3/2010 0117h EPA625 50.7 80.00 0 63.4 10-136 12/3/2010 0117h 1011529-006FMS 2,4,6-Trichlorophenol μg/L 1011529-006FMS 2,4-Dichlorophenol μg/L EPA625 50.9 80.00 0 63.6 10-123 12/3/2010 0117h 1011529-006FMS EPA625 56.5 80.00 0 70.7 10-113 12/3/2010 0117h 2,4-Dimethylphenol μg/L 1011529-006FMS 2,4-Dinitrophenol μg/L EPA625 65.9 80.00 0 82.4 10-175 12/3/2010 0117h 1011529-006FMS 2,4-Dinitrotoluene EPA625 94.7 80.00 0 118 10-175 12/3/2010 0117h μg/L 1011529-006FMS 2,6-Dichlorophenol μg/L EPA625 57.5 80.00 0 71.9 10-148 12/3/2010 0117h 1011529-006FMS 2,6-Dinitrotoluene  $\mu g/L$ EPA625 79.0 80.00 0 98.8 10-175 12/3/2010 0117h 1011529-006FMS 2-Acetylaminofluorene μg/L EPA625 78.6 80.00 0 98.3 10-94 12/3/2010 0117h 1011529-006FMS 2-Chloronaphthalene  $\mu g/L$ EPA625 42.1 80.00 0 52.6 10-93 12/3/2010 0117h

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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**QC Type:** MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD Limit	Qual	Date Analyzed
1011529-006FMS	2-Chlorophenol	μg/L	EPA625	45.6	80.00	0	57.0	10-92			12/3/2010 0117h
1011529-006FMS	2-Methylnaphthalene	$\mu g/L$	EPA625	38.1	80.00	0	47.7	15-78			12/3/2010 0117h
1011529-006FMS	2-Methylphenol	$\mug/L$	EPA625	46.6	80.00	0	58.2	10-83			12/3/2010 0117h
1011529-006FMS	2-Naphthylamine	$\mug/L$	EPA625	59.8	80.00	0	74.8	10-154			12/3/2010 0117h
1011529-006FMS	2-Nitroaniline	$\mug/L$	EPA625	78.5	80.00	0	98.2	10-175			12/3/2010 0117h
1011529-006FMS	2-Nitrophenol	$\mug/L$	EPA625	68.6	80.00	0	85.8	10-175			12/3/2010 0117h
1011529-006FMS	2-Picoline	$\mug/L$	EPA625	22.4	80.00	0	28.0	10-61			12/3/2010 0117h
1011529-006FMS	3&4-Methylphenol	$\mug/L$	EPA625	29.4	80.00	0	36.8	10-80			12/3/2010 0117h
1011529-006FMS	3,3´-Dichlorobenzidine	$\mug/L$	EPA625	73.2	80.00	0	91.5	10-150			12/3/2010 0117h
1011529-006FMS	3,3´-Dimethylbenzidine	$\mug/L$	EPA625	73.3	80.00	0	91.6	10-143			12/3/2010 0117h
1011529-006FMS	3-Methylcholanthrene	$\mug/L$	EPA625	76.8	80.00	0	96.0	32-171			12/3/2010 0117h
1011529-006FMS	3-Nitroaniline	$\mug/L$	EPA625	73.6	80.00	0	92.0	10-175			12/3/2010 0117h
1011529-006FMS	4,6-Dinitro-2-methylphenol	$\mug/L$	EPA625	105	80.00	0	131	10-175			12/3/2010 0117h
1011529-006FMS	4-Aminobiphenyl	$\mug/L$	EPA625	73.7	80.00	0	92.1	10-175			12/3/2010 0117h
1011529-006FMS	4-Bromophenyl phenyl ether	$\mug/L$	EPA625	64.1	80.00	0	80.1	16-138			12/3/2010 0117h
1011529-006FMS	4-Chloro-3-methylphenol	$\mug/L$	EPA625	63.5	80.00	0	79.4	10-131			12/3/2010 0117h
1011529-006FMS	4-Chloroaniline	$\mug/L$	EPA625	50.4	80.00	0	63.0	10-98			12/3/2010 0117h
1011529-006FMS	4-Chlorophenyl phenyl ether	$\mu g/L$	EPA625	60.2	80.00	0	75.3	31-108			12/3/2010 0117h
1011529-006FMS	4-Nitroaniline	$\mug/L$	EPA625	68.7	80.00	0	85.9	10-175			12/3/2010 0117h
1011529-006FMS	4-Nitrophenol	$\mug/L$	EPA625	110	80.00	0	137	10-97		1	12/3/2010 0117h
1011529-006FMS	5-Nitro-o-toluidine	$\mug/L$	EPA625	86.4	80.00	0	108	10-175			12/3/2010 0117h
1011529-006FMS	7,12-Dimethylbenz(a)anthracene	$\mug/L$	EPA625	99.1	80.00	0	124	26-174			12/3/2010 0117h
1011529-006FMS	a,a-Dimethylphenethylamine	$\mu g/L$	EPA625	38.2	80.00	0	47.8	10-175			12/3/2010 0117h
1011529-006FMS	Acenaphthene	$\mu g/L$	EPA625	54.7	80.00	0	68.4	29-97			12/3/2010 0117h
									Papor	t Data: 1	2/6/2010 Page 99 of 16

Report Date: 12/6/2010 Page 99 of 162



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QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limit	Qual	Date Analyzed
1011529-006FMS	Acenaphthylene	μg/L	EPA625	54.6	80.00	0	68.3	37-87			12/3/2010 0117h
1011529-006FMS	Acetophenone	μg/L	EPA625	48.8	80.00	0	61.0	10-96			12/3/2010 0117h
1011529-006FMS	alpha-Terpineol	$\mu g/L$	EPA625	50.0	80.00	0	62.5	10-67			12/3/2010 0117h
1011529-006FMS	Aniline	μg/L	EPA625	33.4	80.00	0	41.8	10-71			12/3/2010 0117h
1011529-006FMS	Anthracene	$\mu g/L$	EPA625	76.0	80.00	0	95.0	53-114			12/3/2010 0117h
1011529-006FMS	Aramite	$\mu g/L$	EPA625	85.4	80.00	0	107	29-160			12/3/2010 0117h
1011529-006FMS	Azobenzene	$\mu g/L$	EPA625	56.8	80.00	0	71.0	15-114			12/3/2010 0117h
1011529-006FMS	Benz(a)anthracene	$\mu g/L$	EPA625	72.6	80.00	0	90.8	39-129			12/3/2010 0117h
1011529-006FMS	Benzidine	μg/L	EPA625	59.7	80.00	0	74.6	10-99			12/3/2010 0117h
1011529-006FMS	Benzo(a)pyrene	μg/L	EPA625	79.5	80.00	0	99.4	29-175			12/3/2010 0117h
1011529-006FMS	Benzo(b)fluoranthene	μg/L	EPA625	69.4	80.00	0	86.8	15-140			12/3/2010 0117h
1011529-006FMS	Benzo(g,h,i)perylene	μg/L	EPA625	65.0	80.00	0	81.3	10-182			12/3/2010 0117h
1011529-006FMS	Benzo(k)fluoranthene	μg/L	EPA625	97.9	80.00	0	122	21-154			12/3/2010 0117h
1011529-006FMS	Benzoic acid	μg/L	EPA625	< 20.0	80.00	0	19.4	10-71			12/3/2010 0117h
1011529-006FMS	Benzyl alcohol	μg/L	EPA625	24.1	80.00	0	30.2	10-69			12/3/2010 0117h
1011529-006FMS	Bis(2-chloroethoxy)methane	μg/L	EPA625	36.6	80.00	0	45.7	10-94			12/3/2010 0117h
1011529-006FMS	Bis(2-chloroethyl) ether	μg/L	EPA625	37.0	80.00	0	46.3	10-70			12/3/2010 0117h
1011529-006FMS	Bis(2-chloroisopropyl) ether	μg/L	EPA625	35.4	80.00	0	44.3	10-71			12/3/2010 0117h
1011529-006FMS	Bis(2-ethylhexyl) phthalate	μg/L	EPA625	90.6	80.00	0	113	10-175			12/3/2010 0117h
1011529-006FMS	bis(2-ethylhexyl)adipate	μg/L	EPA625	90.6	80.00	0	113	10-175			12/3/2010 0117h
1011529-006FMS	Butyl benzyl phthalate	μg/L	EPA625	87.8	80.00	0	110	10-175			12/3/2010 0117h
1011529-006FMS	Carbazole	μg/L	EPA625	73.8	80.00	0	92.3	10-151			12/3/2010 0117h
1011529-006FMS	Chlorobenzilate	μg/L	EPA625	75.8	80.00	0	94.8	18-175			12/3/2010 0117h
1011529-006FMS	Chrysene	μg/L	EPA625	79.5	80.00	0	99.4	38-133			12/3/2010 0117h
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QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV
QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMS	Diallate (cis or trans)	μg/L	EPA625	62.9	80.00	0	78.6	10-157				12/3/2010 0117h
1011529-006FMS	Dibenz(a,h)anthracene	$\mu g/L$	EPA625	67.9	80.00	0	84.9	13-168				12/3/2010 0117h
1011529-006FMS	Dibenzofuran	$\mu  g/L$	EPA625	56.8	80.00	0	71.0	29-103				12/3/2010 0117h
1011529-006FMS	Diethyl phthalate	$\mu  g/L$	EPA625	53.9	80.00	0	67.3	10-139				12/3/2010 0117h
1011529-006FMS	Dimethoate	$\mu  g/L$	EPA625	< 10.0	80.00	0	11.4	10-136				12/3/2010 0117h
1011529-006FMS	Dimethyl phthalate	$\mug/L$	EPA625	31.4	80.00	0	39.2	10-122				12/3/2010 0117h
1011529-006FMS	Dimethylaminoazobenzene	$\mug/L$	EPA625	85.0	80.00	0	106	34-142				12/3/2010 0117h
1011529-006FMS	Di-n-butyl phthalate	$\mu  g/L$	EPA625	75.1	80.00	0	93.8	44-124				12/3/2010 0117h
1011529-006FMS	Di-n-octyl phthalate	$\mug/L$	EPA625	134	80.00	0	167	10-175				12/3/2010 0117h
1011529-006FMS	Dinoseb	$\mu  g/L$	EPA625	119	80.00	0	149	10-175				12/3/2010 0117h
1011529-006FMS	Diphenylamine	$\mu  g/L$	EPA625	67.8	80.00	0	84.8	13-110				12/3/2010 0117h
1011529-006FMS	Disulfoton	$\mug/L$	EPA625	50.6	80.00	0	63.3	10-121				12/3/2010 0117h
1011529-006FMS	Ethyl methanesulfonate	$\mug/L$	EPA625	44.2	80.00	0	55.2	10-99				12/3/2010 0117h
1011529-006FMS	Famphur	$\mug/L$	EPA625	140	80.00	0	174	10-71			1	12/3/2010 0117h
1011529-006FMS	Fluoranthene	$\mug/L$	EPA625	74.9	80.00	0	93.6	23-135				12/3/2010 0117h
1011529-006FMS	Fluorene	$\mug/L$	EPA625	60.4	80.00	0	75.5	34-108				12/3/2010 0117h
1011529-006FMS	Hexachlorobenzene	$\mug/L$	EPA625	65.1	80.00	0	81.4	26-131				12/3/2010 0117h
1011529-006FMS	Hexachlorobutadiene	$\mug/L$	EPA625	19.3	80.00	0	24.1	10-110				12/3/2010 0117h
1011529-006FMS	Hexachlorocyclopentadiene	$\mug/L$	EPA625	11.2	80.00	0	14.0	10-45				12/3/2010 0117h
1011529-006FMS	Hexachloroethane	$\mug/L$	EPA625	14.7	80.00	0	18.3	10-58				12/3/2010 0117h
1011529-006FMS	Hexachlorophene	$\mu g/L$	EPA625	81.4	80.00	0	102	10-168				12/3/2010 0117h
1011529-006FMS	Hexachloropropene	$\mu g/L$	EPA625	15.4	80.00	0	19.2	10-72				12/3/2010 0117h
1011529-006FMS	Indene	$\mu  g/L$	EPA625	23.1	80.00	0	28.9	10-35				12/3/2010 0117h
1011529-006FMS	Indeno(1,2,3-cd)pyrene	$\mug/L$	EPA625	68.6	80.00	0	85.7	10-176				12/3/2010 0117h

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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**QC Type:** MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMS	Isodrin	μg/L	EPA625	69.6	80.00	0	87.1	15-165				12/3/2010 0117h
1011529-006FMS	Isophorone	$\mug/L$	EPA625	50.6	80.00	0	63.2	10-99				12/3/2010 0117h
1011529-006FMS	Isosafrole	$\mug/L$	EPA625	43.6	80.00	0	54.6	10-167				12/3/2010 0117h
1011529-006FMS	Kepone	$\mug/L$	EPA625	217	80.00	0	272	10-175			1	12/3/2010 0117h
1011529-006FMS	Methapyrilene	$\mug/L$	EPA625	85.0	80.00	0	106	10-149				12/3/2010 0117h
1011529-006FMS	Methyl methanesulfonate	$\mug/L$	EPA625	29.7	80.00	0	37.2	10-132				12/3/2010 0117h
1011529-006FMS	Naphthalene	$\mug/L$	EPA625	30.5	80.00	0	38.1	10-82				12/3/2010 0117h
1011529-006FMS	n-Decane	$\mug/L$	EPA625	10.4	80.00	0	13.0	10-27				12/3/2010 0117h
1011529-006FMS	Nitrobenzene	$\mug/L$	EPA625	56.0	80.00	0	70.0	10-119				12/3/2010 0117h
1011529-006FMS	Nitroquinoline-1-oxide	$\mug/L$	EPA625	31.1	80.00	0	38.9	10-170				12/3/2010 0117h
1011529-006FMS	N-Nitrosodiethylamine	$\mug/L$	EPA625	41.3	80.00	0	51.6	10-91				12/3/2010 0117h
1011529-006FMS	N-Nitrosodimethylamine	$\mug/L$	EPA625	17.6	80.00	0	22.1	10-42				12/3/2010 0117h
1011529-006FMS	N-Nitrosodi-n-butylamine	$\mug/L$	EPA625	56.1	80.00	0	70.1	10-175				12/3/2010 0117h
1011529-006FMS	N-Nitrosodiphenylamine	$\mug/L$	EPA625	67.0	80.00	0	83.7	12-112				12/3/2010 0117h
1011529-006FMS	N-Nitrosodi-n-propylamine	$\mug/L$	EPA625	45.4	80.00	0	56.7	10-77				12/3/2010 0117h
1011529-006FMS	N-Nitrosomethylethylamine	$\mug/L$	EPA625	33.1	80.00	0	41.4	10-75				12/3/2010 0117h
1011529-006FMS	N-Nitrosomorpholine	$\mug/L$	EPA625	48.0	80.00	0	60.0	10-175				12/3/2010 0117h
1011529-006FMS	N-Nitrosopiperidine	$\mug/L$	EPA625	54.2	80.00	0	67.7	10-105				12/3/2010 0117h
1011529-006FMS	N-Nitrosopyrrolidine	$\mug/L$	EPA625	54.4	80.00	0	68.0	10-88				12/3/2010 0117h
1011529-006FMS	n-Octadecane	$\mug/L$	EPA625	57.6	80.00	0	72.0	10-121				12/3/2010 0117h
1011529-006FMS	O,O,O-Triethyl phosphorothioate	$\mug/L$	EPA625	51.3	80.00	0	64.1	10-93				12/3/2010 0117h
1011529-006FMS	o-Toluidine	$\mug/L$	EPA625	96.4	80.00	0	120	10-107			1	12/3/2010 0117h
1011529-006FMS	Parathion	$\mug/L$	EPA625	105	80.00	0	131	10-175				12/3/2010 0117h
1011529-006FMS	Methyl parathion	$\mu g/L$	EPA625	96.9	80.00	0	121	10-175				12/3/2010 0117h
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Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMS	Pentachlorobenzene	μg/L	EPA625	55.7	80.00	0	69.6	25-134				12/3/2010 0117h
1011529-006FMS	Pentachloronitrobenzene	μg/L	EPA625	77.6	80.00	0	97.0	10-175				12/3/2010 0117h
1011529-006FMS	Pentachlorophenol	μg/L	EPA625	49.2	80.00	0	61.5	10-163				12/3/2010 0117h
1011529-006FMS	Phenacetin	μg/L	EPA625	72.3	80.00	0	90.4	10-175				12/3/2010 0117h
1011529-006FMS	Phenanthrene	μg/L	EPA625	78.6	80.00	0	98.2	31-126				12/3/2010 0117h
1011529-006FMS	Phenol	$\mu g/L$	EPA625	28.1	80.00	0	35.1	10-175				12/3/2010 0117h
1011529-006FMS	Phorate	μg/L	EPA625	74.7	80.00	0	93.4	10-175				12/3/2010 0117h
1011529-006FMS	Pronamide	μg/L	EPA625	77.1	80.00	0	96.4	10-95			1	12/3/2010 0117h
1011529-006FMS	Pyrene	μg/L	EPA625	87.9	80.00	0	110	51-139				12/3/2010 0117h
1011529-006FMS	Pyridine	μg/L	EPA625	11.7	80.00	0	14.6	10-25				12/3/2010 0117h
1011529-006FMS	Quinoline	μg/L	EPA625	57.8	80.00	0	72.3	10-63			1	12/3/2010 0117h
1011529-006FMS	Safrole	μg/L	EPA625	48.2	80.00	0	60.3	10-120				12/3/2010 0117h
1011529-006FMS	Tetraethyl dithiopyrophosphate	μg/L	EPA625	69.1	80.00	0	86.4	13-160				12/3/2010 0117h
1011529-006FMS	Thionazin	μg/L	EPA625	61.8	80.00	0	77.3	10-139				12/3/2010 0117h
1011529-006FMS	Surr: 2,4,6-Tribromophenol	%REC	EPA625	66.3	80.00		82.9	21-154				12/3/2010 0117h
1011529-006FMS	Surr: 2-Fluorobiphenyl	%REC	EPA625	45.7	40.00		114	10-106			S	12/3/2010 0117h
1011529-006FMS	Surr: 2-Fluorophenol	%REC	EPA625	25.6	80.00		32.0	10-56				12/3/2010 0117h
1011529-006FMS	Surr: Nitrobenzene-d5	%REC	EPA625	50.7	40.00		127	10-101			S	12/3/2010 0117h
1011529-006FMS	Surr: Phenol-d6	%REC	EPA625	19.0	80.00		23.8	10-45				12/3/2010 0117h
1011529-006FMS	Surr: Terphenyl-d14	%REC	EPA625	90.9	40.00		227	10-160			S	12/3/2010 0117h
1012039-001DMS	1,1´-Biphenyl	μg/L	EPA625	64.4	81.63	0	78.8	27-99				12/3/2010 0932h
1012039-001DMS	1,2,4,5-Tetrachlorobenzene	μg/L	EPA625	59.0	81.63	0	72.2	10-119				12/3/2010 0932h
1012039-001DMS	1,2,4-Trichlorobenzene	μg/L	EPA625	46.3	81.63	0	56.7	10-79				12/3/2010 0932h
1012039-001DMS	1,2-Dichlorobenzene	μg/L	EPA625	33.9	81.63	0	41.6	10-59				12/3/2010 0932h
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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**QC Type:** MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMS	1,3,5-Trinitrobenzene	μg/L	EPA625	148	81.63	0	182	10-175			1	12/3/2010 0932h
1012039-001DMS	1,3-Dichlorobenzene	μg/L	EPA625	28.5	81.63	0	34.9	10-56				12/3/2010 0932h
1012039-001DMS	1,3-Dinitrobenzene	μg/L	EPA625	94.8	81.63	0	116	10-175				12/3/2010 0932h
1012039-001DMS	1,4-Dichlorobenzene	μg/L	EPA625	30.6	81.63	0	37.5	10-58				12/3/2010 0932h
1012039-001DMS	1,4-Naphthoquinone	μg/L	EPA625	24.3	81.63	0	29.7	10-177				12/3/2010 0932h
1012039-001DMS	1,4-Phenylenediamine	μg/L	EPA625	52.0	81.63	0	63.7	10-124				12/3/2010 0932h
1012039-001DMS	1-Chloronaphthalene	μg/L	EPA625	55.2	81.63	0	67.6	10-106				12/3/2010 0932h
1012039-001DMS	1-Methylnaphthalene	μg/L	EPA625	65.7	81.63	0	80.5	10-83				12/3/2010 0932h
1012039-001DMS	1-Naphthylamine	μg/L	EPA625	70.1	81.63	0	85.9	10-122				12/3/2010 0932h
1012039-001DMS	2,3,4,6-Tetrachlorophenol	μg/L	EPA625	77.9	81.63	0	95.4	10-157				12/3/2010 0932h
1012039-001DMS	2,4,5-Trichlorophenol	μg/L	EPA625	66.3	81.63	0	81.2	10-148				12/3/2010 0932h
1012039-001DMS	2,4,6-Trichlorophenol	μg/L	EPA625	53.5	81.63	0	65.5	10-136				12/3/2010 0932h
1012039-001DMS	2,4-Dichlorophenol	μg/L	EPA625	56.8	81.63	0	69.6	10-123				12/3/2010 0932h
1012039-001DMS	2,4-Dimethylphenol	μg/L	EPA625	63.6	81.63	0	77.9	10-113				12/3/2010 0932h
1012039-001DMS	2,4-Dinitrophenol	μg/L	EPA625	57.1	81.63	0	70.0	10-175				12/3/2010 0932h
1012039-001DMS	2,4-Dinitrotoluene	μg/L	EPA625	98.9	81.63	0	121	10-175				12/3/2010 0932h
1012039-001DMS	2,6-Dichlorophenol	μg/L	EPA625	66.1	81.63	0	81.0	10-148				12/3/2010 0932h
1012039-001DMS	2,6-Dinitrotoluene	μg/L	EPA625	88.4	81.63	0	108	10-175				12/3/2010 0932h
1012039-001DMS	2-Acetylaminofluorene	μg/L	EPA625	73.9	81.63	0	90.5	10-94				12/3/2010 0932h
1012039-001DMS	2-Chloronaphthalene	μg/L	EPA625	64.6	81.63	0	79.2	10-93				12/3/2010 0932h
1012039-001DMS	2-Chlorophenol	μg/L	EPA625	54.0	81.63	0	66.1	10-92				12/3/2010 0932h
1012039-001DMS	2-Methylnaphthalene	μg/L	EPA625	56.4	81.63	0	69.1	15-78				12/3/2010 0932h
1012039-001DMS	2-Methylphenol	μg/L	EPA625	49.9	81.63	0	61.1	10-83				12/3/2010 0932h
1012039-001DMS	2-Naphthylamine	μg/L	EPA625	64.5	81.63	0	79.0	10-154				12/3/2010 0932h
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## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	% RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMS	2-Nitroaniline	μg/L	EPA625	83.8	81.63	0	103	10-175				12/3/2010 0932h
1012039-001DMS	2-Nitrophenol	μg/L	EPA625	80.6	81.63	0	98.7	10-175				12/3/2010 0932h
1012039-001DMS	2-Picoline	$\mug/L$	EPA625	37.1	81.63	0	45.5	10-61				12/3/2010 0932h
1012039-001DMS	3&4-Methylphenol	$\mug/L$	EPA625	35.2	81.63	0	43.1	10-80				12/3/2010 0932h
1012039-001DMS	3,3´-Dichlorobenzidine	$\mug/L$	EPA625	67.1	81.63	0	82.2	10-150				12/3/2010 0932h
1012039-001DMS	3,3´-Dimethylbenzidine	$\mug/L$	EPA625	71.2	81.63	0	87.2	10-143				12/3/2010 0932h
1012039-001DMS	3-Methylcholanthrene	$\mug/L$	EPA625	70.1	81.63	0	85.8	32-171				12/3/2010 0932h
1012039-001DMS	3-Nitroaniline	$\mu g/L$	EPA625	78.1	81.63	0	95.6	10-175				12/3/2010 0932h
1012039-001DMS	4,6-Dinitro-2-methylphenol	$\mu g/L$	EPA625	112	81.63	0	137	10-175				12/3/2010 0932h
1012039-001DMS	4-Aminobiphenyl	$\mu g/L$	EPA625	79.0	81.63	0	96.8	10-175				12/3/2010 0932h
1012039-001DMS	4-Bromophenyl phenyl ether	$\mu g/L$	EPA625	73.7	81.63	0	90.3	16-138				12/3/2010 0932h
1012039-001DMS	4-Chloro-3-methylphenol	$\mug/L$	EPA625	66.9	81.63	0	82.0	10-131				12/3/2010 0932h
1012039-001DMS	4-Chloroaniline	$\mug/L$	EPA625	56.8	81.63	0	69.6	10-98				12/3/2010 0932h
1012039-001DMS	4-Chlorophenyl phenyl ether	$\mu g/L$	EPA625	69.2	81.63	0	84.8	31-108				12/3/2010 0932h
1012039-001DMS	4-Nitroaniline	$\mu g/L$	EPA625	67.2	81.63	0	82.3	10-175				12/3/2010 0932h
1012039-001DMS	4-Nitrophenol	$\mu g/L$	EPA625	121	81.63	0	148	10-97			1	12/3/2010 0932h
1012039-001DMS	5-Nitro-o-toluidine	$\mu g/L$	EPA625	90.2	81.63	0	111	10-175				12/3/2010 0932h
1012039-001DMS	7,12-Dimethylbenz(a)anthracene	μg/L	EPA625	94.5	81.63	0	116	26-174				12/3/2010 0932h
1012039-001DMS	a,a-Dimethylphenethylamine	$\mu g/L$	EPA625	42.1	81.63	0	51.6	10-175				12/3/2010 0932h
1012039-001DMS	Acenaphthene	$\mu g/L$	EPA625	68.3	81.63	0	83.7	29-97				12/3/2010 0932h
1012039-001DMS	Acenaphthylene	$\mu g/L$	EPA625	67.4	81.63	0	82.6	37-87				12/3/2010 0932h
1012039-001DMS	Acetophenone	$\mu g/L$	EPA625	57.9	81.63	0	71.0	10-96				12/3/2010 0932h
1012039-001DMS	alpha-Terpineol	$\mug/L$	EPA625	55.9	81.63	0	68.5	10-67			1	12/3/2010 0932h
1012039-001DMS	Aniline	μg/L	EPA625	45.7	81.63	0	55.9	10-71				12/3/2010 0932h
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## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limit	Qual	Date Analyzed
1012039-001DMS	Anthracene	μg/L	EPA625	83.3	81.63	0	102	53-114			12/3/2010 0932h
1012039-001DMS	Aramite	μg/L	EPA625	90.6	81.63	0	111	29-160			12/3/2010 0932h
1012039-001DMS	Azobenzene	$\mu g/L$	EPA625	66.0	81.63	0	80.8	15-114			12/3/2010 0932h
1012039-001DMS	Benz(a)anthracene	μg/L	EPA625	65.8	81.63	0	80.6	39-129			12/3/2010 0932h
1012039-001DMS	Benzidine	$\mu g/L$	EPA625	59.5	81.63	0	72.9	10-99			12/3/2010 0932h
1012039-001DMS	Benzo(a)pyrene	$\mu g/L$	EPA625	73.9	81.63	0	90.5	29-175			12/3/2010 0932h
1012039-001DMS	Benzo(b)fluoranthene	$\mu g/L$	EPA625	76.4	81.63	0	93.6	15-140			12/3/2010 0932h
1012039-001DMS	Benzo(g,h,i)perylene	$\mu g/L$	EPA625	61.1	81.63	0	74.9	10-182			12/3/2010 0932h
1012039-001DMS	Benzo(k)fluoranthene	$\mu g/L$	EPA625	85.4	81.63	0	105	21-154			12/3/2010 0932h
1012039-001DMS	Benzoic acid	$\mu g/L$	EPA625	< 20.4	81.63	0	17.8	10-71			12/3/2010 0932h
1012039-001DMS	Benzyl alcohol	$\mu g/L$	EPA625	25.8	81.63	0	31.6	10-69			12/3/2010 0932h
1012039-001DMS	Bis(2-chloroethoxy)methane	$\mu g/L$	EPA625	41.2	81.63	0	50.5	10-94			12/3/2010 0932h
1012039-001DMS	Bis(2-chloroethyl) ether	$\mu g/L$	EPA625	40.9	81.63	0	50.1	10-70			12/3/2010 0932h
1012039-001DMS	Bis(2-chloroisopropyl) ether	$\mu g/L$	EPA625	43.6	81.63	0	53.4	10-71			12/3/2010 0932h
1012039-001DMS	Bis(2-ethylhexyl) phthalate	$\mu g/L$	EPA625	85.2	81.63	0	104	10-175			12/3/2010 0932h
1012039-001DMS	bis(2-ethylhexyl)adipate	$\mu g/L$	EPA625	85.6	81.63	0	105	10-175			12/3/2010 0932h
1012039-001DMS	Butyl benzyl phthalate	$\mu g/L$	EPA625	84.9	81.63	0	104	10-175			12/3/2010 0932h
1012039-001DMS	Carbazole	$\mu g/L$	EPA625	77.0	81.63	0	94.3	10-151			12/3/2010 0932h
1012039-001DMS	Chlorobenzilate	$\mu g/L$	EPA625	81.5	81.63	0	99.8	18-175			12/3/2010 0932h
1012039-001DMS	Chrysene	$\mu g/L$	EPA625	73.5	81.63	0	90.0	38-133			12/3/2010 0932h
1012039-001DMS	Diallate (cis or trans)	$\mug/L$	EPA625	71.1	81.63	0	87.1	10-157			12/3/2010 0932h
1012039-001DMS	Dibenz(a,h)anthracene	$\mug/L$	EPA625	61.8	81.63	0	75.7	13-168			12/3/2010 0932h
1012039-001DMS	Dibenzofuran	$\mu g/L$	EPA625	66.6	81.63	0	81.6	29-103			12/3/2010 0932h
1012039-001DMS	Diethyl phthalate	$\mu g/L$	EPA625	72.7	81.63	0	89.1	10-139			12/3/2010 0932h
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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD Limit	Qual	Date Analyzed
1012039-001DMS	Dimethoate	μg/L	EPA625	33.8	81.63	0	41.4	10-136			12/3/2010 0932h
1012039-001DMS	Dimethyl phthalate	μg/L	EPA625	65.3	81.63	0	80.0	10-122			12/3/2010 0932h
1012039-001DMS	Dimethylaminoazobenzene	$\mug/L$	EPA625	80.6	81.63	0	98.7	34-142			12/3/2010 0932h
1012039-001DMS	Di-n-butyl phthalate	$\mu g/L$	EPA625	82.8	81.63	0	101	44-124			12/3/2010 0932h
1012039-001DMS	Di-n-octyl phthalate	$\mu g/L$	EPA625	123	81.63	0	150	10-175			12/3/2010 0932h
1012039-001DMS	Dinoseb	$\mug/L$	EPA625	125	81.63	0	153	10-175			12/3/2010 0932h
1012039-001DMS	Diphenylamine	$\mug/L$	EPA625	76.8	81.63	0	94.1	13-110			12/3/2010 0932h
1012039-001DMS	Disulfoton	$\mu g/L$	EPA625	67.5	81.63	0	82.7	10-121			12/3/2010 0932h
1012039-001DMS	Ethyl methanesulfonate	$\mu g/L$	EPA625	51.5	81.63	0	63.1	10-99			12/3/2010 0932h
1012039-001DMS	Famphur	$\mu g/L$	EPA625	123	81.63	0	150	10-71		1	12/3/2010 0932h
1012039-001DMS	Fluoranthene	$\mu g/L$	EPA625	76.3	81.63	0	93.5	23-135			12/3/2010 0932h
1012039-001DMS	Fluorene	$\mu g/L$	EPA625	70.8	81.63	0	86.7	34-108			12/3/2010 0932h
1012039-001DMS	Hexachlorobenzene	$\mug/L$	EPA625	74.8	81.63	0	91.7	26-131			12/3/2010 0932h
1012039-001DMS	Hexachlorobutadiene	$\mu g/L$	EPA625	41.6	81.63	0	51.0	10-110			12/3/2010 0932h
1012039-001DMS	Hexachlorocyclopentadiene	$\mu g/L$	EPA625	27.2	81.63	0	33.4	10-45			12/3/2010 0932h
1012039-001DMS	Hexachloroethane	$\mu g/L$	EPA625	30.2	81.63	0	37.0	10-58			12/3/2010 0932h
1012039-001DMS	Hexachlorophene	$\mug/L$	EPA625	75.1	81.63	0	92.0	10-168			12/3/2010 0932h
1012039-001DMS	Hexachloropropene	$\mu g/L$	EPA625	38.1	81.63	0	46.7	10-72			12/3/2010 0932h
1012039-001DMS	Indene	$\mu g/L$	EPA625	40.8	81.63	0	50.0	10-35		1	12/3/2010 0932h
1012039-001DMS	Indeno(1,2,3-cd)pyrene	$\mu g/L$	EPA625	63.4	81.63	0	77.7	10-176			12/3/2010 0932h
1012039-001DMS	Isodrin	$\mug/L$	EPA625	76.9	81.63	0	94.2	15-165			12/3/2010 0932h
1012039-001DMS	Isophorone	$\mug/L$	EPA625	58.2	81.63	0	71.3	10-99			12/3/2010 0932h
1012039-001DMS	Isosafrole	$\mu g/L$	EPA625	59.8	81.63	0	73.3	10-167			12/3/2010 0932h
1012039-001DMS	Kepone	$\mug/L$	EPA625	205	81.63	0	252	10-175		1	12/3/2010 0932h

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Laboratory Director

Jose Rocha

QA Officer

## **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**QC Type:** MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD Limit	Qual	Date Analyzed
1012039-001DMS	Methapyrilene	μg/L	EPA625	88.4	81.63	0	108	10-149			12/3/2010 0932h
1012039-001DMS	Methyl methanesulfonate	μg/L	EPA625	36.0	81.63	0	44.1	10-132			12/3/2010 0932h
1012039-001DMS	Naphthalene	μg/L	EPA625	53.1	81.63	0	65.0	10-82			12/3/2010 0932h
1012039-001DMS	n-Decane	μg/L	EPA625	19.2	81.63	0	23.5	10-27			12/3/2010 0932h
1012039-001DMS	Nitrobenzene	μg/L	EPA625	67.8	81.63	0	83.0	10-119			12/3/2010 0932h
1012039-001DMS	Nitroquinoline-1-oxide	μg/L	EPA625	30.3	81.63	0	37.1	10-170			12/3/2010 0932h
1012039-001DMS	N-Nitrosodiethylamine	$\mu g/L$	EPA625	50.0	81.63	0	61.2	10-91			12/3/2010 0932h
1012039-001DMS	N-Nitrosodimethylamine	$\mu g/L$	EPA625	20.1	81.63	0	24.6	10-42			12/3/2010 0932h
1012039-001DMS	N-Nitrosodi-n-butylamine	μg/L	EPA625	63.7	81.63	0	78.1	10-175			12/3/2010 0932h
1012039-001DMS	N-Nitrosodiphenylamine	μg/L	EPA625	75.6	81.63	0	92.7	12-112			12/3/2010 0932h
1012039-001DMS	N-Nitrosodi-n-propylamine	$\mu g/L$	EPA625	53.9	81.63	0	66.1	10-77			12/3/2010 0932h
1012039-001DMS	N-Nitrosomethylethylamine	$\mu g/L$	EPA625	39.8	81.63	0	48.8	10-75			12/3/2010 0932h
1012039-001DMS	N-Nitrosomorpholine	$\mu g/L$	EPA625	55.8	81.63	0	68.3	10-175			12/3/2010 0932h
1012039-001DMS	N-Nitrosopiperidine	$\mug/L$	EPA625	61.2	81.63	0	75.0	10-105			12/3/2010 0932h
1012039-001DMS	N-Nitrosopyrrolidine	$\mu g/L$	EPA625	63.3	81.63	0	77.6	10-88			12/3/2010 0932h
1012039-001DMS	n-Octadecane	$\mu g/L$	EPA625	67.9	81.63	0	83.2	10-121			12/3/2010 0932h
1012039-001DMS	O,O,O-Triethyl phosphorothioate	$\mu g/L$	EPA625	63.9	81.63	0	78.2	10-93			12/3/2010 0932h
1012039-001DMS	o-Toluidine	$\mu g/L$	EPA625	114	81.63	0	140	10-107		1	12/3/2010 0932h
1012039-001DMS	Parathion	$\mu g/L$	EPA625	110	81.63	0	134	10-175			12/3/2010 0932h
1012039-001DMS	Methyl parathion	$\mu g/L$	EPA625	100	81.63	0	123	10-175			12/3/2010 0932h
1012039-001DMS	Pentachlorobenzene	$\mug/L$	EPA625	66.5	81.63	0	81.4	25-134			12/3/2010 0932h
1012039-001DMS	Pentachloronitrobenzene	$\mug/L$	EPA625	83.7	81.63	0	103	10-175			12/3/2010 0932h
1012039-001DMS	Pentachlorophenol	$\mug/L$	EPA625	44.1	81.63	0	54.1	10-163			12/3/2010 0932h
1012039-001DMS	Phenacetin	$\mu g/L$	EPA625	72.1	81.63	0	88.3	10-175			12/3/2010 0932h
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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMS	Phenanthrene	μg/L	EPA625	88.6	81.63	0	109	31-126				12/3/2010 0932h
1012039-001DMS	Phenol	μg/L	EPA625	32.6	81.63	0	40.0	10-175				12/3/2010 0932h
1012039-001DMS	Phorate	μg/L	EPA625	83.9	81.63	0	103	10-175				12/3/2010 0932h
1012039-001DMS	Pronamide	μg/L	EPA625	78.9	81.63	0	96.6	10-95			1	12/3/2010 0932h
1012039-001DMS	Pyrene	μg/L	EPA625	84.4	81.63	0	103	51-139				12/3/2010 0932h
1012039-001DMS	Pyridine	μg/L	EPA625	20.9	81.63	0	25.6	10-25			1	12/3/2010 0932h
1012039-001DMS	Quinoline	μg/L	EPA625	65.0	81.63	0	79.6	10-63			1	12/3/2010 0932h
1012039-001DMS	Safrole	μg/L	EPA625	58.6	81.63	0	71.8	10-120				12/3/2010 0932h
1012039-001DMS	Tetraethyl dithiopyrophosphate	μg/L	EPA625	77.9	81.63	0	95.5	13-160				12/3/2010 0932h
1012039-001DMS	Thionazin	$\mu g/L$	EPA625	69.5	81.63	0	85.1	10-139				12/3/2010 0932h
1012039-001DMS	Surr: 2,4,6-Tribromophenol	%REC	EPA625	66.9	81.63		82.0	21-154				12/3/2010 0932h
1012039-001DMS	Surr: 2-Fluorobiphenyl	%REC	EPA625	61.7	40.82		151	10-106			S	12/3/2010 0932h
1012039-001DMS	Surr: 2-Fluorophenol	%REC	EPA625	29.2	81.63		35.8	10-56				12/3/2010 0932h
1012039-001DMS	Surr: Nitrobenzene-d5	%REC	EPA625	61.3	40.82		150	10-101			S	12/3/2010 0932h
1012039-001DMS	Surr: Phenol-d6	%REC	EPA625	19.8	81.63		24.2	10-45				12/3/2010 0932h
1012039-001DMS	Surr: Terphenyl-d14	%REC	EPA625	84.2	40.82		206	10-160			S	12/3/2010 0932h
1012057-001BMS	1,1´-Biphenyl	$\mug/L$	EPA625	51.8	80.00	0	64.8	27-99				12/3/2010 1121h
1012057-001BMS	1,2,4,5-Tetrachlorobenzene	$\mug/L$	EPA625	47.4	80.00	0	59.2	10-119				12/3/2010 1121h
1012057-001BMS	1,2,4-Trichlorobenzene	$\mug/L$	EPA625	37.8	80.00	0	47.2	10-79				12/3/2010 1121h
1012057-001BMS	1,2-Dichlorobenzene	$\mug/L$	EPA625	29.2	80.00	0	36.5	10-59				12/3/2010 1121h
1012057-001BMS	1,3,5-Trinitrobenzene	$\mug/L$	EPA625	134	80.00	0	167	10-175				12/3/2010 1121h
1012057-001BMS	1,3-Dichlorobenzene	$\mug/L$	EPA625	24.6	80.00	0	30.8	10-56				12/3/2010 1121h
1012057-001BMS	1,3-Dinitrobenzene	$\mug/L$	EPA625	75.9	80.00	0	94.9	10-175				12/3/2010 1121h
1012057-001BMS	1,4-Dichlorobenzene	$\mug\!/L$	EPA625	25.5	80.00	0	31.9	10-58				12/3/2010 1121h
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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPI %RPD Lim		Date Analyzed
1012057-001BMS	1,4-Naphthoquinone	μg/L	EPA625	20.5	80.00	0	25.6	10-177			12/3/2010 1121h
1012057-001BMS	1,4-Phenylenediamine	$\mu g/L$	EPA625	41.7	80.00	0	52.1	10-124			12/3/2010 1121h
1012057-001BMS	1-Chloronaphthalene	$\mu g/L$	EPA625	42.7	80.00	0	53.3	10-106			12/3/2010 1121h
1012057-001BMS	1-Methylnaphthalene	$\mu g/L$	EPA625	52.8	80.00	0	66.0	10-83			12/3/2010 1121h
1012057-001BMS	1-Naphthylamine	$\mu g/L$	EPA625	59.5	80.00	0	74.4	10-122			12/3/2010 1121h
1012057-001BMS	2,3,4,6-Tetrachlorophenol	$\mu g/L$	EPA625	79.2	80.00	0	99.0	10-157			12/3/2010 1121h
1012057-001BMS	2,4,5-Trichlorophenol	$\mu g/L$	EPA625	59.8	80.00	0	74.8	10-148			12/3/2010 1121h
1012057-001BMS	2,4,6-Trichlorophenol	$\mu g/L$	EPA625	47.0	80.00	0	58.8	10-136			12/3/2010 1121h
1012057-001BMS	2,4-Dichlorophenol	$\mu g/L$	EPA625	49.0	80.00	0	61.3	10-123			12/3/2010 1121h
1012057-001BMS	2,4-Dimethylphenol	$\mu g/L$	EPA625	55.3	80.00	0	69.1	10-113			12/3/2010 1121h
1012057-001BMS	2,4-Dinitrophenol	$\mu g/L$	EPA625	50.2	80.00	0	62.8	10-175			12/3/2010 1121h
1012057-001BMS	2,4-Dinitrotoluene	$\mu g/L$	EPA625	86.3	80.00	0	108	10-175			12/3/2010 1121h
1012057-001BMS	2,6-Dichlorophenol	$\mu g/L$	EPA625	60.4	80.00	0	75.6	10-148			12/3/2010 1121h
1012057-001BMS	2,6-Dinitrotoluene	$\mu g/L$	EPA625	74.4	80.00	0	93.0	10-175			12/3/2010 1121h
1012057-001BMS	2-Acetylaminofluorene	$\mu g/L$	EPA625	68.9	80.00	0	86.1	10-94			12/3/2010 1121h
1012057-001BMS	2-Chloronaphthalene	$\mu g/L$	EPA625	46.3	80.00	0	57.8	10-93			12/3/2010 1121h
1012057-001BMS	2-Chlorophenol	$\mu g/L$	EPA625	45.1	80.00	0	56.4	10-92			12/3/2010 1121h
1012057-001BMS	2-Methylnaphthalene	$\mu g/L$	EPA625	45.3	80.00	0	56.6	15-78			12/3/2010 1121h
1012057-001BMS	2-Methylphenol	$\mu g/L$	EPA625	36.6	80.00	0	45.7	10-83			12/3/2010 1121h
1012057-001BMS	2-Naphthylamine	$\mu g/L$	EPA625	54.0	80.00	0	67.5	10-154			12/3/2010 1121h
1012057-001BMS	2-Nitroaniline	$\mug/L$	EPA625	72.1	80.00	0	90.2	10-175			12/3/2010 1121h
1012057-001BMS	2-Nitrophenol	$\mug/L$	EPA625	74.9	80.00	0	93.6	10-175			12/3/2010 1121h
1012057-001BMS	2-Picoline	$\mug/L$	EPA625	24.2	80.00	0	30.3	10-61			12/3/2010 1121h
1012057-001BMS	3&4-Methylphenol	$\mu g/L$	EPA625	19.9	80.00	0	24.9	10-80			12/3/2010 1121h
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QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	% RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMS	3,3'-Dichlorobenzidine	μg/L	EPA625	63.0	80.00	0	78.7	10-150				12/3/2010 1121h
1012057-001BMS	3,3´-Dimethylbenzidine	μg/L	EPA625	71.7	80.00	0	89.7	10-143				12/3/2010 1121h
1012057-001BMS	3-Methylcholanthrene	$\mug/L$	EPA625	66.4	80.00	0	83.0	32-171				12/3/2010 1121h
1012057-001BMS	3-Nitroaniline	μg/L	EPA625	67.7	80.00	0	84.7	10-175				12/3/2010 1121h
1012057-001BMS	4,6-Dinitro-2-methylphenol	$\mug/L$	EPA625	101	80.00	0	127	10-175				12/3/2010 1121h
1012057-001BMS	4-Aminobiphenyl	$\mug/L$	EPA625	67.1	80.00	0	83.9	10-175				12/3/2010 1121h
1012057-001BMS	4-Bromophenyl phenyl ether	$\mug/L$	EPA625	59.5	80.00	0	74.4	16-138				12/3/2010 1121h
1012057-001BMS	4-Chloro-3-methylphenol	$\mug/L$	EPA625	56.9	80.00	0	71.1	10-131				12/3/2010 1121h
1012057-001BMS	4-Chloroaniline	$\mug/L$	EPA625	47.4	80.00	0	59.3	10-98				12/3/2010 1121h
1012057-001BMS	4-Chlorophenyl phenyl ether	μg/L	EPA625	57.6	80.00	0	72.1	31-108				12/3/2010 1121h
1012057-001BMS	4-Nitroaniline	μg/L	EPA625	59.8	80.00	0	74.7	10-175				12/3/2010 1121h
1012057-001BMS	4-Nitrophenol	$\mug/L$	EPA625	101	80.00	0	126	10-97			1	12/3/2010 1121h
1012057-001BMS	5-Nitro-o-toluidine	$\mug/L$	EPA625	78.7	80.00	0	98.4	10-175				12/3/2010 1121h
1012057-001BMS	7,12-Dimethylbenz(a)anthracene	$\mug/L$	EPA625	89.1	80.00	0	111	26-174				12/3/2010 1121h
1012057-001BMS	a,a-Dimethylphenethylamine	μg/L	EPA625	39.1	80.00	0	48.9	10-175				12/3/2010 1121h
1012057-001BMS	Acenaphthene	μg/L	EPA625	56.0	80.00	0	70.0	29-97				12/3/2010 1121h
1012057-001BMS	Acenaphthylene	$\mu g/L$	EPA625	56.1	80.00	0	70.1	37-87				12/3/2010 1121h
1012057-001BMS	Acetophenone	$\mug/L$	EPA625	46.5	80.00	0	58.1	10-96				12/3/2010 1121h
1012057-001BMS	alpha-Terpineol	$\mug/L$	EPA625	46.2	80.00	0	57.8	10-67				12/3/2010 1121h
1012057-001BMS	Aniline	$\mug/L$	EPA625	32.0	80.00	0	40.0	10-71				12/3/2010 1121h
1012057-001BMS	Anthracene	$\mug/L$	EPA625	69.5	80.00	0	86.9	53-114				12/3/2010 1121h
1012057-001BMS	Aramite	$\mug/L$	EPA625	76.1	80.00	0	95.2	29-160				12/3/2010 1121h
1012057-001BMS	Azobenzene	$\mu g/L$	EPA625	55.7	80.00	0	69.6	15-114				12/3/2010 1121h
1012057-001BMS	Benz(a)anthracene	μg/L	EPA625	63.0	80.00	0	78.7	39-129				12/3/2010 1121h
										Daport	Data: 12/	6/2010 Page 111 of 16

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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limi	Qual	Date Analyzed
1012057-001BMS	Benzidine	μg/L	EPA625	56.9	80.00	0	71.1	10-99			12/3/2010 1121h
1012057-001BMS	Benzo(a)pyrene	$\mug/L$	EPA625	71.0	80.00	0	88.7	29-175			12/3/2010 1121h
1012057-001BMS	Benzo(b)fluoranthene	$\mug/L$	EPA625	61.1	80.00	0	76.3	15-140			12/3/2010 1121h
1012057-001BMS	Benzo(g,h,i)perylene	$\mug/L$	EPA625	57.9	80.00	0	72.4	10-182			12/3/2010 1121h
1012057-001BMS	Benzo(k)fluoranthene	$\mug/L$	EPA625	86.7	80.00	0	108	21-154			12/3/2010 1121h
1012057-001BMS	Benzoic acid	$\mug/L$	EPA625	< 20.0	80.00	0	17.5	10-71			12/3/2010 1121h
1012057-001BMS	Benzyl alcohol	$\mug/L$	EPA625	18.5	80.00	0	23.2	10-69			12/3/2010 1121h
1012057-001BMS	Bis(2-chloroethoxy)methane	$\mug/L$	EPA625	34.3	80.00	0	42.9	10-94			12/3/2010 1121h
1012057-001BMS	Bis(2-chloroethyl) ether	$\mug/L$	EPA625	34.9	80.00	0	43.6	10-70			12/3/2010 1121h
1012057-001BMS	Bis(2-chloroisopropyl) ether	$\mug/L$	EPA625	33.5	80.00	0	41.9	10-71			12/3/2010 1121h
1012057-001BMS	Bis(2-ethylhexyl) phthalate	$\mug/L$	EPA625	77.8	80.00	0	97.2	10-175			12/3/2010 1121h
1012057-001BMS	bis(2-ethylhexyl)adipate	$\mug/L$	EPA625	80.9	80.00	0	101	10-175			12/3/2010 1121h
1012057-001BMS	Butyl benzyl phthalate	$\mug/L$	EPA625	81.1	80.00	0	101	10-175			12/3/2010 1121h
1012057-001BMS	Carbazole	$\mug/L$	EPA625	66.5	80.00	0	83.1	10-151			12/3/2010 1121h
1012057-001BMS	Chlorobenzilate	$\mug/L$	EPA625	77.8	80.00	0	97.2	18-175			12/3/2010 1121h
1012057-001BMS	Chrysene	$\mug/L$	EPA625	69.8	80.00	0	87.2	38-133			12/3/2010 1121h
1012057-001BMS	Diallate (cis or trans)	$\mug/L$	EPA625	58.5	80.00	0	73.1	10-157			12/3/2010 1121h
1012057-001BMS	Dibenz(a,h)anthracene	$\mug/L$	EPA625	59.0	80.00	0	73.8	13-168			12/3/2010 1121h
1012057-001BMS	Dibenzofuran	$\mug/L$	EPA625	55.4	80.00	0	69.3	29-103			12/3/2010 1121h
1012057-001BMS	Diethyl phthalate	$\mug/L$	EPA625	60.1	80.00	0	75.1	10-139			12/3/2010 1121h
1012057-001BMS	Dimethoate	$\mug/L$	EPA625	21.6	80.00	0	27.0	10-136			12/3/2010 1121h
1012057-001BMS	Dimethyl phthalate	$\mug/L$	EPA625	52.8	80.00	0	66.0	10-122			12/3/2010 1121h
1012057-001BMS	Dimethylaminoazobenzene	$\mug/L$	EPA625	74.2	80.00	0	92.8	34-142			12/3/2010 1121h
1012057-001BMS	Di-n-butyl phthalate	$\mug/L$	EPA625	70.7	80.00	0	88.4	44-124			12/3/2010 1121h

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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD Limit	Qual	Date Analyzed
1012057-001BMS	Di-n-octyl phthalate	μg/L	EPA625	116	80.00	0	145	10-175			12/3/2010 1121h
1012057-001BMS	Dinoseb	μg/L	EPA625	114	80.00	0	142	10-175			12/3/2010 1121h
1012057-001BMS	Diphenylamine	$\mug/L$	EPA625	64.9	80.00	0	81.2	13-110			12/3/2010 1121h
1012057-001BMS	Disulfoton	$\mu g/L$	EPA625	48.5	80.00	0	60.7	10-121			12/3/2010 1121h
1012057-001BMS	Ethyl methanesulfonate	μg/L	EPA625	41.2	80.00	0	51.5	10-99			12/3/2010 1121h
1012057-001BMS	Famphur	$\mug/L$	EPA625	119	80.00	0	149	10-71		1	12/3/2010 1121h
1012057-001BMS	Fluoranthene	$\mug/L$	EPA625	66.9	80.00	0	83.6	23-135			12/3/2010 1121h
1012057-001BMS	Fluorene	μg/L	EPA625	57.6	80.00	0	72.0	34-108			12/3/2010 1121h
1012057-001BMS	Hexachlorobenzene	μg/L	EPA625	61.0	80.00	0	76.2	26-131			12/3/2010 1121h
1012057-001BMS	Hexachlorobutadiene	μg/L	EPA625	33.6	80.00	0	41.9	10-110			12/3/2010 1121h
1012057-001BMS	Hexachlorocyclopentadiene	μg/L	EPA625	21.2	80.00	0	26.4	10-45			12/3/2010 1121h
1012057-001BMS	Hexachloroethane	$\mug/L$	EPA625	26.9	80.00	0	33.6	10-58			12/3/2010 1121h
1012057-001BMS	Hexachlorophene	$\mug/L$	EPA625	75.1	80.00	0	93.8	10-168			12/3/2010 1121h
1012057-001BMS	Hexachloropropene	$\mug/L$	EPA625	31.6	80.00	0	39.6	10-72			12/3/2010 1121h
1012057-001BMS	Indene	$\mug/L$	EPA625	33.0	80.00	0	41.2	10-35		1	12/3/2010 1121h
1012057-001BMS	Indeno(1,2,3-cd)pyrene	μg/L	EPA625	60.1	80.00	0	75.2	10-176			12/3/2010 1121h
1012057-001BMS	Isodrin	$\mug/L$	EPA625	65.4	80.00	0	81.7	15-165			12/3/2010 1121h
1012057-001BMS	Isophorone	μg/L	EPA625	45.9	80.00	0	57.4	10-99			12/3/2010 1121h
1012057-001BMS	Isosafrole	μg/L	EPA625	46.4	80.00	0	58.0	10-167			12/3/2010 1121h
1012057-001BMS	Kepone	μg/L	EPA625	211	80.00	0	264	10-175		1	12/3/2010 1121h
1012057-001BMS	Methapyrilene	$\mug/L$	EPA625	73.9	80.00	0	92.4	10-149			12/3/2010 1121h
1012057-001BMS	Methyl methanesulfonate	$\mu g/L$	EPA625	25.8	80.00	0	32.2	10-132			12/3/2010 1121h
1012057-001BMS	Naphthalene	$\mu g/L$	EPA625	41.3	80.00	0	51.7	10-82			12/3/2010 1121h
1012057-001BMS	n-Decane	$\mu g/L$	EPA625	17.4	80.00	0	21.8	10-27			12/3/2010 1121h
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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

**QC Type:** MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limit	Qual	Date Analyzed
1012057-001BMS	Nitrobenzene	μg/L	EPA625	53.9	80.00	0	67.4	10-119			12/3/2010 1121h
1012057-001BMS	Nitroquinoline-1-oxide	μg/L	EPA625	29.3	80.00	0	36.6	10-170			12/3/2010 1121h
1012057-001BMS	N-Nitrosodiethylamine	μg/L	EPA625	40.5	80.00	0	50.6	10-91			12/3/2010 1121h
1012057-001BMS	N-Nitrosodimethylamine	μg/L	EPA625	15.3	80.00	0	19.2	10-42			12/3/2010 1121h
1012057-001BMS	N-Nitrosodi-n-butylamine	μg/L	EPA625	52.5	80.00	0	65.7	10-175			12/3/2010 1121h
1012057-001BMS	N-Nitrosodiphenylamine	μg/L	EPA625	63.9	80.00	0	79.9	12-112			12/3/2010 1121h
1012057-001BMS	N-Nitrosodi-n-propylamine	μg/L	EPA625	41.3	80.00	0	51.6	10-77			12/3/2010 1121h
1012057-001BMS	N-Nitrosomethylethylamine	μg/L	EPA625	30.7	80.00	0	38.4	10-75			12/3/2010 1121h
1012057-001BMS	N-Nitrosomorpholine	μg/L	EPA625	43.8	80.00	0	54.8	10-175			12/3/2010 1121h
1012057-001BMS	N-Nitrosopiperidine	μg/L	EPA625	51.9	80.00	0	64.9	10-105			12/3/2010 1121h
1012057-001BMS	N-Nitrosopyrrolidine	μg/L	EPA625	49.6	80.00	0	62.0	10-88			12/3/2010 1121h
1012057-001BMS	n-Octadecane	μg/L	EPA625	53.7	80.00	0	67.1	10-121			12/3/2010 1121h
1012057-001BMS	O,O,O-Triethyl phosphorothioate	μg/L	EPA625	51.2	80.00	0	64.0	10-93			12/3/2010 1121h
1012057-001BMS	o-Toluidine	μg/L	EPA625	92.1	80.00	0	115	10-107		1	12/3/2010 1121h
1012057-001BMS	Parathion	μg/L	EPA625	94.6	80.00	0	118	10-175			12/3/2010 1121h
1012057-001BMS	Methyl parathion	μg/L	EPA625	87.4	80.00	0	109	10-175			12/3/2010 1121h
1012057-001BMS	Pentachlorobenzene	μg/L	EPA625	55.0	80.00	0	68.8	25-134			12/3/2010 1121h
1012057-001BMS	Pentachloronitrobenzene	μg/L	EPA625	71.5	80.00	0	89.3	10-175			12/3/2010 1121h
1012057-001BMS	Pentachlorophenol	μg/L	EPA625	45.1	80.00	0	56.3	10-163			12/3/2010 1121h
1012057-001BMS	Phenacetin	μg/L	EPA625	63.4	80.00	0	79.2	10-175			12/3/2010 1121h
1012057-001BMS	Phenanthrene	μg/L	EPA625	69.6	80.00	0	87.0	31-126			12/3/2010 1121h
1012057-001BMS	Phenol	μg/L	EPA625	21.3	80.00	0	26.6	10-175			12/3/2010 1121h
1012057-001BMS	Phorate	μg/L	EPA625	71.4	80.00	0	89.2	10-175			12/3/2010 1121h
1012057-001BMS	Pronamide	μg/L	EPA625	71.3	80.00	0	89.1	10-95			12/3/2010 1121h
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QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMS	Pyrene	μg/L	EPA625	77.6	80.00	0	97.0	51-139				12/3/2010 1121h
1012057-001BMS	Pyridine	μg/L	EPA625	13.5	80.00	0	16.8	10-25				12/3/2010 1121h
1012057-001BMS	Quinoline	$\mu$ g/L	EPA625	54.5	80.00	0	68.1	10-63			1	12/3/2010 1121h
1012057-001BMS	Safrole	μg/L	EPA625	46.6	80.00	0	58.3	10-120				12/3/2010 1121h
1012057-001BMS	Tetraethyl dithiopyrophosphate	μg/L	EPA625	62.3	80.00	0	77.9	13-160				12/3/2010 1121h
1012057-001BMS	Thionazin	$\mu$ g/L	EPA625	59.2	80.00	0	74.0	10-139				12/3/2010 1121h
1012057-001BMS	Surr: 2,4,6-Tribromophenol	%REC	EPA625	63.0	80.00		78.8	21-154				12/3/2010 1121h
1012057-001BMS	Surr: 2-Fluorobiphenyl	%REC	EPA625	51.7	40.00		129	10-106			S	12/3/2010 1121h
1012057-001BMS	Surr: 2-Fluorophenol	%REC	EPA625	20.1	80.00		25.2	10-56				12/3/2010 1121h
1012057-001BMS	Surr: Nitrobenzene-d5	%REC	EPA625	49.2	40.00		123	10-101			S	12/3/2010 1121h
1012057-001BMS	Surr: Phenol-d6	%REC	EPA625	12.4	80.00		15.5	10-45				12/3/2010 1121h
1012057-001BMS	Surr: Terphenyl-d14	%REC	EPA625	80.2	40.00		200	10-160			S	12/3/2010 1121h
1011529-006FMS	Acenaphthene	$\mug/L$	SW8270D	54.7	80.00	0	68.4	21-113				12/3/2010 0117h
1011529-006FMS	Benzo(a)pyrene	$\mug/L$	SW8270D	79.5	80.00	0	99.4	15-169				12/3/2010 0117h
1011529-006FMS	Pyrene	$\mug/L$	SW8270D	87.9	80.00	0	110	23-150				12/3/2010 0117h
1011529-006FMS	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	66.3	80.00		82.9	14-159				12/3/2010 0117h
1011529-006FMS	Surr: 2-Fluorobiphenyl	%REC	SW8270D	45.7	40.00		114	10-124				12/3/2010 0117h
1011529-006FMS	Surr: 2-Fluorophenol	%REC	SW8270D	25.6	80.00		32.0	10-106				12/3/2010 0117h
1011529-006FMS	Surr: Nitrobenzene-d5	%REC	SW8270D	50.7	40.00		127	10-180				12/3/2010 0117h
1011529-006FMS	Surr: Phenol-d6	%REC	SW8270D	19.0	80.00		23.8	10-122				12/3/2010 0117h
1011529-006FMS	Surr: Terphenyl-d14	%REC	SW8270D	90.9	40.00		227	10-199			S	12/3/2010 0117h
1012039-001DMS	Acenaphthene	$\mug/L$	SW8270D	68.3	81.63	0	83.7	21-113				12/3/2010 0932h
1012039-001DMS	Benzo(a)pyrene	$\mug/L$	SW8270D	73.9	81.63	0	90.5	15-169				12/3/2010 0932h
1012039-001DMS	Pyrene	μg/L	SW8270D	84.4	81.63	0	103	23-150				12/3/2010 0932h
										D	D-4 12	/6/2010 Dags 115 of 16'

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Jose Rocha

QA Officer

### **OC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMS	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	66.9	81.63		82.0	14-159				12/3/2010 0932h
1012039-001DMS	Surr: 2-Fluorobiphenyl	%REC	SW8270D	61.7	40.82		151	10-124			S	12/3/2010 0932h
1012039-001DMS	Surr: 2-Fluorophenol	%REC	SW8270D	29.2	81.63		35.8	10-106				12/3/2010 0932h
1012039-001DMS	Surr: Nitrobenzene-d5	%REC	SW8270D	61.3	40.82		150	10-180				12/3/2010 0932h
1012039-001DMS	Surr: Phenol-d6	%REC	SW8270D	19.8	81.63		24.2	10-122				12/3/2010 0932h
1012039-001DMS	Surr: Terphenyl-d14	%REC	SW8270D	84.2	40.82		206	10-199			S	12/3/2010 0932h
1012057-001BMS	Acenaphthene	$\mug/L$	SW8270D	56.0	80.00	0	70.0	21-113				12/3/2010 1121h
1012057-001BMS	Benzo(a)pyrene	$\mug/L$	SW8270D	71.0	80.00	0	88.7	15-169				12/3/2010 1121h
1012057-001BMS	Pyrene	$\mug/L$	SW8270D	77.6	80.00	0	97.0	23-150				12/3/2010 1121h
1012057-001BMS	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	63.0	80.00		78.8	14-159				12/3/2010 1121h
1012057-001BMS	Surr: 2-Fluorobiphenyl	%REC	SW8270D	51.7	40.00		129	10-124			S	12/3/2010 1121h
1012057-001BMS	Surr: 2-Fluorophenol	%REC	SW8270D	20.1	80.00		25.2	10-106				12/3/2010 1121h
1012057-001BMS	Surr: Nitrobenzene-d5	%REC	SW8270D	49.2	40.00		123	10-180				12/3/2010 1121h
1012057-001BMS	Surr: Phenol-d6	%REC	SW8270D	12.4	80.00		15.5	10-122				12/3/2010 1121h
1012057-001BMS	Surr: Terphenyl-d14	%REC	SW8270D	80.2	40.00		200	10-199			S	12/3/2010 1121h

<sup>&</sup>lt;sup>1</sup> - Matrix spike recovery indicates matrix interference. The method is in control as indicated by the LCS.

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S - Surrogate recoveries outside the control limits. Repreparation and reanalysis and/or MS samples yielded similar results indicating matrix interference.



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QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMSD	1,1´-Biphenyl	μg/L	EPA625	51.6	80.00	0	64.5	27-99	6.64	49		12/3/2010 0144h
1011529-006FMSD	1,2,4,5-Tetrachlorobenzene	$\mu g/L$	EPA625	45.1	80.00	0	56.3	10-119	12.4	52		12/3/2010 0144h
1011529-006FMSD	1,2,4-Trichlorobenzene	$\mu g/L$	EPA625	28.8	80.00	0	36.0	10-79	17.9	49		12/3/2010 0144h
1011529-006FMSD	1,2-Dichlorobenzene	$\mu g/L$	EPA625	22.0	80.00	0	27.6	10-59	17.9	46		12/3/2010 0144h
1011529-006FMSD	1,3,5-Trinitrobenzene	$\mu g/L$	EPA625	150	80.00	0	188	10-175	8.61	33	1	12/3/2010 0144h
1011529-006FMSD	1,3-Dichlorobenzene	$\mu g/L$	EPA625	18.2	80.00	0	22.8	10-56	19.2	49		12/3/2010 0144h
1011529-006FMSD	1,3-Dinitrobenzene	$\mu g/L$	EPA625	96.6	80.00	0	121	10-175	0.269	29		12/3/2010 0144h
1011529-006FMSD	1,4-Dichlorobenzene	$\mu g/L$	EPA625	19.8	80.00	0	24.7	10-58	18.6	51		12/3/2010 0144h
1011529-006FMSD	1,4-Naphthoquinone	$\mu g/L$	EPA625	12.6	80.00	0	15.8	10-177	30.3	99		12/3/2010 0144h
1011529-006FMSD	1,4-Phenylenediamine	μg/L	EPA625	38.8	80.00	0	48.5	10-124	11.5	48		12/3/2010 0144h
1011529-006FMSD	1-Chloronaphthalene	$\mu g/L$	EPA625	45.1	80.00	0	56.3	10-106	2.11	47		12/3/2010 0144h
1011529-006FMSD	1-Methylnaphthalene	μg/L	EPA625	47.6	80.00	0	59.6	10-83	7.92	44		12/3/2010 0144h
1011529-006FMSD	1-Naphthylamine	$\mug/L$	EPA625	67.7	80.00	0	84.7	10-122	6.62	57		12/3/2010 0144h
1011529-006FMSD	2,3,4,6-Tetrachlorophenol	$\mu g/L$	EPA625	90.7	80.00	0	113	10-157	16.6	39		12/3/2010 0144h
1011529-006FMSD	2,4,5-Trichlorophenol	μg/L	EPA625	65.9	80.00	0	82.4	10-148	9.07	56		12/3/2010 0144h
1011529-006FMSD	2,4,6-Trichlorophenol	$\mu g/L$	EPA625	53.4	80.00	0	66.7	10-136	5.11	52		12/3/2010 0144h
1011529-006FMSD	2,4-Dichlorophenol	$\mug/L$	EPA625	53.8	80.00	0	67.2	10-123	5.58	67		12/3/2010 0144h
1011529-006FMSD	2,4-Dimethylphenol	$\mu g/L$	EPA625	57.2	80.00	0	71.5	10-113	1.18	32		12/3/2010 0144h
1011529-006FMSD	2,4-Dinitrophenol	$\mug/L$	EPA625	73.0	80.00	0	91.2	10-175	10.2	78		12/3/2010 0144h
1011529-006FMSD	2,4-Dinitrotoluene	$\mu g/L$	EPA625	98.4	80.00	0	123	10-175	3.85	26		12/3/2010 0144h
1011529-006FMSD	2,6-Dichlorophenol	$\mug/L$	EPA625	60.0	80.00	0	75.0	10-148	4.31	56		12/3/2010 0144h
1011529-006FMSD	2,6-Dinitrotoluene	$\mug/L$	EPA625	84.2	80.00	0	105	10-175	6.33	74		12/3/2010 0144h
1011529-006FMSD	2-Acetylaminofluorene	$\mug/L$	EPA625	82.4	80.00	0	103	10-94	4.69	52	1	12/3/2010 0144h
1011529-006FMSD	2-Chloronaphthalene	$\mu g/L$	EPA625	49.6	80.00	0	62.0	10-93	16.3	49		12/3/2010 0144h

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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMSD	2-Chlorophenol	μg/L	EPA625	48.7	80.00	0	60.9	10-92	6.60	58		12/3/2010 0144h
1011529-006FMSD	2-Methylnaphthalene	μg/L	EPA625	41.6	80.00	0	52.0	15-78	8.68	39		12/3/2010 0144h
1011529-006FMSD	2-Methylphenol	μg/L	EPA625	48.3	80.00	0	60.4	10-83	3.63	49		12/3/2010 0144h
1011529-006FMSD	2-Naphthylamine	μg/L	EPA625	60.7	80.00	0	75.9	10-154	1.46	37		12/3/2010 0144h
1011529-006FMSD	2-Nitroaniline	μg/L	EPA625	80.6	80.00	0	101	10-175	2.64	87		12/3/2010 0144h
1011529-006FMSD	2-Nitrophenol	μg/L	EPA625	71.5	80.00	0	89.3	10-175	4.05	64		12/3/2010 0144h
1011529-006FMSD	2-Picoline	μg/L	EPA625	26.3	80.00	0	32.9	10-61	16.3	63		12/3/2010 0144h
1011529-006FMSD	3&4-Methylphenol	$\mu$ g/L	EPA625	33.6	80.00	0	42.0	10-80	13.2	99		12/3/2010 0144h
1011529-006FMSD	3,3'-Dichlorobenzidine	μg/L	EPA625	78.5	80.00	0	98.2	10-150	6.98	63		12/3/2010 0144h
1011529-006FMSD	3,3´-Dimethylbenzidine	$\mug/L$	EPA625	81.0	80.00	0	101	10-143	10.0	99		12/3/2010 0144h
1011529-006FMSD	3-Methylcholanthrene	$\mu$ g/L	EPA625	81.8	80.00	0	102	32-171	6.34	46		12/3/2010 0144h
1011529-006FMSD	3-Nitroaniline	$\mug/L$	EPA625	76.7	80.00	0	95.8	10-175	4.03	29		12/3/2010 0144h
1011529-006FMSD	4,6-Dinitro-2-methylphenol	$\mug/L$	EPA625	117	80.00	0	146	10-175	10.4	70		12/3/2010 0144h
1011529-006FMSD	4-Aminobiphenyl	$\mug/L$	EPA625	74.1	80.00	0	92.6	10-175	0.501	79		12/3/2010 0144h
1011529-006FMSD	4-Bromophenyl phenyl ether	$\mu g/L$	EPA625	67.1	80.00	0	83.9	16-138	4.65	31		12/3/2010 0144h
1011529-006FMSD	4-Chloro-3-methylphenol	$\mug/L$	EPA625	65.1	80.00	0	81.4	10-131	2.46	37		12/3/2010 0144h
1011529-006FMSD	4-Chloroaniline	$\mug/L$	EPA625	51.9	80.00	0	64.8	10-98	2.93	41		12/3/2010 0144h
1011529-006FMSD	4-Chlorophenyl phenyl ether	$\mu g/L$	EPA625	63.8	80.00	0	79.7	31-108	5.77	30		12/3/2010 0144h
1011529-006FMSD	4-Nitroaniline	$\mu g/L$	EPA625	71.6	80.00	0	89.5	10-175	4.15	99		12/3/2010 0144h
1011529-006FMSD	4-Nitrophenol	$\mu g/L$	EPA625	128	80.00	0	160	10-97	15.6	69	1	12/3/2010 0144h
1011529-006FMSD	5-Nitro-o-toluidine	$\mu g/L$	EPA625	90.6	80.00	0	113	10-175	4.70	26		12/3/2010 0144h
1011529-006FMSD	7,12-Dimethylbenz(a)anthracene	$\mu g/L$	EPA625	107	80.00	0	133	26-174	7.27	40		12/3/2010 0144h
1011529-006FMSD	a,a-Dimethylphenethylamine	$\mu g/L$	EPA625	41.5	80.00	0	51.8	10-175	8.18	99		12/3/2010 0144h
1011529-006FMSD	Acenaphthene	$\mu g/L$	EPA625	59.1	80.00	0	73.9	29-97	7.73	38		12/3/2010 0144h

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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMSD	Acenaphthylene	μg/L	EPA625	57.8	80.00	0	72.3	37-87	5.62	37		12/3/2010 0144h
1011529-006FMSD	Acetophenone	$\mu g/L$	EPA625	51.4	80.00	0	64.2	10-96	5.05	48		12/3/2010 0144h
1011529-006FMSD	alpha-Terpineol	$\mug/L$	EPA625	50.7	80.00	0	63.4	10-67	1.37	46		12/3/2010 0144h
1011529-006FMSD	Aniline	$\mu g/L$	EPA625	35.8	80.00	0	44.8	10-71	6.91	48		12/3/2010 0144h
1011529-006FMSD	Anthracene	$\mug/L$	EPA625	81.4	80.00	0	102	53-114	6.85	26		12/3/2010 0144h
1011529-006FMSD	Aramite	$\mug/L$	EPA625	89.8	80.00	0	112	29-160	5.00	30		12/3/2010 0144h
1011529-006FMSD	Azobenzene	$\mug/L$	EPA625	61.4	80.00	0	76.7	15-114	7.67	36		12/3/2010 0144h
1011529-006FMSD	Benz(a)anthracene	$\mug/L$	EPA625	75.8	80.00	0	94.8	39-129	4.31	32		12/3/2010 0144h
1011529-006FMSD	Benzidine	$\mug/L$	EPA625	66.5	80.00	0	83.1	10-99	10.7	99		12/3/2010 0144h
1011529-006FMSD	Benzo(a)pyrene	$\mu g/L$	EPA625	85.0	80.00	0	106	29-175	6.66	45		12/3/2010 0144h
1011529-006FMSD	Benzo(b)fluoranthene	$\mu g/L$	EPA625	78.8	80.00	0	98.5	15-140	12.7	44		12/3/2010 0144h
1011529-006FMSD	Benzo(g,h,i)perylene	$\mug/L$	EPA625	69.3	80.00	0	86.7	10-182	6.36	48		12/3/2010 0144h
1011529-006FMSD	Benzo(k)fluoranthene	$\mug/L$	EPA625	100	80.00	0	126	21-154	2.64	52		12/3/2010 0144h
1011529-006FMSD	Benzoic acid	$\mug/L$	EPA625	< 20.0	80.00	0	17.8	10-71	0	78		12/3/2010 0144h
1011529-006FMSD	Benzyl alcohol	$\mug/L$	EPA625	24.6	80.00	0	30.7	10-69	1.68	52		12/3/2010 0144h
1011529-006FMSD	Bis(2-chloroethoxy)methane	$\mu g/L$	EPA625	36.9	80.00	0	46.1	10-94	0.844	45		12/3/2010 0144h
1011529-006FMSD	Bis(2-chloroethyl) ether	$\mu g/L$	EPA625	36.0	80.00	0	45.0	10-70	2.71	47		12/3/2010 0144h
1011529-006FMSD	Bis(2-chloroisopropyl) ether	$\mu g/L$	EPA625	37.5	80.00	0	46.8	10-71	5.63	49		12/3/2010 0144h
1011529-006FMSD	Bis(2-ethylhexyl) phthalate	$\mu g/L$	EPA625	164	80.00	0	204	10-175	57.5	28	1@	12/3/2010 0144h
1011529-006FMSD	bis(2-ethylhexyl)adipate	$\mu g/L$	EPA625	94.3	80.00	0	118	10-175	4.07	25		12/3/2010 0144h
1011529-006FMSD	Butyl benzyl phthalate	$\mug/L$	EPA625	90.3	80.00	0	113	10-175	2.75	99		12/3/2010 0144h
1011529-006FMSD	Carbazole	$\mu g/L$	EPA625	78.3	80.00	0	97.9	10-151	5.93	30		12/3/2010 0144h
1011529-006FMSD	Chlorobenzilate	$\mu g/L$	EPA625	80.8	80.00	0	101	18-175	6.31	25		12/3/2010 0144h
1011529-006FMSD	Chrysene	μg/L	EPA625	81.6	80.00	0	102	38-133	2.53	28		12/3/2010 0144h
										D	D-4 12	/6/2010 Page 110 of 162

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Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMSD	Diallate (cis or trans)	μg/L	EPA625	66.2	80.00	0	82.7	10-157	5.16	29		12/3/2010 0144h
1011529-006FMSD	Dibenz(a,h)anthracene	$\mu g/L$	EPA625	72.4	80.00	0	90.5	13-168	6.37	51		12/3/2010 0144h
1011529-006FMSD	Dibenzofuran	$\mug/L$	EPA625	58.8	80.00	0	73.5	29-103	3.46	34		12/3/2010 0144h
1011529-006FMSD	Diethyl phthalate	$\mu g/L$	EPA625	57.0	80.00	0	71.2	10-139	5.65	32		12/3/2010 0144h
1011529-006FMSD	Dimethoate	$\mug/L$	EPA625	< 10.0	80.00	0	11.9	10-136	0	45		12/3/2010 0144h
1011529-006FMSD	Dimethyl phthalate	$\mug/L$	EPA625	35.1	80.00	0	43.9	10-122	11.2	57		12/3/2010 0144h
1011529-006FMSD	Dimethylaminoazobenzene	$\mug/L$	EPA625	89.8	80.00	0	112	34-142	5.50	26		12/3/2010 0144h
1011529-006FMSD	Di-n-butyl phthalate	$\mug/L$	EPA625	79.7	80.00	0	99.6	44-124	5.93	25		12/3/2010 0144h
1011529-006FMSD	Di-n-octyl phthalate	$\mug/L$	EPA625	143	80.00	0	178	10-175	6.36	50	1	12/3/2010 0144h
1011529-006FMSD	Dinoseb	$\mug/L$	EPA625	128	80.00	0	159	10-175	6.61	42		12/3/2010 0144h
1011529-006FMSD	Diphenylamine	$\mug/L$	EPA625	76.5	80.00	0	95.7	13-110	12.0	34		12/3/2010 0144h
1011529-006FMSD	Disulfoton	$\mug/L$	EPA625	56.9	80.00	0	71.1	10-121	11.7	25		12/3/2010 0144h
1011529-006FMSD	Ethyl methanesulfonate	$\mug/L$	EPA625	46.0	80.00	0	57.4	10-99	3.99	46		12/3/2010 0144h
1011529-006FMSD	Famphur	$\mug/L$	EPA625	139	80.00	0	173	10-71	0.546	25	1	12/3/2010 0144h
1011529-006FMSD	Fluoranthene	$\mug/L$	EPA625	77.3	80.00	0	96.6	23-135	3.17	25		12/3/2010 0144h
1011529-006FMSD	Fluorene	$\mug/L$	EPA625	63.4	80.00	0	79.2	34-108	4.86	28		12/3/2010 0144h
1011529-006FMSD	Hexachlorobenzene	$\mug/L$	EPA625	70.4	80.00	0	88.0	26-131	7.83	28		12/3/2010 0144h
1011529-006FMSD	Hexachlorobutadiene	$\mug/L$	EPA625	23.2	80.00	0	29.0	10-110	18.5	68		12/3/2010 0144h
1011529-006FMSD	Hexachlorocyclopentadiene	$\mug/L$	EPA625	14.7	80.00	0	18.4	10-45	27.4	79		12/3/2010 0144h
1011529-006FMSD	Hexachloroethane	$\mug/L$	EPA625	18.8	80.00	0	23.5	10-58	24.6	42		12/3/2010 0144h
1011529-006FMSD	Hexachlorophene	$\mug/L$	EPA625	89.6	80.00	0	112	10-168	9.60	25		12/3/2010 0144h
1011529-006FMSD	Hexachloropropene	$\mug/L$	EPA625	19.9	80.00	0	24.9	10-72	25.7	63		12/3/2010 0144h
1011529-006FMSD	Indene	$\mug/L$	EPA625	28.2	80.00	0	35.2	10-35	19.7	35	1	12/3/2010 0144h
1011529-006FMSD	Indeno(1,2,3-cd)pyrene	$\mug/L$	EPA625	73.0	80.00	0	91.2	10-176	6.20	48		12/3/2010 0144h
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QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMSD	Isodrin	μg/L	EPA625	76.2	80.00	0	95.2	15-165	8.94	25		12/3/2010 0144h
1011529-006FMSD	Isophorone	μg/L	EPA625	50.0	80.00	0	62.6	10-99	1.01	41		12/3/2010 0144h
1011529-006FMSD	Isosafrole	$\mu g/L$	EPA625	47.8	80.00	0	59.8	10-167	9.18	50		12/3/2010 0144h
1011529-006FMSD	Kepone	μg/L	EPA625	215	80.00	0	269	10-175	1.01	46	1	12/3/2010 0144h
1011529-006FMSD	Methapyrilene	$\mug/L$	EPA625	86.3	80.00	0	108	10-149	1.48	51		12/3/2010 0144h
1011529-006FMSD	Methyl methanesulfonate	$\mug/L$	EPA625	30.8	80.00	0	38.6	10-132	3.70	45		12/3/2010 0144h
1011529-006FMSD	Naphthalene	$\mug/L$	EPA625	35.3	80.00	0	44.1	10-82	14.6	43		12/3/2010 0144h
1011529-006FMSD	n-Decane	$\mug/L$	EPA625	12.8	80.00	0	16.0	10-27	21.1	32		12/3/2010 0144h
1011529-006FMSD	Nitrobenzene	$\mug/L$	EPA625	59.5	80.00	0	74.4	10-119	5.96	49		12/3/2010 0144h
1011529-006FMSD	Nitroquinoline-1-oxide	$\mug/L$	EPA625	36.1	80.00	0	45.1	10-170	14.9	97		12/3/2010 0144h
1011529-006FMSD	N-Nitrosodiethylamine	$\mug/L$	EPA625	43.4	80.00	0	54.3	10-91	5.08	54		12/3/2010 0144h
1011529-006FMSD	N-Nitrosodimethylamine	$\mug/L$	EPA625	19.1	80.00	0	23.9	10-42	8.05	66		12/3/2010 0144h
1011529-006FMSD	N-Nitrosodi-n-butylamine	$\mug/L$	EPA625	57.2	80.00	0	71.5	10-175	1.91	80		12/3/2010 0144h
1011529-006FMSD	N-Nitrosodiphenylamine	$\mug/L$	EPA625	75.2	80.00	0	94.0	12-112	11.5	30		12/3/2010 0144h
1011529-006FMSD	N-Nitrosodi-n-propylamine	$\mug/L$	EPA625	47.1	80.00	0	58.9	10-77	3.80	47		12/3/2010 0144h
1011529-006FMSD	N-Nitrosomethylethylamine	$\mug/L$	EPA625	33.4	80.00	0	41.8	10-75	0.781	60		12/3/2010 0144h
1011529-006FMSD	N-Nitrosomorpholine	$\mug/L$	EPA625	47.8	80.00	0	59.8	10-175	0.313	73		12/3/2010 0144h
1011529-006FMSD	N-Nitrosopiperidine	$\mug/L$	EPA625	55.9	80.00	0	69.8	10-105	3.11	44		12/3/2010 0144h
1011529-006FMSD	N-Nitrosopyrrolidine	$\mug/L$	EPA625	56.7	80.00	0	70.9	10-88	4.28	40		12/3/2010 0144h
1011529-006FMSD	n-Octadecane	$\mug/L$	EPA625	59.2	80.00	0	74.0	10-121	2.81	40		12/3/2010 0144h
1011529-006FMSD	O,O,O-Triethyl phosphorothioate	$\mug/L$	EPA625	53.7	80.00	0	67.1	10-93	4.50	44		12/3/2010 0144h
1011529-006FMSD	o-Toluidine	$\mug/L$	EPA625	95.2	80.00	0	119	10-107	1.22	46	1	12/3/2010 0144h
1011529-006FMSD	Parathion	$\mug/L$	EPA625	111	80.00	0	139	10-175	5.99	28		12/3/2010 0144h
1011529-006FMSD	Methyl parathion	$\mug/L$	EPA625	102	80.00	0	128	10-175	5.65	26		12/3/2010 0144h
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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006FMSD	Pentachlorobenzene	μg/L	EPA625	58.5	80.00	0	73.2	25-134	4.99	35		12/3/2010 0144h
1011529-006FMSD	Pentachloronitrobenzene	$\mug/L$	EPA625	81.0	80.00	0	101	10-175	4.27	25		12/3/2010 0144h
1011529-006FMSD	Pentachlorophenol	$\mug/L$	EPA625	64.0	80.00	0	80.0	10-163	26.3	58		12/3/2010 0144h
1011529-006FMSD	Phenacetin	$\mug/L$	EPA625	77.3	80.00	0	96.6	10-175	6.69	38		12/3/2010 0144h
1011529-006FMSD	Phenanthrene	$\mug/L$	EPA625	83.0	80.00	0	104	31-126	5.53	32		12/3/2010 0144h
1011529-006FMSD	Phenol	$\mug/L$	EPA625	29.8	80.00	0	37.3	10-175	6.08	71		12/3/2010 0144h
1011529-006FMSD	Phorate	$\mug/L$	EPA625	81.0	80.00	0	101	10-175	8.08	42		12/3/2010 0144h
1011529-006FMSD	Pronamide	$\mug/L$	EPA625	79.6	80.00	0	99.5	10-95	3.18	26	1	12/3/2010 0144h
1011529-006FMSD	Pyrene	$\mug/L$	EPA625	91.5	80.00	0	114	51-139	4.01	27		12/3/2010 0144h
1011529-006FMSD	Pyridine	$\mug/L$	EPA625	11.6	80.00	0	14.5	10-25	0.601	61		12/3/2010 0144h
1011529-006FMSD	Quinoline	$\mug/L$	EPA625	59.5	80.00	0	74.4	10-63	2.95	99	1	12/3/2010 0144h
1011529-006FMSD	Safrole	$\mug/L$	EPA625	51.5	80.00	0	64.4	10-120	6.59	51		12/3/2010 0144h
1011529-006FMSD	Tetraethyl dithiopyrophosphate	$\mug/L$	EPA625	73.6	80.00	0	92.0	13-160	6.27	35		12/3/2010 0144h
1011529-006FMSD	Thionazin	$\mug/L$	EPA625	66.2	80.00	0	82.8	10-139	6.93	25		12/3/2010 0144h
1011529-006FMSD	Surr: 2,4,6-Tribromophenol	%REC	EPA625	68.2	80.00		85.2	21-154				12/3/2010 0144h
1011529-006FMSD	Surr: 2-Fluorobiphenyl	%REC	EPA625	49.3	40.00		123	10-106			S	12/3/2010 0144h
1011529-006FMSD	Surr: 2-Fluorophenol	%REC	EPA625	26.8	80.00		33.6	10-56				12/3/2010 0144h
1011529-006FMSD	Surr: Nitrobenzene-d5	%REC	EPA625	53.9	40.00		135	10-101			S	12/3/2010 0144h
1011529-006FMSD	Surr: Phenol-d6	%REC	EPA625	19.7	80.00		24.6	10-45				12/3/2010 0144h
1011529-006FMSD	Surr: Terphenyl-d14	%REC	EPA625	92.3	40.00		231	10-160			S	12/3/2010 0144h
1012039-001DMSD	1,1´-Biphenyl	$\mu g/L$	EPA625	64.2	83.33	0	77.1	27-99	0.183	49		12/3/2010 0959h
1012039-001DMSD	1,2,4,5-Tetrachlorobenzene	$\mug/L$	EPA625	58.8	83.33	0	70.5	10-119	0.302	52		12/3/2010 0959h
1012039-001DMSD	1,2,4-Trichlorobenzene	$\mug/L$	EPA625	44.5	83.33	0	53.4	10-79	4.00	49		12/3/2010 0959h
1012039-001DMSD	1,2-Dichlorobenzene	$\mu g/L$	EPA625	32.6	83.33	0	39.1	10-59	4.05	46		12/3/2010 0959h
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Jose Rocha

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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMSD	1,3,5-Trinitrobenzene	μg/L	EPA625	159	83.33	0	191	10-175	6.83	33	1	12/3/2010 0959h
1012039-001DMSD	1,3-Dichlorobenzene	μg/L	EPA625	25.7	83.33	0	30.9	10-56	10.2	49		12/3/2010 0959h
1012039-001DMSD	1,3-Dinitrobenzene	μg/L	EPA625	108	83.33	0	130	10-175	13.1	29		12/3/2010 0959h
1012039-001DMSD	1,4-Dichlorobenzene	μg/L	EPA625	27.5	83.33	0	33.0	10-58	10.6	51		12/3/2010 0959h
1012039-001DMSD	1,4-Naphthoquinone	μg/L	EPA625	19.9	83.33	0	23.9	10-177	19.6	99		12/3/2010 0959h
1012039-001DMSD	1,4-Phenylenediamine	$\mu g/L$	EPA625	51.9	83.33	0	62.3	10-124	0.201	48		12/3/2010 0959h
1012039-001DMSD	1-Chloronaphthalene	μg/L	EPA625	67.4	83.33	0	80.9	10-106	20.0	47		12/3/2010 0959h
1012039-001DMSD	1-Methylnaphthalene	μg/L	EPA625	65.2	83.33	0	78.2	10-83	0.758	44		12/3/2010 0959h
1012039-001DMSD	1-Naphthylamine	μg/L	EPA625	71.6	83.33	0	85.9	10-122	2.11	57		12/3/2010 0959h
1012039-001DMSD	2,3,4,6-Tetrachlorophenol	μg/L	EPA625	87.7	83.33	0	105	10-157	11.8	39		12/3/2010 0959h
1012039-001DMSD	2,4,5-Trichlorophenol	μg/L	EPA625	75.2	83.33	0	90.3	10-148	12.7	56		12/3/2010 0959h
1012039-001DMSD	2,4,6-Trichlorophenol	$\mu g/L$	EPA625	60.2	83.33	0	72.3	10-136	11.8	52		12/3/2010 0959h
1012039-001DMSD	2,4-Dichlorophenol	$\mu g/L$	EPA625	60.4	83.33	0	72.5	10-123	6.16	67		12/3/2010 0959h
1012039-001DMSD	2,4-Dimethylphenol	$\mu g/L$	EPA625	62.9	83.33	0	75.5	10-113	1.05	32		12/3/2010 0959h
1012039-001DMSD	2,4-Dinitrophenol	μg/L	EPA625	75.0	83.33	0	90.0	10-175	27.1	78		12/3/2010 0959h
1012039-001DMSD	2,4-Dinitrotoluene	μg/L	EPA625	105	83.33	0	126	10-175	5.54	26		12/3/2010 0959h
1012039-001DMSD	2,6-Dichlorophenol	μg/L	EPA625	70.2	83.33	0	84.3	10-148	6.04	56		12/3/2010 0959h
1012039-001DMSD	2,6-Dinitrotoluene	$\mu g/L$	EPA625	88.7	83.33	0	106	10-175	0.304	74		12/3/2010 0959h
1012039-001DMSD	2-Acetylaminofluorene	$\mu g/L$	EPA625	78.0	83.33	0	93.6	10-94	5.43	52		12/3/2010 0959h
1012039-001DMSD	2-Chloronaphthalene	μg/L	EPA625	56.6	83.33	0	67.9	10-93	13.3	49		12/3/2010 0959h
1012039-001DMSD	2-Chlorophenol	$\mu g/L$	EPA625	57.1	83.33	0	68.5	10-92	5.61	58		12/3/2010 0959h
1012039-001DMSD	2-Methylnaphthalene	$\mug/L$	EPA625	55.7	83.33	0	66.8	15-78	1.25	39		12/3/2010 0959h
1012039-001DMSD	2-Methylphenol	$\mug/L$	EPA625	53.3	83.33	0	63.9	10-83	6.62	49		12/3/2010 0959h
1012039-001DMSD	2-Naphthylamine	$\mug/L$	EPA625	67.4	83.33	0	80.8	10-154	4.39	37		12/3/2010 0959h
										Danart	Doto: 12	/6/2010 Page 123 of 16

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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMSD	2-Nitroaniline	μg/L	EPA625	89.1	83.33	0	107	10-175	6.12	87		12/3/2010 0959h
1012039-001DMSD	2-Nitrophenol	μg/L	EPA625	83.2	83.33	0	99.8	10-175	3.14	64		12/3/2010 0959h
1012039-001DMSD	2-Picoline	$\mug/L$	EPA625	37.9	83.33	0	45.5	10-61	2.06	63		12/3/2010 0959h
1012039-001DMSD	3&4-Methylphenol	$\mug/L$	EPA625	30.8	83.33	0	36.9	10-80	13.5	99		12/3/2010 0959h
1012039-001DMSD	3,3´-Dichlorobenzidine	$\mug/L$	EPA625	71.2	83.33	0	85.5	10-150	5.97	63		12/3/2010 0959h
1012039-001DMSD	3,3´-Dimethylbenzidine	$\mug/L$	EPA625	80.9	83.33	0	97.1	10-143	12.8	99		12/3/2010 0959h
1012039-001DMSD	3-Methylcholanthrene	$\mug/L$	EPA625	72.0	83.33	0	86.4	32-171	2.77	46		12/3/2010 0959h
1012039-001DMSD	3-Nitroaniline	$\mug/L$	EPA625	83.1	83.33	0	99.8	10-175	6.27	29		12/3/2010 0959h
1012039-001DMSD	4,6-Dinitro-2-methylphenol	$\mug/L$	EPA625	115	83.33	0	138	10-175	2.73	70		12/3/2010 0959h
1012039-001DMSD	4-Aminobiphenyl	$\mug/L$	EPA625	82.3	83.33	0	98.7	10-175	4.00	79		12/3/2010 0959h
1012039-001DMSD	4-Bromophenyl phenyl ether	$\mug/L$	EPA625	74.7	83.33	0	89.7	16-138	1.40	31		12/3/2010 0959h
1012039-001DMSD	4-Chloro-3-methylphenol	$\mug/L$	EPA625	70.9	83.33	0	85.0	10-131	5.70	37		12/3/2010 0959h
1012039-001DMSD	4-Chloroaniline	$\mug/L$	EPA625	59.4	83.33	0	71.3	10-98	4.42	41		12/3/2010 0959h
1012039-001DMSD	4-Chlorophenyl phenyl ether	$\mug/L$	EPA625	70.9	83.33	0	85.1	31-108	2.41	30		12/3/2010 0959h
1012039-001DMSD	4-Nitroaniline	$\mug/L$	EPA625	73.3	83.33	0	88.0	10-175	8.71	99		12/3/2010 0959h
1012039-001DMSD	4-Nitrophenol	$\mug/L$	EPA625	148	83.33	0	177	10-97	20.1	69	1	12/3/2010 0959h
1012039-001DMSD	5-Nitro-o-toluidine	$\mug/L$	EPA625	93.4	83.33	0	112	10-175	3.39	26		12/3/2010 0959h
1012039-001DMSD	7,12-Dimethylbenz(a)anthracene	$\mug/L$	EPA625	97.1	83.33	0	117	26-174	2.79	40		12/3/2010 0959h
1012039-001DMSD	a,a-Dimethylphenethylamine	$\mug/L$	EPA625	47.2	83.33	0	56.6	10-175	11.5	99		12/3/2010 0959h
1012039-001DMSD	Acenaphthene	$\mug/L$	EPA625	67.6	83.33	0	81.2	29-97	0.972	38		12/3/2010 0959h
1012039-001DMSD	Acenaphthylene	$\mug/L$	EPA625	69.1	83.33	0	82.9	37-87	2.41	37		12/3/2010 0959h
1012039-001DMSD	Acetophenone	$\mug/L$	EPA625	57.4	83.33	0	68.9	10-96	0.816	48		12/3/2010 0959h
1012039-001DMSD	alpha-Terpineol	$\mug/L$	EPA625	57.1	83.33	0	68.5	10-67	2.03	46	1	12/3/2010 0959h
1012039-001DMSD	Aniline	$\mug/L$	EPA625	46.6	83.33	0	55.9	10-71	1.93	48		12/3/2010 0959h
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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMSD	Anthracene	μg/L	EPA625	84.9	83.33	0	102	53-114	1.99	26		12/3/2010 0959h
1012039-001DMSD	Aramite	$\mu g/L$	EPA625	92.0	83.33	0	110	29-160	1.47	30		12/3/2010 0959h
1012039-001DMSD	Azobenzene	$\mug/L$	EPA625	67.3	83.33	0	80.7	15-114	1.94	36		12/3/2010 0959h
1012039-001DMSD	Benz(a)anthracene	$\mug/L$	EPA625	70.2	83.33	0	84.2	39-129	6.43	32		12/3/2010 0959h
1012039-001DMSD	Benzidine	$\mug/L$	EPA625	65.9	83.33	0	79.1	10-99	10.2	99		12/3/2010 0959h
1012039-001DMSD	Benzo(a)pyrene	$\mug/L$	EPA625	76.3	83.33	0	91.6	29-175	3.31	45		12/3/2010 0959h
1012039-001DMSD	Benzo(b)fluoranthene	$\mug/L$	EPA625	65.8	83.33	0	79.0	15-140	14.8	44		12/3/2010 0959h
1012039-001DMSD	Benzo(g,h,i)perylene	$\mug/L$	EPA625	61.8	83.33	0	74.2	10-182	1.12	48		12/3/2010 0959h
1012039-001DMSD	Benzo(k)fluoranthene	$\mug/L$	EPA625	93.6	83.33	0	112	21-154	9.10	52		12/3/2010 0959h
1012039-001DMSD	Benzoic acid	$\mug/L$	EPA625	< 20.8	83.33	0	13.5	10-71	0	78		12/3/2010 0959h
1012039-001DMSD	Benzyl alcohol	$\mug/L$	EPA625	28.7	83.33	0	34.4	10-69	10.7	52		12/3/2010 0959h
1012039-001DMSD	Bis(2-chloroethoxy)methane	$\mug/L$	EPA625	42.7	83.33	0	51.2	10-94	3.46	45		12/3/2010 0959h
1012039-001DMSD	Bis(2-chloroethyl) ether	$\mug/L$	EPA625	46.5	83.33	0	55.8	10-70	12.8	47		12/3/2010 0959h
1012039-001DMSD	Bis(2-chloroisopropyl) ether	$\mug/L$	EPA625	43.3	83.33	0	52.0	10-71	0.689	49		12/3/2010 0959h
1012039-001DMSD	Bis(2-ethylhexyl) phthalate	$\mug/L$	EPA625	85.3	83.33	0	102	10-175	0.187	28		12/3/2010 0959h
1012039-001DMSD	bis(2-ethylhexyl)adipate	$\mug/L$	EPA625	88.4	83.33	0	106	10-175	3.20	25		12/3/2010 0959h
1012039-001DMSD	Butyl benzyl phthalate	$\mug/L$	EPA625	86.8	83.33	0	104	10-175	2.15	99		12/3/2010 0959h
1012039-001DMSD	Carbazole	$\mug/L$	EPA625	79.2	83.33	0	95.1	10-151	2.88	30		12/3/2010 0959h
1012039-001DMSD	Chlorobenzilate	$\mug/L$	EPA625	84.3	83.33	0	101	18-175	3.44	25		12/3/2010 0959h
1012039-001DMSD	Chrysene	$\mug/L$	EPA625	77.4	83.33	0	92.8	38-133	5.19	28		12/3/2010 0959h
1012039-001DMSD	Diallate (cis or trans)	$\mug/L$	EPA625	71.6	83.33	0	85.9	10-157	0.703	29		12/3/2010 0959h
1012039-001DMSD	Dibenz(a,h)anthracene	$\mug/L$	EPA625	63.5	83.33	0	76.2	13-168	2.79	51		12/3/2010 0959h
1012039-001DMSD	Dibenzofuran	$\mug/L$	EPA625	69.3	83.33	0	83.2	29-103	3.91	34		12/3/2010 0959h
1012039-001DMSD	Diethyl phthalate	$\mug/L$	EPA625	72.8	83.33	0	87.4	10-139	0.0783	32		12/3/2010 0959h
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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMSD	Dimethoate	μg/L	EPA625	32.1	83.33	0	38.6	10-136	4.94	45		12/3/2010 0959h
1012039-001DMSD	Dimethyl phthalate	$\mug/L$	EPA625	64.5	83.33	0	77.4	10-122	1.26	57		12/3/2010 0959h
1012039-001DMSD	Dimethylaminoazobenzene	$\mu g/L$	EPA625	83.0	83.33	0	99.6	34-142	2.96	26		12/3/2010 0959h
1012039-001DMSD	Di-n-butyl phthalate	$\mug/L$	EPA625	82.0	83.33	0	98.4	44-124	1.04	25		12/3/2010 0959h
1012039-001DMSD	Di-n-octyl phthalate	$\mug/L$	EPA625	124	83.33	0	149	10-175	1.28	50		12/3/2010 0959h
1012039-001DMSD	Dinoseb	$\mug/L$	EPA625	135	83.33	0	162	10-175	7.70	42		12/3/2010 0959h
1012039-001DMSD	Diphenylamine	$\mug/L$	EPA625	78.0	83.33	0	93.6	13-110	1.53	34		12/3/2010 0959h
1012039-001DMSD	Disulfoton	$\mug/L$	EPA625	59.1	83.33	0	71.0	10-121	13.3	25		12/3/2010 0959h
1012039-001DMSD	Ethyl methanesulfonate	$\mug/L$	EPA625	54.1	83.33	0	64.9	10-99	4.85	46		12/3/2010 0959h
1012039-001DMSD	Famphur	$\mug/L$	EPA625	123	83.33	0	148	10-71	0.671	25	1	12/3/2010 0959h
1012039-001DMSD	Fluoranthene	$\mug/L$	EPA625	78.7	83.33	0	94.5	23-135	3.14	25		12/3/2010 0959h
1012039-001DMSD	Fluorene	$\mug/L$	EPA625	70.9	83.33	0	85.0	34-108	0.140	28		12/3/2010 0959h
1012039-001DMSD	Hexachlorobenzene	$\mug/L$	EPA625	75.2	83.33	0	90.3	26-131	0.496	28		12/3/2010 0959h
1012039-001DMSD	Hexachlorobutadiene	$\mug/L$	EPA625	39.2	83.33	0	47.1	10-110	5.79	68		12/3/2010 0959h
1012039-001DMSD	Hexachlorocyclopentadiene	$\mug/L$	EPA625	28.0	83.33	0	33.6	10-45	2.73	79		12/3/2010 0959h
1012039-001DMSD	Hexachloroethane	$\mug/L$	EPA625	26.3	83.33	0	31.6	10-58	13.7	42		12/3/2010 0959h
1012039-001DMSD	Hexachlorophene	$\mug/L$	EPA625	84.6	83.33	0	101	10-168	11.9	25		12/3/2010 0959h
1012039-001DMSD	Hexachloropropene	$\mu g/L$	EPA625	35.3	83.33	0	42.4	10-72	7.54	63		12/3/2010 0959h
1012039-001DMSD	Indene	$\mug/L$	EPA625	38.3	83.33	0	46.0	10-35	6.28	35	1	12/3/2010 0959h
1012039-001DMSD	Indeno(1,2,3-cd)pyrene	$\mug/L$	EPA625	65.1	83.33	0	78.1	10-176	2.58	48		12/3/2010 0959h
1012039-001DMSD	Isodrin	$\mug/L$	EPA625	77.3	83.33	0	92.7	15-165	0.537	25		12/3/2010 0959h
1012039-001DMSD	Isophorone	$\mug/L$	EPA625	58.0	83.33	0	69.6	10-99	0.281	41		12/3/2010 0959h
1012039-001DMSD	Isosafrole	$\mug/L$	EPA625	58.8	83.33	0	70.6	10-167	1.64	50		12/3/2010 0959h
1012039-001DMSD	Kepone	$\mug/L$	EPA625	217	83.33	0	261	10-175	5.68	46	1	12/3/2010 0959h
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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMSD	Methapyrilene	μg/L	EPA625	90.9	83.33	0	109	10-149	2.76	51		12/3/2010 0959h
1012039-001DMSD	Methyl methanesulfonate	$\mug/L$	EPA625	38.9	83.33	0	46.6	10-132	7.74	45		12/3/2010 0959h
1012039-001DMSD	Naphthalene	$\mug/L$	EPA625	51.0	83.33	0	61.2	10-82	3.96	43		12/3/2010 0959h
1012039-001DMSD	n-Decane	$\mug/L$	EPA625	15.5	83.33	0	18.6	10-27	21.4	32		12/3/2010 0959h
1012039-001DMSD	Nitrobenzene	$\mug/L$	EPA625	68.1	83.33	0	81.8	10-119	0.530	49		12/3/2010 0959h
1012039-001DMSD	Nitroquinoline-1-oxide	$\mug/L$	EPA625	36.8	83.33	0	44.2	10-170	19.4	97		12/3/2010 0959h
1012039-001DMSD	N-Nitrosodiethylamine	$\mug/L$	EPA625	51.5	83.33	0	61.8	10-91	3.04	54		12/3/2010 0959h
1012039-001DMSD	N-Nitrosodimethylamine	$\mug/L$	EPA625	21.6	83.33	0	25.9	10-42	7.11	66		12/3/2010 0959h
1012039-001DMSD	N-Nitrosodi-n-butylamine	$\mug/L$	EPA625	64.5	83.33	0	77.4	10-175	1.24	80		12/3/2010 0959h
1012039-001DMSD	N-Nitrosodiphenylamine	$\mug/L$	EPA625	77.1	83.33	0	92.5	12-112	1.87	30		12/3/2010 0959h
1012039-001DMSD	N-Nitrosodi-n-propylamine	$\mug/L$	EPA625	54.2	83.33	0	65.1	10-77	0.613	47		12/3/2010 0959h
1012039-001DMSD	N-Nitrosomethylethylamine	$\mug/L$	EPA625	43.3	83.33	0	52.0	10-75	8.36	60		12/3/2010 0959h
1012039-001DMSD	N-Nitrosomorpholine	$\mug/L$	EPA625	58.9	83.33	0	70.6	10-175	5.41	73		12/3/2010 0959h
1012039-001DMSD	N-Nitrosopiperidine	$\mug/L$	EPA625	63.0	83.33	0	75.6	10-105	2.81	44		12/3/2010 0959h
1012039-001DMSD	N-Nitrosopyrrolidine	$\mug/L$	EPA625	65.7	83.33	0	78.9	10-88	3.74	40		12/3/2010 0959h
1012039-001DMSD	n-Octadecane	$\mug/L$	EPA625	68.0	83.33	0	81.6	10-121	0.0591	40		12/3/2010 0959h
1012039-001DMSD	O,O,O-Triethyl phosphorothioate	$\mug/L$	EPA625	60.5	83.33	0	72.6	10-93	5.35	44		12/3/2010 0959h
1012039-001DMSD	o-Toluidine	$\mug/L$	EPA625	121	83.33	0	145	10-107	5.43	46	1	12/3/2010 0959h
1012039-001DMSD	Parathion	$\mug/L$	EPA625	111	83.33	0	134	10-175	1.55	28		12/3/2010 0959h
1012039-001DMSD	Methyl parathion	$\mug/L$	EPA625	103	83.33	0	123	10-175	2.57	26		12/3/2010 0959h
1012039-001DMSD	Pentachlorobenzene	$\mug/L$	EPA625	67.7	83.33	0	81.3	25-134	1.88	35		12/3/2010 0959h
1012039-001DMSD	Pentachloronitrobenzene	$\mug/L$	EPA625	84.8	83.33	0	102	10-175	1.38	25		12/3/2010 0959h
1012039-001DMSD	Pentachlorophenol	$\mug/L$	EPA625	47.5	83.33	0	57.0	10-163	7.31	58		12/3/2010 0959h
1012039-001DMSD	Phenacetin	$\mug/L$	EPA625	78.0	83.33	0	93.6	10-175	7.89	38		12/3/2010 0959h
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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMSD	Phenanthrene	μg/L	EPA625	88.2	83.33	0	106	31-126	0.433	32		12/3/2010 0959h
1012039-001DMSD	Phenol	$\mug/L$	EPA625	33.7	83.33	0	40.5	10-175	3.30	71		12/3/2010 0959h
1012039-001DMSD	Phorate	$\mug/L$	EPA625	86.5	83.33	0	104	10-175	3.13	42		12/3/2010 0959h
1012039-001DMSD	Pronamide	$\mug/L$	EPA625	85.7	83.33	0	103	10-95	8.30	26	1	12/3/2010 0959h
1012039-001DMSD	Pyrene	$\mug/L$	EPA625	86.5	83.33	0	104	51-139	2.39	27		12/3/2010 0959h
1012039-001DMSD	Pyridine	$\mug/L$	EPA625	21.9	83.33	0	26.2	10-25	4.47	61	1	12/3/2010 0959h
1012039-001DMSD	Quinoline	$\mug/L$	EPA625	67.3	83.33	0	80.8	10-63	3.53	99	1	12/3/2010 0959h
1012039-001DMSD	Safrole	$\mug/L$	EPA625	60.7	83.33	0	72.9	10-120	3.53	51		12/3/2010 0959h
1012039-001DMSD	Tetraethyl dithiopyrophosphate	$\mug/L$	EPA625	77.3	83.33	0	92.8	13-160	0.833	35		12/3/2010 0959h
1012039-001DMSD	Thionazin	$\mug/L$	EPA625	71.5	83.33	0	85.8	10-139	2.92	25		12/3/2010 0959h
1012039-001DMSD	Surr: 2,4,6-Tribromophenol	%REC	EPA625	70.6	83.33		84.7	21-154				12/3/2010 0959h
1012039-001DMSD	Surr: 2-Fluorobiphenyl	%REC	EPA625	63.1	41.67		152	10-106			S	12/3/2010 0959h
1012039-001DMSD	Surr: 2-Fluorophenol	%REC	EPA625	30.6	83.33		36.8	10-56				12/3/2010 0959h
1012039-001DMSD	Surr: Nitrobenzene-d5	%REC	EPA625	60.8	41.67		146	10-101			S	12/3/2010 0959h
1012039-001DMSD	Surr: Phenol-d6	%REC	EPA625	20.9	83.33		25.1	10-45				12/3/2010 0959h
1012039-001DMSD	Surr: Terphenyl-d14	%REC	EPA625	88.3	41.67		212	10-160			S	12/3/2010 0959h
1012057-001BMSD	1,1´-Biphenyl	$\mug/L$	EPA625	65.5	80.00	0	81.9	27-99	23.3	49		12/3/2010 1149h
1012057-001BMSD	1,2,4,5-Tetrachlorobenzene	$\mug/L$	EPA625	63.2	80.00	0	79.0	10-119	28.7	52		12/3/2010 1149h
1012057-001BMSD	1,2,4-Trichlorobenzene	$\mug/L$	EPA625	47.4	80.00	0	59.2	10-79	22.5	49		12/3/2010 1149h
1012057-001BMSD	1,2-Dichlorobenzene	$\mug/L$	EPA625	36.0	80.00	0	45.0	10-59	20.7	46		12/3/2010 1149h
1012057-001BMSD	1,3,5-Trinitrobenzene	$\mug/L$	EPA625	155	80.00	0	193	10-175	14.4	33	1	12/3/2010 1149h
1012057-001BMSD	1,3-Dichlorobenzene	$\mug/L$	EPA625	29.5	80.00	0	36.9	10-56	18.2	49		12/3/2010 1149h
1012057-001BMSD	1,3-Dinitrobenzene	$\mug/L$	EPA625	107	80.00	0	134	10-175	34.0	29	@	12/3/2010 1149h
1012057-001BMSD	1,4-Dichlorobenzene	$\mug/L$	EPA625	31.7	80.00	0	39.6	10-58	21.7	51		12/3/2010 1149h
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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMSD	1,4-Naphthoquinone	μg/L	EPA625	17.2	80.00	0	21.5	10-177	17.4	99		12/3/2010 1149h
1012057-001BMSD	1,4-Phenylenediamine	μg/L	EPA625	54.6	80.00	0	68.3	10-124	26.9	48		12/3/2010 1149h
1012057-001BMSD	1-Chloronaphthalene	μg/L	EPA625	64.8	80.00	0	81.0	10-106	41.2	47		12/3/2010 1149h
1012057-001BMSD	1-Methylnaphthalene	μg/L	EPA625	66.8	80.00	0	83.5	10-83	23.4	44	1	12/3/2010 1149h
1012057-001BMSD	1-Naphthylamine	μg/L	EPA625	70.8	80.00	0	88.5	10-122	17.2	57		12/3/2010 1149h
1012057-001BMSD	2,3,4,6-Tetrachlorophenol	μg/L	EPA625	84.8	80.00	0	106	10-157	6.87	39		12/3/2010 1149h
1012057-001BMSD	2,4,5-Trichlorophenol	μg/L	EPA625	71.6	80.00	0	89.6	10-148	18.0	56		12/3/2010 1149h
1012057-001BMSD	2,4,6-Trichlorophenol	μg/L	EPA625	57.2	80.00	0	71.6	10-136	19.6	52		12/3/2010 1149h
1012057-001BMSD	2,4-Dichlorophenol	μg/L	EPA625	58.1	80.00	0	72.6	10-123	17.0	67		12/3/2010 1149h
1012057-001BMSD	2,4-Dimethylphenol	μg/L	EPA625	63.7	80.00	0	79.6	10-113	14.1	32		12/3/2010 1149h
1012057-001BMSD	2,4-Dinitrophenol	μg/L	EPA625	73.7	80.00	0	92.1	10-175	37.9	78		12/3/2010 1149h
1012057-001BMSD	2,4-Dinitrotoluene	μg/L	EPA625	103	80.00	0	128	10-175	17.2	26		12/3/2010 1149h
1012057-001BMSD	2,6-Dichlorophenol	μg/L	EPA625	67.8	80.00	0	84.8	10-148	11.5	56		12/3/2010 1149h
1012057-001BMSD	2,6-Dinitrotoluene	μg/L	EPA625	87.3	80.00	0	109	10-175	15.9	74		12/3/2010 1149h
1012057-001BMSD	2-Acetylaminofluorene	μg/L	EPA625	79.2	80.00	0	99.0	10-94	14.0	52	1	12/3/2010 1149h
1012057-001BMSD	2-Chloronaphthalene	μg/L	EPA625	59.3	80.00	0	74.1	10-93	24.7	49		12/3/2010 1149h
1012057-001BMSD	2-Chlorophenol	μg/L	EPA625	55.1	80.00	0	68.9	10-92	20.0	58		12/3/2010 1149h
1012057-001BMSD	2-Methylnaphthalene	μg/L	EPA625	55.8	80.00	0	69.7	15-78	20.7	39		12/3/2010 1149h
1012057-001BMSD	2-Methylphenol	μg/L	EPA625	44.7	80.00	0	55.8	10-83	19.9	49		12/3/2010 1149h
1012057-001BMSD	2-Naphthylamine	μg/L	EPA625	64.4	80.00	0	80.5	10-154	17.5	37		12/3/2010 1149h
1012057-001BMSD	2-Nitroaniline	μg/L	EPA625	87.8	80.00	0	110	10-175	19.6	87		12/3/2010 1149h
1012057-001BMSD	2-Nitrophenol	μg/L	EPA625	82.1	80.00	0	103	10-175	9.21	64		12/3/2010 1149h
1012057-001BMSD	2-Picoline	μg/L	EPA625	40.8	80.00	0	51.0	10-61	51.0	63		12/3/2010 1149h
1012057-001BMSD	3&4-Methylphenol	μg/L	EPA625	26.9	80.00	0	33.6	10-80	29.9	99		12/3/2010 1149h

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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMSD	3,3´-Dichlorobenzidine	μg/L	EPA625	73.6	80.00	0	92.0	10-150	15.6	63		12/3/2010 1149h
1012057-001BMSD	3,3'-Dimethylbenzidine	μg/L	EPA625	82.1	80.00	0	103	10-143	13.5	99		12/3/2010 1149h
1012057-001BMSD	3-Methylcholanthrene	$\mu g/L$	EPA625	78.3	80.00	0	97.8	32-171	16.4	46		12/3/2010 1149h
1012057-001BMSD	3-Nitroaniline	μg/L	EPA625	80.6	80.00	0	101	10-175	17.4	29		12/3/2010 1149h
1012057-001BMSD	4,6-Dinitro-2-methylphenol	$\mu g/L$	EPA625	119	80.00	0	149	10-175	15.8	70		12/3/2010 1149h
1012057-001BMSD	4-Aminobiphenyl	$\mug/L$	EPA625	78.6	80.00	0	98.3	10-175	15.9	79		12/3/2010 1149h
1012057-001BMSD	4-Bromophenyl phenyl ether	$\mu g/L$	EPA625	70.5	80.00	0	88.2	16-138	17.0	31		12/3/2010 1149h
1012057-001BMSD	4-Chloro-3-methylphenol	$\mu g/L$	EPA625	65.1	80.00	0	81.4	10-131	13.5	37		12/3/2010 1149h
1012057-001BMSD	4-Chloroaniline	$\mu g/L$	EPA625	59.3	80.00	0	74.1	10-98	22.2	41		12/3/2010 1149h
1012057-001BMSD	4-Chlorophenyl phenyl ether	$\mu g/L$	EPA625	70.9	80.00	0	88.6	31-108	20.6	30		12/3/2010 1149h
1012057-001BMSD	4-Nitroaniline	$\mug/L$	EPA625	71.7	80.00	0	89.7	10-175	18.2	99		12/3/2010 1149h
1012057-001BMSD	4-Nitrophenol	$\mug/L$	EPA625	139	80.00	0	174	10-97	31.7	69	1	12/3/2010 1149h
1012057-001BMSD	5-Nitro-o-toluidine	$\mug/L$	EPA625	92.1	80.00	0	115	10-175	15.7	26		12/3/2010 1149h
1012057-001BMSD	7,12-Dimethylbenz(a)anthracene	$\mug/L$	EPA625	106	80.00	0	132	26-174	17.0	40		12/3/2010 1149h
1012057-001BMSD	a,a-Dimethylphenethylamine	$\mu g/L$	EPA625	43.4	80.00	0	54.2	10-175	10.4	99		12/3/2010 1149h
1012057-001BMSD	Acenaphthene	$\mu g/L$	EPA625	68.6	80.00	0	85.7	29-97	20.2	38		12/3/2010 1149h
1012057-001BMSD	Acenaphthylene	$\mug/L$	EPA625	69.5	80.00	0	86.9	37-87	21.5	37		12/3/2010 1149h
1012057-001BMSD	Acetophenone	$\mug/L$	EPA625	61.2	80.00	0	76.5	10-96	27.2	48		12/3/2010 1149h
1012057-001BMSD	alpha-Terpineol	$\mug/L$	EPA625	56.7	80.00	0	70.8	10-67	20.3	46	1	12/3/2010 1149h
1012057-001BMSD	Aniline	$\mug/L$	EPA625	46.7	80.00	0	58.4	10-71	37.5	48		12/3/2010 1149h
1012057-001BMSD	Anthracene	$\mug/L$	EPA625	81.8	80.00	0	102	53-114	16.2	26		12/3/2010 1149h
1012057-001BMSD	Aramite	$\mug/L$	EPA625	89.0	80.00	0	111	29-160	15.6	30		12/3/2010 1149h
1012057-001BMSD	Azobenzene	$\mug/L$	EPA625	65.8	80.00	0	82.2	15-114	16.6	36		12/3/2010 1149h
1012057-001BMSD	Benz(a)anthracene	$\mug/L$	EPA625	73.2	80.00	0	91.5	39-129	15.0	32		12/3/2010 1149h
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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMSD	Benzidine	μg/L	EPA625	66.6	80.00	0	83.2	10-99	15.7	99		12/3/2010 1149h
1012057-001BMSD	Benzo(a)pyrene	$\mu g/L$	EPA625	84.4	80.00	0	105	29-175	17.2	45		12/3/2010 1149h
1012057-001BMSD	Benzo(b)fluoranthene	$\mug/L$	EPA625	73.3	80.00	0	91.7	15-140	18.2	44		12/3/2010 1149h
1012057-001BMSD	Benzo(g,h,i)perylene	$\mug/L$	EPA625	68.3	80.00	0	85.4	10-182	16.6	48		12/3/2010 1149h
1012057-001BMSD	Benzo(k)fluoranthene	$\mug/L$	EPA625	101	80.00	0	127	21-154	15.6	52		12/3/2010 1149h
1012057-001BMSD	Benzoic acid	$\mug/L$	EPA625	< 20.0	80.00	0	18.2	10-71	0	78		12/3/2010 1149h
1012057-001BMSD	Benzyl alcohol	$\mug/L$	EPA625	26.0	80.00	0	32.5	10-69	33.4	52		12/3/2010 1149h
1012057-001BMSD	Bis(2-chloroethoxy)methane	$\mug/L$	EPA625	44.0	80.00	0	54.9	10-94	24.7	45		12/3/2010 1149h
1012057-001BMSD	Bis(2-chloroethyl) ether	$\mug/L$	EPA625	50.5	80.00	0	63.1	10-70	36.5	47		12/3/2010 1149h
1012057-001BMSD	Bis(2-chloroisopropyl) ether	$\mug/L$	EPA625	47.3	80.00	0	59.2	10-71	34.2	49		12/3/2010 1149h
1012057-001BMSD	Bis(2-ethylhexyl) phthalate	$\mug/L$	EPA625	88.0	80.00	0	110	10-175	12.3	28		12/3/2010 1149h
1012057-001BMSD	bis(2-ethylhexyl)adipate	$\mug/L$	EPA625	89.9	80.00	0	112	10-175	10.5	25		12/3/2010 1149h
1012057-001BMSD	Butyl benzyl phthalate	$\mug/L$	EPA625	89.7	80.00	0	112	10-175	10.0	99		12/3/2010 1149h
1012057-001BMSD	Carbazole	$\mug/L$	EPA625	77.4	80.00	0	96.8	10-151	15.2	30		12/3/2010 1149h
1012057-001BMSD	Chlorobenzilate	$\mug/L$	EPA625	87.0	80.00	0	109	18-175	11.2	25		12/3/2010 1149h
1012057-001BMSD	Chrysene	$\mug/L$	EPA625	80.0	80.00	0	100	38-133	13.6	28		12/3/2010 1149h
1012057-001BMSD	Diallate (cis or trans)	$\mug/L$	EPA625	69.2	80.00	0	86.5	10-157	16.8	29		12/3/2010 1149h
1012057-001BMSD	Dibenz(a,h)anthracene	$\mug/L$	EPA625	70.3	80.00	0	87.9	13-168	17.4	51		12/3/2010 1149h
1012057-001BMSD	Dibenzofuran	$\mug/L$	EPA625	69.1	80.00	0	86.4	29-103	22.0	34		12/3/2010 1149h
1012057-001BMSD	Diethyl phthalate	$\mug/L$	EPA625	69.6	80.00	0	87.0	10-139	14.7	32		12/3/2010 1149h
1012057-001BMSD	Dimethoate	$\mug/L$	EPA625	26.7	80.00	0	33.3	10-136	21.2	45		12/3/2010 1149h
1012057-001BMSD	Dimethyl phthalate	$\mug/L$	EPA625	61.8	80.00	0	77.3	10-122	15.8	57		12/3/2010 1149h
1012057-001BMSD	Dimethylaminoazobenzene	$\mug/L$	EPA625	87.7	80.00	0	110	34-142	16.7	26		12/3/2010 1149h
1012057-001BMSD	Di-n-butyl phthalate	$\mug/L$	EPA625	79.0	80.00	0	98.8	44-124	11.1	25		12/3/2010 1149h
										D	D-4 10	/6/2010 Page 121 of 162

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Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMSD	Di-n-octyl phthalate	μg/L	EPA625	137	80.00	0	172	10-175	16.5	50		12/3/2010 1149h
1012057-001BMSD	Dinoseb	$\mug/L$	EPA625	128	80.00	0	160	10-175	11.6	42		12/3/2010 1149h
1012057-001BMSD	Diphenylamine	$\mug\!/L$	EPA625	76.0	80.00	0	95.0	13-110	15.7	34		12/3/2010 1149h
1012057-001BMSD	Disulfoton	$\mug/L$	EPA625	56.3	80.00	0	70.4	10-121	14.9	25		12/3/2010 1149h
1012057-001BMSD	Ethyl methanesulfonate	$\mug/L$	EPA625	55.8	80.00	0	69.8	10-99	30.1	46		12/3/2010 1149h
1012057-001BMSD	Famphur	$\mug/L$	EPA625	126	80.00	0	158	10-71	5.66	25	1	12/3/2010 1149h
1012057-001BMSD	Fluoranthene	$\mug/L$	EPA625	77.3	80.00	0	96.6	23-135	14.4	25		12/3/2010 1149h
1012057-001BMSD	Fluorene	$\mug/L$	EPA625	69.7	80.00	0	87.1	34-108	19.0	28		12/3/2010 1149h
1012057-001BMSD	Hexachlorobenzene	$\mug/L$	EPA625	72.0	80.00	0	90.1	26-131	16.7	28		12/3/2010 1149h
1012057-001BMSD	Hexachlorobutadiene	$\mug/L$	EPA625	41.9	80.00	0	52.4	10-110	22.2	68		12/3/2010 1149h
1012057-001BMSD	Hexachlorocyclopentadiene	$\mug/L$	EPA625	31.7	80.00	0	39.6	10-45	39.8	79		12/3/2010 1149h
1012057-001BMSD	Hexachloroethane	$\mug/L$	EPA625	30.4	80.00	0	38.0	10-58	12.1	42		12/3/2010 1149h
1012057-001BMSD	Hexachlorophene	$\mug/L$	EPA625	84.3	80.00	0	105	10-168	11.6	25		12/3/2010 1149h
1012057-001BMSD	Hexachloropropene	$\mug/L$	EPA625	40.2	80.00	0	50.3	10-72	23.9	63		12/3/2010 1149h
1012057-001BMSD	Indene	$\mug/L$	EPA625	41.9	80.00	0	52.4	10-35	23.8	35	1	12/3/2010 1149h
1012057-001BMSD	Indeno(1,2,3-cd)pyrene	$\mug/L$	EPA625	71.3	80.00	0	89.2	10-176	17.1	48		12/3/2010 1149h
1012057-001BMSD	Isodrin	$\mug/L$	EPA625	76.2	80.00	0	95.2	15-165	15.3	25		12/3/2010 1149h
1012057-001BMSD	Isophorone	$\mug/L$	EPA625	58.2	80.00	0	72.8	10-99	23.6	41		12/3/2010 1149h
1012057-001BMSD	Isosafrole	$\mu g/L$	EPA625	60.5	80.00	0	75.6	10-167	26.4	50		12/3/2010 1149h
1012057-001BMSD	Kepone	$\mu g/L$	EPA625	212	80.00	0	265	10-175	0.383	46	1	12/3/2010 1149h
1012057-001BMSD	Methapyrilene	$\mug/L$	EPA625	84.3	80.00	0	105	10-149	13.1	51		12/3/2010 1149h
1012057-001BMSD	Methyl methanesulfonate	$\mug/L$	EPA625	35.2	80.00	0	44.0	10-132	31.1	45		12/3/2010 1149h
1012057-001BMSD	Naphthalene	$\mu g/L$	EPA625	53.0	80.00	0	66.3	10-82	24.8	43		12/3/2010 1149h
1012057-001BMSD	n-Decane	$\mu g/L$	EPA625	17.0	80.00	0	21.2	10-27	2.67	32		12/3/2010 1149h
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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMSD	Nitrobenzene	μg/L	EPA625	72.7	80.00	0	90.9	10-119	29.7	49		12/3/2010 1149h
1012057-001BMSD	Nitroquinoline-1-oxide	$\mu g/L$	EPA625	36.9	80.00	0	46.1	10-170	22.9	97		12/3/2010 1149h
1012057-001BMSD	N-Nitrosodiethylamine	$\mug/L$	EPA625	56.2	80.00	0	70.2	10-91	32.4	54		12/3/2010 1149h
1012057-001BMSD	N-Nitrosodimethylamine	$\mug/L$	EPA625	21.8	80.00	0	27.2	10-42	34.8	66		12/3/2010 1149h
1012057-001BMSD	N-Nitrosodi-n-butylamine	$\mug/L$	EPA625	63.8	80.00	0	79.8	10-175	19.4	80		12/3/2010 1149h
1012057-001BMSD	N-Nitrosodiphenylamine	$\mug/L$	EPA625	75.1	80.00	0	93.8	12-112	16.1	30		12/3/2010 1149h
1012057-001BMSD	N-Nitrosodi-n-propylamine	$\mug/L$	EPA625	54.4	80.00	0	68.0	10-77	27.4	47		12/3/2010 1149h
1012057-001BMSD	N-Nitrosomethylethylamine	$\mug/L$	EPA625	44.0	80.00	0	55.0	10-75	35.5	60		12/3/2010 1149h
1012057-001BMSD	N-Nitrosomorpholine	$\mug/L$	EPA625	56.8	80.00	0	71.0	10-175	25.7	73		12/3/2010 1149h
1012057-001BMSD	N-Nitrosopiperidine	$\mug/L$	EPA625	64.5	80.00	0	80.6	10-105	21.7	44		12/3/2010 1149h
1012057-001BMSD	N-Nitrosopyrrolidine	$\mug/L$	EPA625	61.2	80.00	0	76.5	10-88	21.0	40		12/3/2010 1149h
1012057-001BMSD	n-Octadecane	$\mug/L$	EPA625	65.2	80.00	0	81.5	10-121	19.4	40		12/3/2010 1149h
1012057-001BMSD	O,O,O-Triethyl phosphorothioate	$\mug/L$	EPA625	64.1	80.00	0	80.1	10-93	22.3	44		12/3/2010 1149h
1012057-001BMSD	o-Toluidine	$\mug/L$	EPA625	125	80.00	0	157	10-107	30.7	46	1	12/3/2010 1149h
1012057-001BMSD	Parathion	$\mug/L$	EPA625	110	80.00	0	138	10-175	15.5	28		12/3/2010 1149h
1012057-001BMSD	Methyl parathion	$\mug/L$	EPA625	101	80.00	0	126	10-175	14.0	26		12/3/2010 1149h
1012057-001BMSD	Pentachlorobenzene	$\mug/L$	EPA625	68.0	80.00	0	85.1	25-134	21.1	35		12/3/2010 1149h
1012057-001BMSD	Pentachloronitrobenzene	$\mug/L$	EPA625	82.9	80.00	0	104	10-175	14.8	25		12/3/2010 1149h
1012057-001BMSD	Pentachlorophenol	$\mug/L$	EPA625	45.7	80.00	0	57.2	10-163	1.45	58		12/3/2010 1149h
1012057-001BMSD	Phenacetin	$\mug/L$	EPA625	75.2	80.00	0	94.0	10-175	17.1	38		12/3/2010 1149h
1012057-001BMSD	Phenanthrene	$\mug/L$	EPA625	84.6	80.00	0	106	31-126	19.4	32		12/3/2010 1149h
1012057-001BMSD	Phenol	$\mug/L$	EPA625	27.3	80.00	0	34.2	10-175	25.0	71		12/3/2010 1149h
1012057-001BMSD	Phorate	$\mug/L$	EPA625	82.6	80.00	0	103	10-175	14.7	42		12/3/2010 1149h
1012057-001BMSD	Pronamide	$\mug/L$	EPA625	81.6	80.00	0	102	10-95	13.5	26	1	12/3/2010 1149h

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Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012057-001BMSD	Pyrene	μg/L	EPA625	89.0	80.00	0	111	51-139	13.7	27		12/3/2010 1149h
1012057-001BMSD	Pyridine	μg/L	EPA625	23.1	80.00	0	28.9	10-25	52.6	61	1	12/3/2010 1149h
1012057-001BMSD	Quinoline	μg/L	EPA625	67.0	80.00	0	83.7	10-63	20.5	99	1	12/3/2010 1149h
1012057-001BMSD	Safrole	μg/L	EPA625	61.5	80.00	0	76.9	10-120	27.6	51		12/3/2010 1149h
1012057-001BMSD	Tetraethyl dithiopyrophosphate	μg/L	EPA625	74.2	80.00	0	92.8	13-160	17.4	35		12/3/2010 1149h
1012057-001BMSD	Thionazin	$\mu$ g/L	EPA625	68.5	80.00	0	85.6	10-139	14.5	25		12/3/2010 1149h
1012057-001BMSD	Surr: 2,4,6-Tribromophenol	%REC	EPA625	67.9	80.00		84.8	21-154				12/3/2010 1149h
1012057-001BMSD	Surr: 2-Fluorobiphenyl	%REC	EPA625	64.6	40.00		161	10-106			S	12/3/2010 1149h
1012057-001BMSD	Surr: 2-Fluorophenol	%REC	EPA625	25.0	80.00		31.3	10-56				12/3/2010 1149h
1012057-001BMSD	Surr: Nitrobenzene-d5	%REC	EPA625	64.0	40.00		160	10-101			S	12/3/2010 1149h
1012057-001BMSD	Surr: Phenol-d6	%REC	EPA625	15.0	80.00		18.8	10-45				12/3/2010 1149h
1012057-001BMSD	Surr: Terphenyl-d14	%REC	EPA625	89.1	40.00		223	10-160			S	12/3/2010 1149h
1011529-006FMSD	Acenaphthene	$\mug/L$	SW8270D	59.1	80.00	0	73.9	21-113	7.73	25		12/3/2010 0144h
1011529-006FMSD	Benzo(a)pyrene	μg/L	SW8270D	85.0	80.00	0	106	15-169	6.66	25		12/3/2010 0144h
1011529-006FMSD	Pyrene	μg/L	SW8270D	91.5	80.00	0	114	23-150	4.01	25		12/3/2010 0144h
1011529-006FMSD	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	68.2	80.00		85.2	14-159				12/3/2010 0144h
1011529-006FMSD	Surr: 2-Fluorobiphenyl	%REC	SW8270D	49.3	40.00		123	10-124				12/3/2010 0144h
1011529-006FMSD	Surr: 2-Fluorophenol	%REC	SW8270D	26.8	80.00		33.6	10-106				12/3/2010 0144h
1011529-006FMSD	Surr: Nitrobenzene-d5	%REC	SW8270D	53.9	40.00		135	10-180				12/3/2010 0144h
1011529-006FMSD	Surr: Phenol-d6	%REC	SW8270D	19.7	80.00		24.6	10-122				12/3/2010 0144h
1011529-006FMSD	Surr: Terphenyl-d14	%REC	SW8270D	92.3	40.00		231	10-199			S	12/3/2010 0144h
1012039-001DMSD	Acenaphthene	$\mu$ g/L	SW8270D	67.6	83.33	0	81.2	21-113	0.972	25		12/3/2010 0959h
1012039-001DMSD	Benzo(a)pyrene	μg/L	SW8270D	76.3	83.33	0	91.6	15-169	3.31	25		12/3/2010 0959h
1012039-001DMSD	Pyrene	μg/L	SW8270D	86.5	83.33	0	104	23-150	2.39	25		12/3/2010 0959h
										Danort	Data: 12	/6/2010 Page 13/Lof 16'

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Jose Rocha

QA Officer

### **OC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSSV

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012039-001DMSD	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	70.6	83.33		84.7	14-159				12/3/2010 0959h
1012039-001DMSD	Surr: 2-Fluorobiphenyl	%REC	SW8270D	63.1	41.67		152	10-124			S	12/3/2010 0959h
1012039-001DMSD	Surr: 2-Fluorophenol	%REC	SW8270D	30.6	83.33		36.8	10-106				12/3/2010 0959h
1012039-001DMSD	Surr: Nitrobenzene-d5	%REC	SW8270D	60.8	41.67		146	10-180				12/3/2010 0959h
1012039-001DMSD	Surr: Phenol-d6	%REC	SW8270D	20.9	83.33		25.1	10-122				12/3/2010 0959h
1012039-001DMSD	Surr: Terphenyl-d14	%REC	SW8270D	88.3	41.67		212	10-199			S	12/3/2010 0959h
1012057-001BMSD	Acenaphthene	$\mug/L$	SW8270D	68.6	80.00	0	85.7	21-113	20.2	25		12/3/2010 1149h
1012057-001BMSD	Benzo(a)pyrene	$\mug/L$	SW8270D	84.4	80.00	0	105	15-169	17.2	25		12/3/2010 1149h
1012057-001BMSD	Pyrene	$\mug/L$	SW8270D	89.0	80.00	0	111	23-150	13.7	25		12/3/2010 1149h
1012057-001BMSD	Surr: 2,4,6-Tribromophenol	%REC	SW8270D	67.9	80.00		84.8	14-159				12/3/2010 1149h
1012057-001BMSD	Surr: 2-Fluorobiphenyl	%REC	SW8270D	64.6	40.00		161	10-124			S	12/3/2010 1149h
1012057-001BMSD	Surr: 2-Fluorophenol	%REC	SW8270D	25.0	80.00		31.3	10-106				12/3/2010 1149h
1012057-001BMSD	Surr: Nitrobenzene-d5	%REC	SW8270D	64.0	40.00		160	10-180				12/3/2010 1149h
1012057-001BMSD	Surr: Phenol-d6	%REC	SW8270D	15.0	80.00		18.8	10-122				12/3/2010 1149h
1012057-001BMSD	Surr: Terphenyl-d14	%REC	SW8270D	89.1	40.00		223	10-199			S	12/3/2010 1149h

<sup>@ -</sup> High RPD due to suspected sample non-homogeneity or matrix interference.

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<sup>&</sup>lt;sup>1</sup> - Matrix spike recovery indicates matrix interference. The method is in control as indicated by the LCS.

S - Surrogate recoveries outside the control limits. Repreparation and reanalysis and/or MS samples yielded similar results indicating matrix interference.



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Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
LCS VOC 120210A	1,1,1,2-Tetrachloroethane	μg/L	EPA624	20.0	20.00	0	99.8	74-117				12/2/2010 0711h
LCS VOC 120210A	1,1,1-Trichloroethane	μg/L	EPA624	20.0	20.00	0	100	49-140				12/2/2010 0711h
LCS VOC 120210A	1,1,2,2-Tetrachloroethane	μg/L	EPA624	15.8	20.00	0	79.0	67-119				12/2/2010 0711h
LCS VOC 120210A	1,1,2-Trichloro-1,2,2-trifluoroetha	μg/L	EPA624	19.2	20.00	0	96.2	21-206				12/2/2010 0711h
LCS VOC 120210A	1,1,2-Trichloroethane	μg/L	EPA624	18.5	20.00	0	92.6	80-123				12/2/2010 0711h
LCS VOC 120210A	1,1-Dichloropropene	μg/L	EPA624	41.8	40.00	0	105	10-140				12/2/2010 0711h
LCS VOC 120210A	1,1-Dichloroethane	μg/L	EPA624	21.2	20.00	0	106	70-130				12/2/2010 0711h
LCS VOC 120210A	1,1-Dichloroethene	μg/L	EPA624	23.7	20.00	0	119	52-171				12/2/2010 0711h
LCS VOC 120210A	1,2,3-Trichlorobenzene	μg/L	EPA624	19.9	20.00	0	99.6	67-131				12/2/2010 0711h
LCS VOC 120210A	1,2,3-Trichloropropane	μg/L	EPA624	16.3	20.00	0	81.6	62-116				12/2/2010 0711h
LCS VOC 120210A	1,2,3-Trimethylbenzene	μg/L	EPA624	21.2	20.00	0	106	76-140				12/2/2010 0711h
LCS VOC 120210A	1,2,4-Trichlorobenzene	μg/L	EPA624	19.9	20.00	0	99.4	58-133				12/2/2010 0711h
LCS VOC 120210A	1,2,4-Trimethylbenzene	μg/L	EPA624	20.3	20.00	0	102	79-151				12/2/2010 0711h
LCS VOC 120210A	1,2-Dibromo-3-chloropropane	μg/L	EPA624	15.8	20.00	0	79.1	64-129				12/2/2010 0711h
LCS VOC 120210A	1,2-Dibromoethane	μg/L	EPA624	18.6	20.00	0	92.9	70-126				12/2/2010 0711h
LCS VOC 120210A	1,2-Dichlorobenzene	μg/L	EPA624	20.8	20.00	0	104	67-135				12/2/2010 0711h
LCS VOC 120210A	1,2-Dichloroethane	μg/L	EPA624	19.1	20.00	0	95.3	60-137				12/2/2010 0711h
LCS VOC 120210A	1,2-Dichloropropane	μg/L	EPA624	20.5	20.00	0	103	59-135				12/2/2010 0711h
LCS VOC 120210A	1,3,5-Trimethylbenzene	μg/L	EPA624	20.3	20.00	0	101	77-151				12/2/2010 0711h
LCS VOC 120210A	1,3-Dichlorobenzene	μg/L	EPA624	21.0	20.00	0	105	78-134				12/2/2010 0711h
LCS VOC 120210A	1,3-Dichloropropane	μg/L	EPA624	18.4	20.00	0	91.9	78-116				12/2/2010 0711h
LCS VOC 120210A	1,4-Dichlorobenzene	μg/L	EPA624	20.8	20.00	0	104	72-139				12/2/2010 0711h
LCS VOC 120210A	1,4-Dioxane	μg/L	EPA624	184	200.0	0	92.1	33-149				12/2/2010 0711h
LCS VOC 120210A	2,2-Dichloropropane	μg/L	EPA624	20.1	30.00	0	67.1	13-180				12/2/2010 0711h
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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
LCS VOC 120210A	2-Butanone	μg/L	EPA624	24.1	20.00	0	120	10-217				12/2/2010 0711h
LCS VOC 120210A	2-Chloroethyl vinyl ether	μg/L	EPA624	41.0	40.00	0	102	32-163				12/2/2010 0711h
LCS VOC 120210A	2-Chlorotoluene	$\mu g/L$	EPA624	20.2	20.00	0	101	79-142				12/2/2010 0711h
LCS VOC 120210A	2-Hexanone	μg/L	EPA624	15.9	20.00	0	79.7	50-156				12/2/2010 0711h
LCS VOC 120210A	2-Nitropropane	$\mu g/L$	EPA624	16.6	20.00	0	82.8	10-243				12/2/2010 0711h
LCS VOC 120210A	4-Chlorotoluene	$\mu g/L$	EPA624	21.2	20.00	0	106	68-128				12/2/2010 0711h
LCS VOC 120210A	4-Isopropyltoluene	$\mug/L$	EPA624	18.8	20.00	0	93.9	73-156				12/2/2010 0711h
LCS VOC 120210A	4-Methyl-2-pentanone	$\mu g/L$	EPA624	15.6	20.00	0	77.8	10-214				12/2/2010 0711h
LCS VOC 120210A	Acetone	$\mu g/L$	EPA624	18.5	20.00	0	92.4	10-313				12/2/2010 0711h
LCS VOC 120210A	Acetonitrile	$\mu g/L$	EPA624	33.1	40.00	0	82.8	37-159				12/2/2010 0711h
LCS VOC 120210A	Acrolein	$\mu g/L$	EPA624	44.9	40.00	0	112	10-325				12/2/2010 0711h
LCS VOC 120210A	Acrylonitrile	$\mu g/L$	EPA624	15.3	20.00	0	76.7	53-134				12/2/2010 0711h
LCS VOC 120210A	Allyl chloride	$\mu g/L$	EPA624	20.3	20.00	0	102	10-243				12/2/2010 0711h
LCS VOC 120210A	Benzene	$\mu g/L$	EPA624	24.0	20.00	0	120	62-127				12/2/2010 0711h
LCS VOC 120210A	Benzyl chloride	$\mu g/L$	EPA624	17.5	20.00	0	87.7	40-146				12/2/2010 0711h
LCS VOC 120210A	Bis(2-chloroisopropyl) ether	$\mu g/L$	EPA624	15.8	20.00	0	79.0	54-146				12/2/2010 0711h
LCS VOC 120210A	Bromobenzene	$\mu g/L$	EPA624	19.2	20.00	0	95.9	78-148				12/2/2010 0711h
LCS VOC 120210A	Bromochloromethane	$\mu g/L$	EPA624	21.2	20.00	0	106	75-134				12/2/2010 0711h
LCS VOC 120210A	Bromodichloromethane	$\mu g/L$	EPA624	19.8	20.00	0	99.0	74-121				12/2/2010 0711h
LCS VOC 120210A	Bromoform	$\mu g/L$	EPA624	17.7	20.00	0	88.5	68-131				12/2/2010 0711h
LCS VOC 120210A	Bromomethane	$\mu g/L$	EPA624	14.2	20.00	0	71.0	10-185				12/2/2010 0711h
LCS VOC 120210A	Butyl acetate	$\mu g/L$	EPA624	18.5	20.00	0	92.3	46-178				12/2/2010 0711h
LCS VOC 120210A	Carbon disulfide	$\mu g/L$	EPA624	32.5	20.00	0	163	21-224				12/2/2010 0711h
LCS VOC 120210A	Carbon tetrachloride	$\mu g/L$	EPA624	20.2	20.00	0	101	60-157				12/2/2010 0711h
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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: LCS

a		** *.	36.0	<b>.</b>	Amount	Original	a/ PEG	<b>.</b>	0 / DDD	RPD		D
Sample ID	Analyte	Units	Method	Result	Spiked	Amount	% REC	Limits	%RPD	Limit	Qual	Date Analyzed
LCS VOC 120210A	Chlorobenzene	$\mug/L$	EPA624	20.8	20.00	0	104	63-140				12/2/2010 0711h
LCS VOC 120210A	Chloroethane	$\mug/L$	EPA624	20.6	20.00	0	103	41-173				12/2/2010 0711h
LCS VOC 120210A	Chloroform	$\mug/L$	EPA624	20.4	20.00	0	102	67-132				12/2/2010 0711h
LCS VOC 120210A	Chloromethane	$\mug/L$	EPA624	16.0	20.00	0	79.8	10-138				12/2/2010 0711h
LCS VOC 120210A	Chloroprene	$\mug/L$	EPA624	19.7	20.00	0	98.5	10-161				12/2/2010 0711h
LCS VOC 120210A	cis-1,2-Dichloroethene	$\mug/L$	EPA624	21.8	20.00	0	109	72-137				12/2/2010 0711h
LCS VOC 120210A	cis-1,3-Dichloropropene	$\mug/L$	EPA624	21.5	40.00	0	53.8	10-134				12/2/2010 0711h
LCS VOC 120210A	Cyclohexane	$\mu g/L$	EPA624	22.6	20.00	0	113	35-230				12/2/2010 0711h
LCS VOC 120210A	Cyclohexanone	$\mu g/L$	EPA624	< 50.0	40.00	0	60.4	10-374				12/2/2010 0711h
LCS VOC 120210A	Dibromochloromethane	$\mu g/L$	EPA624	19.3	20.00	0	96.5	68-135				12/2/2010 0711h
LCS VOC 120210A	Dibromomethane	$\mu g/L$	EPA624	18.8	20.00	0	93.8	74-120				12/2/2010 0711h
LCS VOC 120210A	Dichlorodifluoromethane	$\mu g/L$	EPA624	11.9	20.00	0	59.4	10-150				12/2/2010 0711h
LCS VOC 120210A	Ethyl acetate	$\mu g/L$	EPA624	35.3	40.00	0	88.3	50-155				12/2/2010 0711h
LCS VOC 120210A	Ethyl ether	$\mu g/L$	EPA624	21.5	20.00	0	108	45-146				12/2/2010 0711h
LCS VOC 120210A	Ethyl methacrylate	$\mu g/L$	EPA624	17.7	20.00	0	88.4	64-113				12/2/2010 0711h
LCS VOC 120210A	Ethylbenzene	$\mu g/L$	EPA624	20.7	20.00	0	104	55-133				12/2/2010 0711h
LCS VOC 120210A	Hexachlorobutadiene	$\mu g/L$	EPA624	16.8	20.00	0	84.0	35-213				12/2/2010 0711h
LCS VOC 120210A	Iodomethane	μg/L	EPA624	16.4	20.00	0	81.8	10-233				12/2/2010 0711h
LCS VOC 120210A	Isobutyl alcohol	μg/L	EPA624	< 100	80.00	0	89.4	12-202				12/2/2010 0711h
LCS VOC 120210A	Isopropyl acetate	μg/L	EPA624	17.8	20.00	0	89.2	55-145				12/2/2010 0711h
LCS VOC 120210A	Isopropyl alcohol	μg/L	EPA624	73.6	80.00	0	91.9	12-250				12/2/2010 0711h
LCS VOC 120210A	Isopropylbenzene	$\mu g/L$	EPA624	19.6	20.00	0	98.1	60-147				12/2/2010 0711h
LCS VOC 120210A	Isopropyltoluene	$\mu g/L$	EPA624	18.8	20.00	0	93.9	73-156				12/2/2010 0711h
LCS VOC 120210A	m,p-Xylene	$\mug/L$	EPA624	43.6	40.00	0	109	70-130				12/2/2010 0711h

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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
LCS VOC 120210A	Methacrylonitrile	μg/L	EPA624	16.2	20.00	0	81.0	79-123				12/2/2010 0711h
LCS VOC 120210A	Methyl Acetate	$\mu g/L$	EPA624	34.8	20.00	0	174	5-398				12/2/2010 0711h
LCS VOC 120210A	Methyl methacrylate	$\mu g/L$	EPA624	16.8	20.00	0	84.1	55-128				12/2/2010 0711h
LCS VOC 120210A	Methyl tert-butyl ether	$\mug/L$	EPA624	19.4	20.00	0	96.8	37-189				12/2/2010 0711h
LCS VOC 120210A	Methylcyclohexane	$\mug/L$	EPA624	20.7	20.00	0	103	65-175				12/2/2010 0711h
LCS VOC 120210A	Methylene chloride	$\mu g/L$	EPA624	23.3	20.00	0	117	55-138				12/2/2010 0711h
LCS VOC 120210A	n-Amyl acetate	$\mu g/L$	EPA624	12.9	20.00	0	64.6	10-187				12/2/2010 0711h
LCS VOC 120210A	Naphthalene	$\mug/L$	EPA624	19.1	20.00	0	95.7	41-131				12/2/2010 0711h
LCS VOC 120210A	n-Butyl alcohol	$\mu g/L$	EPA624	68.6	80.00	0	85.8	10-226				12/2/2010 0711h
LCS VOC 120210A	n-Butylbenzene	$\mu g/L$	EPA624	18.6	20.00	0	93.1	40-158				12/2/2010 0711h
LCS VOC 120210A	n-Hexane	$\mug/L$	EPA624	22.9	20.00	0	114	10-277				12/2/2010 0711h
LCS VOC 120210A	n-Octane	$\mug/L$	EPA624	16.0	20.00	0	80.0	45-158				12/2/2010 0711h
LCS VOC 120210A	n-Propylbenzene	$\mu g/L$	EPA624	19.8	20.00	0	99.0	67-131				12/2/2010 0711h
LCS VOC 120210A	o-Xylene	$\mu g/L$	EPA624	20.8	20.00	0	104	70-130				12/2/2010 0711h
LCS VOC 120210A	Pentachloroethane	$\mu g/L$	EPA624	15.5	20.00	0	77.4	10-314				12/2/2010 0711h
LCS VOC 120210A	Propionitrile	$\mu g/L$	EPA624	< 25.0	20.00	0	77.5	60-132				12/2/2010 0711h
LCS VOC 120210A	Propyl acetate	$\mu g/L$	EPA624	18.2	20.00	0	90.8	48-143				12/2/2010 0711h
LCS VOC 120210A	sec-Butylbenzene	$\mug/L$	EPA624	19.1	20.00	0	95.4	72-157				12/2/2010 0711h
LCS VOC 120210A	Styrene	$\mu g/L$	EPA624	20.9	20.00	0	104	81-125				12/2/2010 0711h
LCS VOC 120210A	tert-Butyl alcohol	$\mu g/L$	EPA624	32.8	40.00	0	82.0	50-286				12/2/2010 0711h
LCS VOC 120210A	tert-Butylbenzene	$\mug/L$	EPA624	18.6	20.00	0	92.8	75-157				12/2/2010 0711h
LCS VOC 120210A	Tetrachloroethene	$\mu g/L$	EPA624	24.7	20.00	0	124	49-163				12/2/2010 0711h
LCS VOC 120210A	Tetrahydrofuran	$\mug/L$	EPA624	14.8	20.00	0	74.0	43-146				12/2/2010 0711h
LCS VOC 120210A	Toluene	$\mug/L$	EPA624	21.3	20.00	0	107	67-128				12/2/2010 0711h
										Domont	Data: 12	6/2010 Page 120 of 162

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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
LCS VOC 120210A	trans-1,2-Dichloroethene	μg/L	EPA624	24.7	20.00	0	123	47-146				12/2/2010 0711h
LCS VOC 120210A	trans-1,3-Dichloropropene	$\mu g/L$	EPA624	20.4	20.00	0	102	29-143				12/2/2010 0711h
LCS VOC 120210A	trans-1,4-Dichloro-2-butene	$\mug/L$	EPA624	21.2	20.00	0	106	20-214				12/2/2010 0711h
LCS VOC 120210A	Trichloroethene	$\mu g/L$	EPA624	21.5	20.00	0	108	54-152				12/2/2010 0711h
LCS VOC 120210A	Trichlorofluoromethane	$\mug/L$	EPA624	16.2	20.00	0	80.9	56-166				12/2/2010 0711h
LCS VOC 120210A	Vinyl acetate	$\mug/L$	EPA624	27.5	40.00	0	68.7	38-121				12/2/2010 0711h
LCS VOC 120210A	Vinyl chloride	$\mug/L$	EPA624	17.7	20.00	0	88.4	13-155				12/2/2010 0711h
LCS VOC 120210A	Xylenes, Total	$\mu g/L$	EPA624	64.4	60.00	0	107	52-130				12/2/2010 0711h
LCS VOC 120210A	Surr: 1,2-Dichloroethane-d4	%REC	EPA624	47.0	50.00		94.1	69-132				12/2/2010 0711h
LCS VOC 120210A	Surr: 4-Bromofluorobenzene	%REC	EPA624	49.5	50.00		99.0	85-118				12/2/2010 0711h
LCS VOC 120210A	Surr: Dibromofluoromethane	%REC	EPA624	49.3	50.00		98.7	80-120				12/2/2010 0711h
LCS VOC 120210A	Surr: Toluene-d8	%REC	EPA624	49.5	50.00		99.0	81-120				12/2/2010 0711h

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QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA

QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB VOC 120210A	1,1,1,2-Tetrachloroethane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,1,1-Trichloroethane	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,1,2,2-Tetrachloroethane	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,1,2-Trichloro-1,2,2-trifluoroetha	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,1,2-Trichloroethane	$\mu\text{g}/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,1-Dichloropropene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,1-Dichloroethane	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,1-Dichloroethene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2,3-Trichlorobenzene	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2,3-Trichloropropane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2,3-Trimethylbenzene	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2,4-Trichlorobenzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2,4-Trimethylbenzene	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2-Dibromo-3-chloropropane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2-Dibromoethane	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2-Dichlorobenzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2-Dichloroethane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,2-Dichloropropane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,3,5-Trimethylbenzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,3-Dichlorobenzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,3-Dichloropropane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,4-Dichlorobenzene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	1,4-Dioxane	$\mu g/L$	EPA624	< 40.0				-				12/2/2010 0750h
MB VOC 120210A	2,2-Dichloropropane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
										Report	Date: 12/	6/2010 Page 141 of 162

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Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA

QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB VOC 120210A	2-Butanone	μg/L	EPA624	< 10.0				-				12/2/2010 0750h
MB VOC 120210A	2-Chloroethyl vinyl ether	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	2-Chlorotoluene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	2-Hexanone	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	2-Nitropropane	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	4-Chlorotoluene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	4-Isopropyltoluene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	4-Methyl-2-pentanone	$\mu g/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Acetone	μg/L	EPA624	< 10.0				-				12/2/2010 0750h
MB VOC 120210A	Acetonitrile	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Acrolein	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Acrylonitrile	μg/L	EPA624	< 10.0				-				12/2/2010 0750h
MB VOC 120210A	Allyl chloride	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Benzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Benzyl chloride	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Bis(2-chloroisopropyl) ether	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Bromobenzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Bromochloromethane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Bromodichloromethane	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Bromoform	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Bromomethane	$\mug/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Butyl acetate	μg/L	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Carbon disulfide	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Carbon tetrachloride	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
										Report	Date: 12/	6/2010 Page 142 of 1

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# **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB VOC 120210A	Chlorobenzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Chloroethane	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Chloroform	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Chloromethane	$\mu g/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Chloroprene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	cis-1,2-Dichloroethene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	cis-1,3-Dichloropropene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Cyclohexane	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Cyclohexanone	$\mu g/L$	EPA624	< 50.0				-				12/2/2010 0750h
MB VOC 120210A	Dibromochloromethane	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Dibromomethane	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Dichlorodifluoromethane	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Ethyl acetate	$\mu g/L$	EPA624	< 10.0				-				12/2/2010 0750h
MB VOC 120210A	Ethyl ether	$\mu g/L$	EPA624	< 10.0				-				12/2/2010 0750h
MB VOC 120210A	Ethyl methacrylate	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Ethylbenzene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Hexachlorobutadiene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Iodomethane	$\mu g/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Isobutyl alcohol	$\mu g/L$	EPA624	< 100				-				12/2/2010 0750h
MB VOC 120210A	Isopropyl acetate	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Isopropyl alcohol	$\mug/L$	EPA624	< 25.0				-				12/2/2010 0750h
MB VOC 120210A	Isopropylbenzene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Isopropyltoluene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	m,p-Xylene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
										Report	Date: 12/	6/2010 Page 1/13 of 162

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Kyle F. Gross

Laboratory Director

Jose Rocha

QA Officer

# **QC SUMMARY REPORT**

**Client:** Utah DEQ DERR

**Lab Set ID:** 1011529 Red Butte **Project:** 

Jim Harris **Contact: MSVOA** Dept:

QC Type: MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB VOC 120210A	Methacrylonitrile	$\mug/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Methyl Acetate	$\mug/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Methyl methacrylate	$\mug/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Methyl tert-butyl ether	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Methylcyclohexane	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Methylene chloride	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	n-Amyl acetate	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Naphthalene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	n-Butyl alcohol	$\mug/L$	EPA624	< 25.0				-				12/2/2010 0750h
MB VOC 120210A	n-Butylbenzene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	n-Hexane	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	n-Octane	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	n-Propylbenzene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	o-Xylene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Pentachloroethane	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Propionitrile	$\mug/L$	EPA624	< 25.0				-				12/2/2010 0750h
MB VOC 120210A	Propyl acetate	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	sec-Butylbenzene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Styrene	μg/L	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	tert-Butyl alcohol	$\mug/L$	EPA624	< 20.0				-				12/2/2010 0750h
MB VOC 120210A	tert-Butylbenzene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Tetrachloroethene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Tetrahydrofuran	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Toluene	$\mug/L$	EPA624	< 2.00				-				12/2/2010 0750h
										Report	Date: 12/	6/2010 Page 144 of 162



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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MBLK

g l ID	A 14	<b>T</b> T •/	36.0	<b>D</b> 1/	Amount	Original	A/ DEG	Ŧ,	a/ DDD	RPD	0 1	B
Sample ID	Analyte	Units	Method	Result	Spiked	Amount	% REC	Limits	%RPD	Limit	Qual	Date Analyzed
MB VOC 120210A	TPH C11-C15 (DRO)	$\mug/L$	EPA624	< 20.0				-				12/2/2010 0750h
MB VOC 120210A	TPH C6-C10 (GRO)	$\mug\!/L$	EPA624	< 20.0				-				12/2/2010 0750h
MB VOC 120210A	trans-1,2-Dichloroethene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	trans-1,3-Dichloropropene	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	trans-1,4-Dichloro-2-butene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Trichloroethene	$\mu g/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Trichlorofluoromethane	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Vinyl acetate	$\mug\!/L$	EPA624	< 5.00				-				12/2/2010 0750h
MB VOC 120210A	Vinyl chloride	$\mug\!/L$	EPA624	< 1.00				-				12/2/2010 0750h
MB VOC 120210A	Xylenes, Total	$\mug\!/L$	EPA624	< 2.00				-				12/2/2010 0750h
MB VOC 120210A	Surr: 1,2-Dichloroethane-d4	%REC	EPA624	49.2	50.00		98.5	69-132				12/2/2010 0750h
MB VOC 120210A	Surr: 4-Bromofluorobenzene	%REC	EPA624	52.2	50.00		104	85-118				12/2/2010 0750h
MB VOC 120210A	Surr: Dibromofluoromethane	%REC	EPA624	49.0	50.00		97.9	80-120				12/2/2010 0750h
MB VOC 120210A	Surr: Toluene-d8	%REC	EPA624	50.4	50.00		101	81-120				12/2/2010 0750h

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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD Limit	Qual	Date Analyzed
1011529-006BMS	1,1,1,2-Tetrachloroethane	μg/L	EPA624	17.2	20.00	0	86.2	74-117			12/2/2010 1021h
1011529-006BMS	1,1,1-Trichloroethane	μg/L	EPA624	21.8	20.00	0	109	67-147			12/2/2010 1021h
1011529-006BMS	1,1,2,2-Tetrachloroethane	μg/L	EPA624	16.0	20.00	0	80.2	67-119			12/2/2010 1021h
1011529-006BMS	1,1,2-Trichloro-1,2,2-trifluoroetha	μg/L	EPA624	24.2	20.00	0	121	21-206			12/2/2010 1021h
1011529-006BMS	1,1,2-Trichloroethane	μg/L	EPA624	17.3	20.00	0	86.3	80-123			12/2/2010 1021h
1011529-006BMS	1,1-Dichloropropene	μg/L	EPA624	46.1	40.00	0	115	10-140			12/2/2010 1021h
1011529-006BMS	1,1-Dichloroethane	μg/L	EPA624	20.3	20.00	0	101	70-130			12/2/2010 1021h
1011529-006BMS	1,1-Dichloroethene	μg/L	EPA624	24.8	20.00	0	124	62-152			12/2/2010 1021h
1011529-006BMS	1,2,3-Trichlorobenzene	μg/L	EPA624	15.8	20.00	0	78.9	67-131			12/2/2010 1021h
1011529-006BMS	1,2,3-Trichloropropane	μg/L	EPA624	17.1	20.00	0	85.4	62-116			12/2/2010 1021h
1011529-006BMS	1,2,3-Trimethylbenzene	μg/L	EPA624	18.7	20.00	0	93.5	76-140			12/2/2010 1021h
1011529-006BMS	1,2,4-Trichlorobenzene	μg/L	EPA624	15.3	20.00	0	76.4	58-133			12/2/2010 1021h
1011529-006BMS	1,2,4-Trimethylbenzene	μg/L	EPA624	17.6	20.00	0	87.8	79-151			12/2/2010 1021h
1011529-006BMS	1,2-Dibromo-3-chloropropane	μg/L	EPA624	15.8	20.00	0	79.2	64-129			12/2/2010 1021h
1011529-006BMS	1,2-Dibromoethane	μg/L	EPA624	17.3	20.00	0	86.4	70-126			12/2/2010 1021h
1011529-006BMS	1,2-Dichlorobenzene	μg/L	EPA624	17.5	20.00	0	87.4	70-130			12/2/2010 1021h
1011529-006BMS	1,2-Dichloroethane	μg/L	EPA624	20.5	20.00	0	102	39-162			12/2/2010 1021h
1011529-006BMS	1,2-Dichloropropane	μg/L	EPA624	18.6	20.00	0	93.0	59-135			12/2/2010 1021h
1011529-006BMS	1,3,5-Trimethylbenzene	μg/L	EPA624	17.9	20.00	0	89.6	77-151			12/2/2010 1021h
1011529-006BMS	1,3-Dichlorobenzene	μg/L	EPA624	17.3	20.00	0	86.7	78-134			12/2/2010 1021h
1011529-006BMS	1,3-Dichloropropane	μg/L	EPA624	17.1	20.00	0	85.7	78-116			12/2/2010 1021h
1011529-006BMS	1,4-Dichlorobenzene	μg/L	EPA624	16.6	20.00	0	83.1	72-139			12/2/2010 1021h
1011529-006BMS	1,4-Dioxane	μg/L	EPA624	179	200.0	0	89.5	33-149			12/2/2010 1021h
1011529-006BMS	2,2-Dichloropropane	μg/L	EPA624	21.2	30.00	0	70.5	13-180			12/2/2010 1021h
									Report D	ate: 12/	6/2010 Page 146 of 1

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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA

**QC Type:** MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMS	2-Butanone	μg/L	EPA624	26.4	20.00	0	132	10-217				12/2/2010 1021h
1011529-006BMS	2-Chloroethyl vinyl ether	μg/L	EPA624	< 5.00	40.00	0	0	32-163			1	12/2/2010 1021h
1011529-006BMS	2-Chlorotoluene	μg/L	EPA624	17.2	20.00	0	86.0	79-142				12/2/2010 1021h
1011529-006BMS	2-Hexanone	μg/L	EPA624	16.0	20.00	0	79.8	50-156				12/2/2010 1021h
1011529-006BMS	2-Nitropropane	μg/L	EPA624	18.3	20.00	0	91.4	10-243				12/2/2010 1021h
1011529-006BMS	4-Chlorotoluene	μg/L	EPA624	17.5	20.00	0	87.4	68-128				12/2/2010 1021h
1011529-006BMS	4-Isopropyltoluene	μg/L	EPA624	17.2	20.00	0	85.9	73-156				12/2/2010 1021h
1011529-006BMS	4-Methyl-2-pentanone	μg/L	EPA624	17.2	20.00	0	86.1	10-214				12/2/2010 1021h
1011529-006BMS	Acetone	μg/L	EPA624	12.8	20.00	0	64.1	10-313				12/2/2010 1021h
1011529-006BMS	Acetonitrile	μg/L	EPA624	38.1	40.00	0	95.2	37-159				12/2/2010 1021h
1011529-006BMS	Acrolein	μg/L	EPA624	63.8	40.00	0	160	10-325				12/2/2010 1021h
1011529-006BMS	Acrylonitrile	$\mu g/L$	EPA624	17.4	20.00	0	86.9	53-134				12/2/2010 1021h
1011529-006BMS	Allyl chloride	$\mu g/L$	EPA624	19.4	20.00	0	97.2	10-243				12/2/2010 1021h
1011529-006BMS	Benzene	$\mu g/L$	EPA624	21.2	20.00	0	106	66-145				12/2/2010 1021h
1011529-006BMS	Benzyl chloride	$\mu g/L$	EPA624	16.8	20.00	0	84.0	40-146				12/2/2010 1021h
1011529-006BMS	Bis(2-chloroisopropyl) ether	$\mu g/L$	EPA624	16.9	20.00	0	84.5	54-146				12/2/2010 1021h
1011529-006BMS	Bromobenzene	$\mu g/L$	EPA624	16.7	20.00	0	83.6	78-148				12/2/2010 1021h
1011529-006BMS	Bromochloromethane	$\mu g/L$	EPA624	19.5	20.00	0	97.3	75-134				12/2/2010 1021h
1011529-006BMS	Bromodichloromethane	$\mu g/L$	EPA624	19.3	20.00	0	96.6	74-121				12/2/2010 1021h
1011529-006BMS	Bromoform	$\mu g/L$	EPA624	16.6	20.00	0	82.8	68-131				12/2/2010 1021h
1011529-006BMS	Bromomethane	$\mug/L$	EPA624	13.5	20.00	0	67.4	10-185				12/2/2010 1021h
1011529-006BMS	Butyl acetate	$\mug/L$	EPA624	17.6	20.00	0	87.8	46-178				12/2/2010 1021h
1011529-006BMS	Carbon disulfide	$\mug/L$	EPA624	32.9	20.00	0	164	21-224				12/2/2010 1021h
1011529-006BMS	Carbon tetrachloride	$\mu g/L$	EPA624	23.0	20.00	0	115	60-157				12/2/2010 1021h
										Daport	Date: 12	6/2010 Page 1/17 of 16

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Kyle F. Gross

Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

**Client:** Utah DEQ DERR

**Lab Set ID:** 1011529 Red Butte **Project:** 

Jim Harris **Contact: MSVOA** Dept:

QC Type: MS **RPD** Amount Original

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Amount	%REC	Limits		mit Qual	Date Analyzed
1011529-006BMS	Chlorobenzene	μg/L	EPA624	17.6	20.00	0	88.2	63-140			12/2/2010 1021h
1011529-006BMS	Chloroethane	$\mu g/L$	EPA624	19.3	20.00	0	96.4	41-173			12/2/2010 1021h
1011529-006BMS	Chloroform	$\mug/L$	EPA624	19.9	20.00	0	99.4	50-146			12/2/2010 1021h
1011529-006BMS	Chloromethane	$\mu g/L$	EPA624	16.0	20.00	0	80.2	10-138			12/2/2010 1021h
1011529-006BMS	Chloroprene	$\mug/L$	EPA624	21.3	20.00	0	107	10-161			12/2/2010 1021h
1011529-006BMS	cis-1,2-Dichloroethene	$\mug/L$	EPA624	19.7	20.00	0	98.3	72-137			12/2/2010 1021h
1011529-006BMS	cis-1,3-Dichloropropene	$\mug/L$	EPA624	19.0	40.00	0	47.4	10-134			12/2/2010 1021h
1011529-006BMS	Cyclohexane	$\mug/L$	EPA624	25.3	20.00	0	127	35-230			12/2/2010 1021h
1011529-006BMS	Cyclohexanone	$\mug/L$	EPA624	< 50.0	40.00	0	50.9	10-374			12/2/2010 1021h
1011529-006BMS	Dibromochloromethane	$\mug/L$	EPA624	17.0	20.00	0	85.0	68-135			12/2/2010 1021h
1011529-006BMS	Dibromomethane	$\mug/L$	EPA624	18.6	20.00	0	93.0	74-120			12/2/2010 1021h
1011529-006BMS	Dichlorodifluoromethane	$\mug/L$	EPA624	15.2	20.00	0	76.1	10-150			12/2/2010 1021h
1011529-006BMS	Ethyl acetate	$\mug/L$	EPA624	39.6	40.00	0	99.0	50-155			12/2/2010 1021h
1011529-006BMS	Ethyl ether	$\mug/L$	EPA624	20.2	20.00	0	101	45-146			12/2/2010 1021h
1011529-006BMS	Ethyl methacrylate	$\mug/L$	EPA624	16.4	20.00	0	82.2	77-151			12/2/2010 1021h
1011529-006BMS	Ethylbenzene	$\mu g/L$	EPA624	18.2	20.00	0	90.8	69-133			12/2/2010 1021h
1011529-006BMS	Hexachlorobutadiene	$\mu g/L$	EPA624	15.8	20.00	0	79.0	35-213			12/2/2010 1021h
1011529-006BMS	Iodomethane	$\mu g/L$	EPA624	17.5	20.00	0	87.7	10-233			12/2/2010 1021h
1011529-006BMS	Isobutyl alcohol	$\mug/L$	EPA624	< 100	80.00	0	97.6	12-202			12/2/2010 1021h
1011529-006BMS	Isopropyl acetate	$\mu g/L$	EPA624	19.2	20.00	0	96.1	55-145			12/2/2010 1021h
1011529-006BMS	Isopropyl alcohol	$\mug/L$	EPA624	62.1	80.00	0	77.6	12-250			12/2/2010 1021h
1011529-006BMS	Isopropylbenzene	$\mug/L$	EPA624	18.0	20.00	0	90.2	60-147			12/2/2010 1021h
1011529-006BMS	Isopropyltoluene	$\mug/L$	EPA624	17.2	20.00	0	85.9	73-156			12/2/2010 1021h
1011529-006BMS	m,p-Xylene	μg/L	EPA624	37.4	40.00	0	93.5	70-130			12/2/2010 1021h
									т.	) D-4 10	/C/2010 D 140 -£160

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QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMS	Methacrylonitrile	μg/L	EPA624	16.8	20.00	0	83.8	79-123	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,		Q	12/2/2010 1021h
	•											
1011529-006BMS	Methyl Acetate	μg/L	EPA624	28.9	20.00	0	145	5-398				12/2/2010 1021h
1011529-006BMS	Methyl methacrylate	μg/L	EPA624	16.7	20.00	0	83.6	55-128				12/2/2010 1021h
1011529-006BMS	Methyl tert-butyl ether	$\mu g/L$	EPA624	18.6	20.00	0	93.0	37-189				12/2/2010 1021h
1011529-006BMS	Methylcyclohexane	$\mu g/L$	EPA624	25.0	20.00	0	125	65-175				12/2/2010 1021h
1011529-006BMS	Methylene chloride	$\mug/L$	EPA624	14.2	20.00	0	71.1	55-138				12/2/2010 1021h
1011529-006BMS	n-Amyl acetate	$\mug/L$	EPA624	11.3	20.00	0	56.6	10-187				12/2/2010 1021h
1011529-006BMS	Naphthalene	$\mu g/L$	EPA624	16.1	20.00	0	80.5	41-131				12/2/2010 1021h
1011529-006BMS	n-Butyl alcohol	$\mug/L$	EPA624	60.3	80.00	0	75.4	10-226				12/2/2010 1021h
1011529-006BMS	n-Butylbenzene	$\mug/L$	EPA624	16.8	20.00	0	83.8	40-158				12/2/2010 1021h
1011529-006BMS	n-Hexane	$\mu  g/L$	EPA624	24.5	20.00	0	122	10-277				12/2/2010 1021h
1011529-006BMS	n-Octane	$\mu  g/L$	EPA624	16.8	20.00	0	84.1	45-158				12/2/2010 1021h
1011529-006BMS	n-Propylbenzene	$\mug/L$	EPA624	17.7	20.00	0	88.3	67-131				12/2/2010 1021h
1011529-006BMS	o-Xylene	$\mu g/L$	EPA624	17.8	20.00	0	89.0	70-130				12/2/2010 1021h
1011529-006BMS	Pentachloroethane	$\mu g/L$	EPA624	12.2	20.00	0	60.8	10-314				12/2/2010 1021h
1011529-006BMS	Propionitrile	$\mu g/L$	EPA624	< 25.0	20.00	0	81.3	60-132				12/2/2010 1021h
1011529-006BMS	Propyl acetate	$\mu g/L$	EPA624	18.5	20.00	0	92.4	48-143				12/2/2010 1021h
1011529-006BMS	sec-Butylbenzene	$\mug/L$	EPA624	18.2	20.00	0	91.2	72-157				12/2/2010 1021h
1011529-006BMS	Styrene	$\mug/L$	EPA624	17.2	20.00	0	85.8	81-125				12/2/2010 1021h
1011529-006BMS	tert-Butyl alcohol	$\mug/L$	EPA624	37.6	40.00	0	94.0	50-286				12/2/2010 1021h
1011529-006BMS	tert-Butylbenzene	$\mug/L$	EPA624	17.5	20.00	0	87.3	75-157				12/2/2010 1021h
1011529-006BMS	Tetrachloroethene	$\mu  g/L$	EPA624	22.3	20.00	0	111	49-163				12/2/2010 1021h
1011529-006BMS	Tetrahydrofuran	$\mu  g/L$	EPA624	16.6	20.00	0	83.2	43-146				12/2/2010 1021h
1011529-006BMS	Toluene	$\mug/L$	EPA624	18.6	20.00	0	92.9	18-192				12/2/2010 1021h

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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMS	trans-1,2-Dichloroethene	μg/L	EPA624	19.6	20.00	0	98.0	47-146				12/2/2010 1021h
1011529-006BMS	trans-1,3-Dichloropropene	$\mu g/L$	EPA624	19.1	20.00	0	95.6	29-143				12/2/2010 1021h
1011529-006BMS	trans-1,4-Dichloro-2-butene	$\mu g/L$	EPA624	19.0	20.00	0	95.1	20-214				12/2/2010 1021h
1011529-006BMS	Trichloroethene	$\mu g/L$	EPA624	21.0	20.00	0	105	61-153				12/2/2010 1021h
1011529-006BMS	Trichlorofluoromethane	$\mu g/L$	EPA624	20.2	20.00	0	101	56-166				12/2/2010 1021h
1011529-006BMS	Vinyl acetate	$\mu g/L$	EPA624	26.8	40.00	0	67.0	38-121				12/2/2010 1021h
1011529-006BMS	Vinyl chloride	$\mu g/L$	EPA624	18.1	20.00	0	90.6	13-155				12/2/2010 1021h
1011529-006BMS	Xylenes, Total	$\mu g/L$	EPA624	55.2	60.00	0	92.0	42-167				12/2/2010 1021h
1011529-006BMS	Surr: 1,2-Dichloroethane-d4	%REC	EPA624	56.2	50.00		112	77-144				12/2/2010 1021h
1011529-006BMS	Surr: 4-Bromofluorobenzene	%REC	EPA624	47.8	50.00		95.7	80-123				12/2/2010 1021h
1011529-006BMS	Surr: Dibromofluoromethane	%REC	EPA624	52.4	50.00		105	80-124				12/2/2010 1021h
1011529-006BMS	Surr: Toluene-d8	%REC	EPA624	47.0	50.00		94.0	80-125				12/2/2010 1021h

<sup>&</sup>lt;sup>1</sup> - Matrix spike recovery indicates matrix interference. The method is in control as indicated by the LCS.

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### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMSD	1,1,1,2-Tetrachloroethane	μg/L	EPA624	18.5	20.00	0	92.5	74-117	6.94	25		12/2/2010 1040h
1011529-006BMSD	1,1,1-Trichloroethane	$\mu g/L$	EPA624	23.3	20.00	0	116	67-147	6.65	25		12/2/2010 1040h
1011529-006BMSD	1,1,2,2-Tetrachloroethane	$\mu g/L$	EPA624	17.0	20.00	0	85.0	67-119	5.81	25		12/2/2010 1040h
1011529-006BMSD	1,1,2-Trichloro-1,2,2-trifluoroetha	μg/L	EPA624	24.5	20.00	0	123	21-206	1.27	25		12/2/2010 1040h
1011529-006BMSD	1,1,2-Trichloroethane	μg/L	EPA624	18.0	20.00	0	90.2	80-123	4.36	25		12/2/2010 1040h
1011529-006BMSD	1,1-Dichloropropene	μg/L	EPA624	48.1	40.00	0	120	10-140	4.18	25		12/2/2010 1040h
1011529-006BMSD	1,1-Dichloroethane	$\mu g/L$	EPA624	21.5	20.00	0	108	70-130	5.98	25		12/2/2010 1040h
1011529-006BMSD	1,1-Dichloroethene	μg/L	EPA624	26.1	20.00	0	131	62-152	5.27	25		12/2/2010 1040h
1011529-006BMSD	1,2,3-Trichlorobenzene	$\mu g/L$	EPA624	16.7	20.00	0	83.7	67-131	5.90	25		12/2/2010 1040h
1011529-006BMSD	1,2,3-Trichloropropane	μg/L	EPA624	18.0	20.00	0	89.8	62-116	4.97	25		12/2/2010 1040h
1011529-006BMSD	1,2,3-Trimethylbenzene	$\mu$ g/L	EPA624	19.3	20.00	0	96.5	76-140	3.21	25		12/2/2010 1040h
1011529-006BMSD	1,2,4-Trichlorobenzene	μg/L	EPA624	15.5	20.00	0	77.3	58-133	1.24	25		12/2/2010 1040h
1011529-006BMSD	1,2,4-Trimethylbenzene	$\mu g/L$	EPA624	18.6	20.00	0	93.0	79-151	5.81	25		12/2/2010 1040h
1011529-006BMSD	1,2-Dibromo-3-chloropropane	μg/L	EPA624	16.5	20.00	0	82.4	64-129	3.90	25		12/2/2010 1040h
1011529-006BMSD	1,2-Dibromoethane	μg/L	EPA624	18.6	20.00	0	93.3	70-126	7.63	25		12/2/2010 1040h
1011529-006BMSD	1,2-Dichlorobenzene	μg/L	EPA624	18.1	20.00	0	90.7	70-130	3.76	25		12/2/2010 1040h
1011529-006BMSD	1,2-Dichloroethane	μg/L	EPA624	21.5	20.00	0	107	39-162	4.72	25		12/2/2010 1040h
1011529-006BMSD	1,2-Dichloropropane	μg/L	EPA624	19.1	20.00	0	95.7	59-135	2.86	25		12/2/2010 1040h
1011529-006BMSD	1,3,5-Trimethylbenzene	μg/L	EPA624	18.7	20.00	0	93.6	77-151	4.42	25		12/2/2010 1040h
1011529-006BMSD	1,3-Dichlorobenzene	μg/L	EPA624	18.0	20.00	0	90.0	78-134	3.79	25		12/2/2010 1040h
1011529-006BMSD	1,3-Dichloropropane	μg/L	EPA624	18.5	20.00	0	92.5	78-116	7.64	25		12/2/2010 1040h
1011529-006BMSD	1,4-Dichlorobenzene	μg/L	EPA624	17.5	20.00	0	87.5	72-139	5.10	25		12/2/2010 1040h
1011529-006BMSD	1,4-Dioxane	μg/L	EPA624	156	200.0	0	78.2	33-149	13.4	25		12/2/2010 1040h
1011529-006BMSD	2,2-Dichloropropane	μg/L	EPA624	22.4	30.00	0	74.5	13-180	5.56	25		12/2/2010 1040h
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Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMSD	2-Butanone	μg/L	EPA624	27.3	20.00	0	137	10-217	3.65	25		12/2/2010 1040h
1011529-006BMSD	2-Chloroethyl vinyl ether	$\mu g/L$	EPA624	< 5.00	40.00	0	0	32-163	0	25	1	12/2/2010 1040h
1011529-006BMSD	2-Chlorotoluene	$\mug/L$	EPA624	18.5	20.00	0	92.6	79-142	7.45	25		12/2/2010 1040h
1011529-006BMSD	2-Hexanone	$\mug/L$	EPA624	17.0	20.00	0	85.0	50-156	6.37	25		12/2/2010 1040h
1011529-006BMSD	2-Nitropropane	$\mug/L$	EPA624	19.8	20.00	0	99.0	10-243	8.09	25		12/2/2010 1040h
1011529-006BMSD	4-Chlorotoluene	$\mug/L$	EPA624	18.3	20.00	0	91.7	68-128	4.80	25		12/2/2010 1040h
1011529-006BMSD	4-Isopropyltoluene	$\mug/L$	EPA624	17.6	20.00	0	88.1	73-156	2.59	25		12/2/2010 1040h
1011529-006BMSD	4-Methyl-2-pentanone	$\mug/L$	EPA624	18.1	20.00	0	90.4	10-214	4.82	25		12/2/2010 1040h
1011529-006BMSD	Acetone	$\mug/L$	EPA624	13.2	20.00	0	66.2	10-313	3.30	25		12/2/2010 1040h
1011529-006BMSD	Acetonitrile	$\mug/L$	EPA624	37.9	40.00	0	94.7	37-159	0.474	25		12/2/2010 1040h
1011529-006BMSD	Acrolein	$\mug/L$	EPA624	68.0	40.00	0	170	10-325	6.36	25		12/2/2010 1040h
1011529-006BMSD	Acrylonitrile	$\mug/L$	EPA624	19.5	20.00	0	97.5	53-134	11.4	25		12/2/2010 1040h
1011529-006BMSD	Allyl chloride	$\mug/L$	EPA624	20.3	20.00	0	101	10-243	4.23	25		12/2/2010 1040h
1011529-006BMSD	Benzene	$\mug/L$	EPA624	22.0	20.00	0	110	66-145	3.65	25		12/2/2010 1040h
1011529-006BMSD	Benzyl chloride	$\mug/L$	EPA624	17.3	20.00	0	86.7	40-146	3.22	25		12/2/2010 1040h
1011529-006BMSD	Bis(2-chloroisopropyl) ether	$\mug/L$	EPA624	17.7	20.00	0	88.6	54-146	4.79	25		12/2/2010 1040h
1011529-006BMSD	Bromobenzene	$\mug/L$	EPA624	17.5	20.00	0	87.7	78-148	4.73	25		12/2/2010 1040h
1011529-006BMSD	Bromochloromethane	$\mug/L$	EPA624	20.4	20.00	0	102	75-134	4.72	25		12/2/2010 1040h
1011529-006BMSD	Bromodichloromethane	$\mug/L$	EPA624	19.6	20.00	0	98.2	74-121	1.64	25		12/2/2010 1040h
1011529-006BMSD	Bromoform	$\mug/L$	EPA624	17.5	20.00	0	87.4	68-131	5.41	25		12/2/2010 1040h
1011529-006BMSD	Bromomethane	$\mug/L$	EPA624	15.6	20.00	0	77.8	10-185	14.3	25		12/2/2010 1040h
1011529-006BMSD	Butyl acetate	$\mug/L$	EPA624	18.5	20.00	0	92.4	46-178	5.11	25		12/2/2010 1040h
1011529-006BMSD	Carbon disulfide	$\mug/L$	EPA624	34.6	20.00	0	173	21-224	5.10	25		12/2/2010 1040h
1011529-006BMSD	Carbon tetrachloride	$\mug/L$	EPA624	23.8	20.00	0	119	60-157	3.59	25		12/2/2010 1040h
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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMSD	Chlorobenzene	μg/L	EPA624	18.3	20.00	0	91.5	63-140	3.67	25		12/2/2010 1040h
1011529-006BMSD	Chloroethane	μg/L	EPA624	20.1	20.00	0	100	41-173	4.02	25		12/2/2010 1040h
1011529-006BMSD	Chloroform	μg/L	EPA624	20.8	20.00	0	104	50-146	4.38	25		12/2/2010 1040h
1011529-006BMSD	Chloromethane	μg/L	EPA624	16.3	20.00	0	81.7	10-138	1.79	25		12/2/2010 1040h
1011529-006BMSD	Chloroprene	$\mu g/L$	EPA624	22.0	20.00	0	110	10-161	3.00	25		12/2/2010 1040h
1011529-006BMSD	cis-1,2-Dichloroethene	$\mu g/L$	EPA624	21.0	20.00	0	105	72-137	6.50	25		12/2/2010 1040h
1011529-006BMSD	cis-1,3-Dichloropropene	$\mu g/L$	EPA624	19.8	40.00	0	49.6	10-134	4.59	25		12/2/2010 1040h
1011529-006BMSD	Cyclohexane	$\mu g/L$	EPA624	25.7	20.00	0	129	35-230	1.72	25		12/2/2010 1040h
1011529-006BMSD	Cyclohexanone	$\mu g/L$	EPA624	< 50.0	40.00	0	45.4	10-374	0	25		12/2/2010 1040h
1011529-006BMSD	Dibromochloromethane	$\mu g/L$	EPA624	18.5	20.00	0	92.5	68-135	8.40	25		12/2/2010 1040h
1011529-006BMSD	Dibromomethane	$\mug/L$	EPA624	19.5	20.00	0	97.6	74-120	4.83	25		12/2/2010 1040h
1011529-006BMSD	Dichlorodifluoromethane	$\mug/L$	EPA624	15.7	20.00	0	78.7	10-150	3.36	25		12/2/2010 1040h
1011529-006BMSD	Ethyl acetate	$\mug/L$	EPA624	41.1	40.00	0	103	50-155	3.74	25		12/2/2010 1040h
1011529-006BMSD	Ethyl ether	$\mug/L$	EPA624	21.3	20.00	0	107	45-146	5.59	25		12/2/2010 1040h
1011529-006BMSD	Ethyl methacrylate	$\mug/L$	EPA624	17.1	20.00	0	85.4	77-151	3.88	25		12/2/2010 1040h
1011529-006BMSD	Ethylbenzene	$\mug/L$	EPA624	19.1	20.00	0	95.4	69-133	5.05	25		12/2/2010 1040h
1011529-006BMSD	Hexachlorobutadiene	$\mug/L$	EPA624	15.9	20.00	0	79.6	35-213	0.631	25		12/2/2010 1040h
1011529-006BMSD	Iodomethane	$\mug/L$	EPA624	18.2	20.00	0	91.0	10-233	3.69	25		12/2/2010 1040h
1011529-006BMSD	Isobutyl alcohol	$\mug/L$	EPA624	< 100	80.00	0	73.4	12-202	0	25		12/2/2010 1040h
1011529-006BMSD	Isopropyl acetate	$\mug/L$	EPA624	20.0	20.00	0	100	55-145	4.08	25		12/2/2010 1040h
1011529-006BMSD	Isopropyl alcohol	$\mug/L$	EPA624	52.6	80.00	0	65.8	12-250	16.5	25		12/2/2010 1040h
1011529-006BMSD	Isopropylbenzene	$\mug/L$	EPA624	18.9	20.00	0	94.6	60-147	4.76	25		12/2/2010 1040h
1011529-006BMSD	Isopropyltoluene	$\mug/L$	EPA624	17.6	20.00	0	88.1	73-156	2.59	25		12/2/2010 1040h
1011529-006BMSD	m,p-Xylene	$\mug/L$	EPA624	38.9	40.00	0	97.2	70-130	3.85	25		12/2/2010 1040h
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### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMSD	Methacrylonitrile	μg/L	EPA624	18.9	20.00	0	94.4	79-123	12.0	25		12/2/2010 1040h
1011529-006BMSD	Methyl Acetate	μg/L	EPA624	25.4	20.00	0	127	5-398	13.0	25		12/2/2010 1040h
1011529-006BMSD	Methyl methacrylate	μg/L	EPA624	17.4	20.00	0	87.1	55-128	3.98	25		12/2/2010 1040h
1011529-006BMSD	Methyl tert-butyl ether	μg/L	EPA624	15.7	20.00	0	78.5	37-189	16.9	25		12/2/2010 1040h
1011529-006BMSD	Methylcyclohexane	μg/L	EPA624	24.2	20.00	0	121	65-175	3.26	25		12/2/2010 1040h
1011529-006BMSD	Methylene chloride	$\mu g/L$	EPA624	14.6	20.00	0	73.0	55-138	2.64	25		12/2/2010 1040h
1011529-006BMSD	n-Amyl acetate	$\mu g/L$	EPA624	11.8	20.00	0	59.1	10-187	4.41	25		12/2/2010 1040h
1011529-006BMSD	Naphthalene	$\mu g/L$	EPA624	17.0	20.00	0	85.2	41-131	5.61	25		12/2/2010 1040h
1011529-006BMSD	n-Butyl alcohol	$\mu g/L$	EPA624	61.3	80.00	0	76.6	10-226	1.69	25		12/2/2010 1040h
1011529-006BMSD	n-Butylbenzene	$\mu g/L$	EPA624	17.3	20.00	0	86.3	40-158	2.94	25		12/2/2010 1040h
1011529-006BMSD	n-Hexane	$\mu g/L$	EPA624	18.1	20.00	0	90.7	10-277	29.8	25	@	12/2/2010 1040h
1011529-006BMSD	n-Octane	$\mu g/L$	EPA624	16.2	20.00	0	81.0	45-158	3.82	25		12/2/2010 1040h
1011529-006BMSD	n-Propylbenzene	$\mug/L$	EPA624	18.4	20.00	0	91.8	67-131	3.89	25		12/2/2010 1040h
1011529-006BMSD	o-Xylene	$\mug/L$	EPA624	18.5	20.00	0	92.3	70-130	3.64	25		12/2/2010 1040h
1011529-006BMSD	Pentachloroethane	$\mug/L$	EPA624	13.4	20.00	0	66.8	10-314	9.25	25		12/2/2010 1040h
1011529-006BMSD	Propionitrile	$\mug/L$	EPA624	< 25.0	20.00	0	85.7	60-132	0	25		12/2/2010 1040h
1011529-006BMSD	Propyl acetate	$\mug/L$	EPA624	19.2	20.00	0	96.1	48-143	3.98	25		12/2/2010 1040h
1011529-006BMSD	sec-Butylbenzene	$\mug/L$	EPA624	18.7	20.00	0	93.6	72-157	2.60	25		12/2/2010 1040h
1011529-006BMSD	Styrene	$\mug/L$	EPA624	17.7	20.00	0	88.5	81-125	3.16	25		12/2/2010 1040h
1011529-006BMSD	tert-Butyl alcohol	$\mug/L$	EPA624	35.8	40.00	0	89.5	50-286	4.85	25		12/2/2010 1040h
1011529-006BMSD	tert-Butylbenzene	$\mug/L$	EPA624	18.0	20.00	0	90.0	75-157	2.99	25		12/2/2010 1040h
1011529-006BMSD	Tetrachloroethene	$\mug/L$	EPA624	23.2	20.00	0	116	49-163	3.87	25		12/2/2010 1040h
1011529-006BMSD	Tetrahydrofuran	$\mug/L$	EPA624	18.6	20.00	0	93.1	43-146	11.2	25		12/2/2010 1040h
1011529-006BMSD	Toluene	$\mug/L$	EPA624	19.4	20.00	0	96.9	18-192	4.16	25		12/2/2010 1040h
										Report	Date: 12	/6/2010 Page 15/Lof 160

Report Date: 12/6/2010 Page 154 of 162



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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

Contact: Jim Harris
Dept: MSVOA
QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006BMSD	trans-1,2-Dichloroethene	μg/L	EPA624	20.8	20.00	0	104	47-146	6.14	25		12/2/2010 1040h
1011529-006BMSD	trans-1,3-Dichloropropene	μg/L	EPA624	20.0	20.00	0	100	29-143	4.60	25		12/2/2010 1040h
1011529-006BMSD	trans-1,4-Dichloro-2-butene	$\mug/L$	EPA624	19.7	20.00	0	98.6	20-214	3.62	25		12/2/2010 1040h
1011529-006BMSD	Trichloroethene	$\mu g/L$	EPA624	21.7	20.00	0	108	61-153	2.95	25		12/2/2010 1040h
1011529-006BMSD	Trichlorofluoromethane	$\mug/L$	EPA624	21.5	20.00	0	107	56-166	5.85	25		12/2/2010 1040h
1011529-006BMSD	Vinyl acetate	$\mug/L$	EPA624	27.8	40.00	0	69.4	38-121	3.63	25		12/2/2010 1040h
1011529-006BMSD	Vinyl chloride	$\mug/L$	EPA624	19.4	20.00	0	96.8	13-155	6.62	25		12/2/2010 1040h
1011529-006BMSD	Xylenes, Total	$\mug/L$	EPA624	57.3	60.00	0	95.6	42-167	3.79	25		12/2/2010 1040h
1011529-006BMSD	Surr: 1,2-Dichloroethane-d4	%REC	EPA624	56.7	50.00		113	77-144				12/2/2010 1040h
1011529-006BMSD	Surr: 4-Bromofluorobenzene	%REC	EPA624	48.9	50.00		97.7	80-123				12/2/2010 1040h
1011529-006BMSD	Surr: Dibromofluoromethane	%REC	EPA624	52.6	50.00		105	80-124				12/2/2010 1040h
1011529-006BMSD	Surr: Toluene-d8	%REC	EPA624	48.0	50.00		95.9	80-125				12/2/2010 1040h

<sup>@ -</sup> High RPD due to suspected sample non-homogeneity or matrix interference.

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<sup>&</sup>lt;sup>1</sup> - Matrix spike recovery indicates matrix interference. The method is in control as indicated by the LCS.



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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

Dept: WC

QC Type: DUP

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1012041-001BDUP	Chemical Oxygen Demand	mg/L	HACH 8000	419		449.0		-	6.91	20		12/2/2010 1200h
1012042-001BDUP	Chemical Oxygen Demand	mg/L	HACH 8000	482		477.0		-	1.04	20		12/2/2010 1200h

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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

Dept: WC

QC Type: LCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	RI %RPD Li	PD nit Qual	Date Analyzed
LCS-R21077	Chemical Oxygen Demand	mg/L	HACH 8000	306	300.0	0	102	85-115			12/2/2010 1200h
LCS1-R21077	Chemical Oxygen Demand	mg/L	HACH 8000	108	100.0	0	108	85-115			12/2/2010 1200h
LCS2-R21077	Chemical Oxygen Demand	mg/L	HACH 8000	1,050	1,000	0	105	85-115			12/2/2010 1200h
LCS3-R21077	Chemical Oxygen Demand	mg/L	HACH 8000	< 10.0	10.00	0	90.0	85-115			12/2/2010 1200h
LCS-R21036	Total Recoverable Petroleum Hydr	mg/L	E1664A-SG	19.1	20.00	2.400	83.5	64-132			12/1/2010 1354h

Report Date: 12/6/2010 Page 157 of 162



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Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

**Dept:** WC

**QC Type:** MBLK

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
MB-R21077	Chemical Oxygen Demand	mg/L	HACH 8000	< 10.0				-				12/2/2010 1200h
MB-R21036	Total Recoverable Petroleum Hydr	mg/L	E1664A-SG	< 3.00				-				12/1/2010 1354h

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Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

Dept: WC

QC Type: MS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	RPD %RPD Limi	Date Analyzed
1011529-006CMS	Chemical Oxygen Demand	mg/L	HACH 8000	53.0	50.00	0	106	85-115		12/2/2010 1200h
1012039-003AMS	Chemical Oxygen Demand	mg/L	HACH 8000	55.0	50.00	6.000	98.0	85-115		12/2/2010 1200h
1011529-006AMS	Total Recoverable Petroleum Hydr	mg/L	E1664A-SG	20.4	20.00	0	102	64-132		12/1/2010 1354h

Report Date: 12/6/2010 Page 159 of 162



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Laboratory Director

Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

**Dept:** WC

QC Type: MSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
1011529-006CMSD	Chemical Oxygen Demand	mg/L	HACH 8000	53.0	50.00	0	106	85-115	0	10		12/2/2010 1200h
1012039-003AMSD	Chemical Oxygen Demand	mg/L	HACH 8000	52.0	50.00	6.000	92.0	85-115	5.61	10		12/2/2010 1200h
1011529-006AMSD	Total Recoverable Petroleum Hydr	mg/L	E1664A-SG	21.0	20.00	0	105	64-132	2.91	34		12/1/2010 1354h

Report Date: 12/6/2010 Page 160 of 162



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QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

Dept: WC

QC Type: QCS

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	%REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
QCS-R21036	Total Recoverable Petroleum Hyd	r mg/L	E1664A-SG	19.5	20.00	2.400	85.5	64-132				12/1/2010 1354h

Report Date: 12/6/2010 Page 161 of 162



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Jose Rocha

QA Officer

### **QC SUMMARY REPORT**

Client: Utah DEQ DERR

**Lab Set ID:** 1011529 **Project:** Red Butte

**Contact:** Jim Harris

**Dept:** WC

QC Type: QCSD

Sample ID	Analyte	Units	Method	Result	Amount Spiked	Original Amount	% REC	Limits	%RPD	RPD Limit	Qual	Date Analyzed
QCSD-R21036	Total Recoverable Petroleum Hydr	mg/L	E1664A-SG	19.9	20.00	2.400	87.5	64-132	2.03	34		12/1/2010 1354h

Report Date: 12/6/2010 Page 162 of 162

## **American West Analytical Laboratories**



**WORK ORDER Summary** 

Client:

Utah DEQ DERR

Client ID: Project:

Comments:

UTD100

Red Butte

Contact:

Jim Harris QC Level:

LEVEL II+

WO Type:

Work Order:

12/2/2010 Standard

Page 1 of 3

1011529

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					<u> </u>			<u> </u>	Δ
Sample ID	Client Sample ID	Collected Date	Received Date	Date Due	Matrix	Test Code	Sel	Storage	~
1011529-001A	4992095 Red Butte AB Garden	11/30/2010 9:50:00 AM	11/30/2010 3:38:00 PM	12/3/2010	Aqueous	OGF-W-1664SGT		OGFFridge	1
1011529-001B				12/3/2010		624-W	<b>✓</b>	VOCFridge	3
1011529-001C		<u> </u>		12/3/2010		COD-HACH8000		ww - cod	1
1011529-001D				12/3/2010		3510-ORO-PR		Hall-TPH	2
				12/3/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-001E				12/3/2010		3510-TPH-PR		Hall-TPH	
				12/3/2010		8015-W-TPH(1L)	✓	Hall-TPH	
1011529-001F				12/3/2010		3510-SVOA-PR		Walkin- Semi	3
•				12/3/2010		625-W	✓	Walkin- Semi	
		•		12/3/2010		8270-W	V	Walkin- Semi	
1011529-002A	4992085 Red Butte BL Garden	11/30/2010 10:30:00 AM		12/3/2010		OGF-W-1664SGT		OGFFridge	1
1011529-002B			-	12/3/2010		624-W	<b>V</b>	VOCFridge	3
1011529-002C				12/3/2010		COD-HACH8000		ww - cod	1
1011529-002D				12/3/2010		3510-ORO-PR		Hall-TPH	2
				12/3/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-002E	•			12/3/2010		3510-TPH-PR		Hall-TPH	
				12/3/2010		8015-W-TPH(1L)	<b>V</b>	Hall-TPH	
1011529-002F		-		12/3/2010		3510-SVOA-PR		Walkin- Semi	3
				12/3/2010		625-W	V	Walkin- Semi	
				12/3/2010		8270-W	V	Walkin- Semi	
1011529-003A	4992084 Red Butte @ 1500 E.	11/30/2010 12:15:00 PM		12/3/2010		OGF-W-1664SGT		OGFFridge	1
1011529-003B		4		12/3/2010		624-W	<b>V</b>	VOCFridge	3
1011529-003C				12/3/2010		COD-HACH8000		ww - cod	1
1011529-003D				12/3/2010		3510-ORO-PR		Hail-TPH	2
				12/3/2010		8015-W-ORO(1L)		Hall-TPH	-

**WORK ORDER Summary** 

Client:

Utah DEQ DERR

Page 2 of 3

1011529 12/2/2010

Client ID: Project:

UTD100

Contact:

Jim Harris

WO Type:

Work Order:

Standard

Comments:

QC Level: LEVEL II+ Red Butte No Hard Copies. QC 2+ / Report out TIC's for Semi's - Changed to a RUSH, due tomorrow, except for the ORO. Do the best we can for the ORO.;

Sample ID	Client Sample ID	Collected Date	Received Date	Date Due	Matrix	Test Code	Sel	Storage	
1011529-003E	4992084 Red Butte @ 1500 E.	11/30/2010 12:15:00 PM	11/30/2010 3:38:00 PM	12/3/2010	Aqueous	3510-TPH-PR		Hall-TPH	2
				12/3/2010		8015-W-TPH(1L)	V	Hall-TPH	
1011529-003F				12/3/2010		3510-SVOA-PR		Walkin- Semi	3
				12/3/2010		625-W	<b>V</b>	Walkin- Semi	
				12/3/2010		8270-W	<b>V</b>	Walkin- Semi	
1011529-004A	4992050 J.R. Below 900 S. Storm Sewer	11/30/2010 2:05:00 PM		12/3/2010		OGF-W-1664SGT		OGFFridge	1
1011529-004B				12/3/2010		624-W	V	VOCFridge	3
1011529-004C				12/3/2010		COD-HACH8000		ww - cod	1
1011529-004D				12/3/2010		3510-ORO-PR		Hall-TPH	2
				12/3/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-004E				12/3/2010		3510-TPH-PR		Hall-TPH	
				12/3/2010		8015-W-TPH(1L)	<b>✓</b>	Hall-TPH	
1011529-004F				12/3/2010		3510-SVOA-PR		Walkin- Semi	3
				12/3/2010		625-W	<b>✓</b>	Walkin- Semi	
				12/3/2010		8270-W	<b>V</b>	Walkin- Semi	
1011529-005A	4992290 J.R. @ 1700 S.	11/30/2010 2:30:00 PM		12/3/2010		OGF-W-1664SGT		OGFFridge	1
1011529-005B				12/3/2010		624-W	V	VOCFridge	3
1011529-005C				12/3/2010		COD-HACH8000		ww - cod	1
1011529-005D				12/3/2010		3510-ORO-PR		Hall-TPH	2
				12/3/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-005E				12/3/2010		3510-TPH-PR		Hall-TPH	
				12/3/2010		8015-W-TPH(1L)	<b>V</b>	Hall-TPH	
1011529-005F				12/3/2010		3510-SVOA-PR		Walkin- Semi	3
				12/3/2010		625-W	<b>V</b>	Walkin- Semi	
				12/3/2010		8270-W	<b>V</b>	Walkin- Semi	

WORK ORDER Summary

Client:

Utah DEQ DERR

Client ID: Project:

UTD100

Red Butte

Contact: QC Level: LEVEL II+

Jim Harris

WO Type:

Work Order:

12/2/2010 Standard

1011529 Page 3 of 3

Comments:

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Sample ID	Client Sample ID	Collected Date	Received Date	Date Due	Matrix	Test Code	Sel	Storage	
1011529-006A	4992083 Red Butte @ 1100 E.	11/30/2010 1:15:00 PM	11/30/2010 3:38:00 PM	12/3/2010	Aqueous	OGF-W-1664SGT		OGFFridge	3
1011529-006B				12/3/2010		624-W	V	VOCFridge	6
1011529-006C				12/3/2010		COD-HACH8000		ww - cod	1
1011529-006D				12/3/2010		3510-ORO-PR		Hall-TPH	3
				12/3/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-006E				12/3/2010		3510-TPH-PR		Hall-TPH	2
				12/3/2010		8015-W-TPH(1L)	V	Hall-TPH	
1011529-006F				12/3/2010		3510-SVOA-PR		Walkin- Semi	4
				12/3/2010		625-W	<b>V</b>	Walkin- Semi	
				12/3/2010		8270-W	<b>V</b>	Walkin- Semi	

# **American West Analytical Laboratories**

WORK ORDER Summary

Client:

Utah DEQ DERR

Client ID:

UTD100

Project: Comments: Red Butte

No Hard Copies. QC 2+/Report out TIC's for Semi's;

Work Order:

1011529

Page 1 of 3

11/30/2010

WO Type: Standard

ple ID	Collected Date			3.5	m 4 C 1 -	0.1		
	Conected Date	Received Date	Date Due	Matrix	Test Code	Sei	Storage	
d Butte AB Garden	11/30/2010 9:50:00 AM	11/30/2010 3:38:00 PM	12/14/2010	Aqueous	OGF-W-1664SGT		OGFFridge	1
		•	12/14/2010		624-W	<b>✓</b>	VOCFridge	3
		•	12/14/2010		COD-HACH8000		ww - cod	1
			12/14/2010		3510-ORO-PR		Hall-TPH	2
			12/14/2010		8015-W-ORO(1L)		Hall-TPH	
			12/14/2010		3510-TPH-PR		Hall-TPH	
			12/14/2010		8015-W-TPH(1L)	<b>✓</b>	Hall-TPH	
			12/14/2010		3510-SVOA-PR		Walkin- Semi	3
	,		12/14/2010		625-W	<b>✓</b>	Walkin- Semi	
			12/14/2010		8270-W	<b>V</b>	Walkin- Semi	
d Butte BL Garden	11/30/2010 10:30:00 AM		12/14/2010		OGF-W-1664SGT		OGFFridge	1
			12/14/2010		624-W	V	VOCFridge	3
			12/14/2010		COD-HACH8000		ww - cod	1
			12/14/2010		3510-ORO-PR		Hall-TPH	2
		-	12/14/2010		8015-W-ORO(1L)		Hall-TPH	
			12/14/2010		3510-TPH-PR		Hall-TPH	
			12/14/2010		8015-W-TPH(1L)	<b>V</b>	Hall-TPH	
			12/14/2010		3510-SVOA-PR		Walkin- Semi	3
			12/14/2010		625-W	<b>✓</b>	Walkin- Semi	
			12/14/2010		8270-W	✓	Walkin- Semi	
ed Butte @ 1500 E.	11/30/2010 12:15:00 PM		12/14/2010		OGF-W-1664SGT		OGFFridge	1
			12/14/2010		624-W	✓	VOCFridge	3
			12/14/2010		COD-HACH8000		ww - cod	1
			12/14/2010		3510-ORO-PR		Hall-TPH	2
			12/14/2010		8015-W-ORO(1L)		Hall-TPH	
	ed Butte BL Garden	ed Butte BL Garden 11/30/2010 10:30:00 AM	2d Butte BL Garden 11/30/2010 10:30:00 AM	12/14/2010 12/14/2010	12/14/2010   12/	12/14/2010   624-W     12/14/2010   COD-HACH8000     12/14/2010   3510-ORO-PR     12/14/2010   8015-W-ORO(IL)     12/14/2010   3510-TPI-PR     12/14/2010   3510-TPI-PR     12/14/2010   3510-TPI-PR     12/14/2010   3510-SVOA-PR     12/14/2010   625-W     12/14/2010   625-W     12/14/2010   625-W     12/14/2010   624-W     12/14/2010   624-W     12/14/2010   624-W     12/14/2010   3510-ORO-PR     12/14/2010   3510-TPI-PR     12/14/2010   3510-TPI-PR	12/14/2010   624-W	12/14/2010   624-W

Contact:

QC Level:

Jim Harris

ĻEVEL II+

# WORK ORDER Summary

Client:

Utah DEQ DERR

Client ID: Project:

Comments:

UTD100

Red Butte

No Hard Copies. QC 2+ / Report out TIC's for Semi's;

Work Order:

1011529

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11/30/2010

WO Type: Standard

Sample ID	Client Sample ID	Collected Date	Received Date	Date Due	Matrix	Test Code	Sel	Storage	
1011529-003E	4992084 Red Butte @ 1500 E.	11/30/2010 12:15:00 PM	11/30/2010 3:38:00 PM	12/14/2010	Aqueous	3510-TPH-PR		Hall-TPH	2
				12/14/2010		8015-W-TPH(1L)	<b>V</b>	Hall-TPH	
1011529-003F				12/14/2010	,	3510-SVOA-PR		Walkin- Semi	3
				12/14/2010		625-W	✓	Walkin- Semi	
				12/14/2010		8270-W	<b>V</b>	Walkin- Semi	
1011529-004A	4992050 J.R. Below 900 S. Storm Sewer	11/30/2010 2:05:00 PM		12/14/2010		OGF-W-1664SGT		OGFFridge	1
1011529-004B				12/14/2010		624 <b>-</b> W	✓	VOCFridge	3
1011529-004C				12/14/2010		COD-HACH8000		ww - cod	1
1011529-004D				12/14/2010		3510-ORO-PR		Hall-TPH	2
				12/14/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-004E				12/14/2010		3510-TPH-PR		Hall-TPH	
				12/14/2010		8015-W-TPH(1L)	<b>V</b>	Hall-TPH	
1011529-004F				12/14/2010		3510-SVOA-PR		Walkin- Semi	3
				12/14/2010		625-W	<b>✓</b>	Walkin- Semi	,
				12/14/2010		8270-W	<b>✓</b>	Walkin- Semi	
1011529-005A	4992290 J.R. @ 1700 S.	11/30/2010 2:30:00 PM		12/14/2010		OGF-W-1664SGT		OGFFridge	1
1011529-005B	9			12/14/2010		624-W	V	VOCFridge	3
1011529-005C	·			12/14/2010		COD-HACH8000		ww - cod	1
1011529-005D				12/14/2010		3510-ORO-PR		Hall-TPH	2
				12/14/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-005E	,			12/14/2010		3510-TPH-PR		Hall-TPH	
				12/14/2010		8015-W-TPH(1L)	V	Hall-TPH	
1011529-005F				12/14/2010		3510-SVOA-PR		Walkin- Semi	3
	i			12/14/2010		625-W	<b>V</b>	Walkin- Semi	
				12/14/2010		8270-W	V	Walkin- Semi	

Jim Harris

LEVEL II+

Contact:

QC Level:

**WORK ORDER Summary** 

Client:

Utah DEQ DERR

Client ID: Project:

Comments:

UTD100

Red Butte

No Hard Copies. QC 2+/Report out TIC's for Semi's;

Work Order:

1011529

Page 3 of 3

11/30/2010

WO Type:

Standard

Sample ID	Client Sample ID	Collected Date	Received Date	Date Due	Matrix	Test Code	Sel	Storage	
1011529-006A	4992083 Red Butte @ 1100 E.	11/30/2010 1:15:00 PM	11/30/2010 3:38:00 PM	12/14/2010	Aqueous	OGF-W-1664SGT		OGFFridge	3
1011529-006B				12/14/2010		624-W	<b>✓</b>	VOCFridge	6
1011529-006C				12/14/2010		COD-HACH8000		ww - cod	1
1011529-006D				12/14/2010		3510-ORO-PR		Hall-TPH	3
				12/14/2010		8015-W-ORO(1L)		Hall-TPH	
1011529-006E				12/14/2010		3510-TPH-PR		Hall-TPH	2
				12/14/2010		8015-W-TPH(1L)	<b>✓</b>	Hall-TPH	
1011529-006F				12/14/2010		3510-SVOA-PR		Walkin- Semi	4
				12/14/2010		625-W	<b>✓</b>	Walkin- Semi	
				12/14/2010		8270-W	✓	Walkin- Semi	

Contact:

QC Level:

Jim Harris

LEVEL II+

City State  Phone 801 536 - 9360 Fax				4	<b>\</b>	4	A LAB 63 W Salt L	NAI ORA est 3	W LYTI ATOI 600 S City,	outh	Fax Em	(80) (88) x (80)	HA US: 1) 26: 8) 26: 1) 26: wal@	TO 3-868 3-868 3-868	<b>DY</b> 6 6 7		Page	n Arou	ınd	Time	_ of _ (Circ	/ 52 le One	)	ndard
Contact T. HARRIS							T	EST	'S R	EQL	JIRE	D			(	QC LE	VE	L	LA	BORATO	RY USE	ONLY		
E-mail james harris Project Name Un ONTE	=	000				otal)								× 0				_	_	$\sim$	SAI	MPLES W	ERE:	livered
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Sampler Name JAMES HARRIS				Date/Time Collected			*	× (c	13		×O	<b>!</b> 'ا	×		1		3	3+		4		Ambient o Notes:	Chilled	/
Sample ID					Matrix	Number of	TROH	3	580	T1C5	0 120	DR	\ \frac{1}{3}	Frac			C	ОММІ	ENT	ΓS	Ш	Temperatu Received I	•	- 1
	AG GAR	060	11/30 0950			H	1	ì	-	1				ī			1,6	1,65 cfs			1	(Improperly S		
	BL GA		11/30	1030														80			-	Notes:		
4992084 Red Butte @ 15	OOE		4/80	12:15	11	_											1:	76 0	43	; ·		Properly P (Y) Checked a		
4992050 J. R. Below 9005, Stoven San				_		_						$\perp$					9				- 11 - 1	Y Notes:	N	
4992290 J.R.at 1700			11/30														£20				4	*		
499 2083 RED BUTTE @ 1100 E				01315	14								4	_			1,	15	ch'	5	$ \Big \Big _6$	Received \	Vithin	
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### **Preservation Check Sheet**

Sample Set Extension and pH

Sample Set Extension and pri																	
Bottle Type	Preservative	All	Except	Except	Except	Except	Except	Except	Except	Except	Except	Except	Except	Except	Except	Except	Except
	i	OK															
Ammonia	pH <2 H <sub>2</sub> SO <sub>4</sub>																<u> </u>
COD	pH <2 H <sub>2</sub> SO <sub>4</sub>	1					·										
Cyanide	PH >12 NaOH												, ,				
Metals	pH <2 HNO₃	· .															
NO <sub>2</sub> & NO <sub>3</sub>	pH <2 H <sub>2</sub> SO <sub>4</sub>																
Nutrients	pH <2 H <sub>2</sub> SO <sub>4</sub>																
O&G	pH <2 HCL	V											·				
Phenols	pH <2 H <sub>2</sub> SO <sub>4</sub>																
Sulfide	pH > 9NaOH,																
	Zn Acetate					<u> </u>											
TKN	pH <2 H <sub>2</sub> SO <sub>4</sub>					<u> </u>											
TOC	pH <2 H <sub>3</sub> PO <sub>4</sub>							,								•	
TOX	pH <2 H <sub>2</sub> SO <sub>4</sub>													<u> </u>			
T PO <sub>4</sub>	pH <2 H <sub>2</sub> SO <sub>4</sub>				_												
TPH	pH <2 HCL	_							,								
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Procedure:

- Pour a small amount of sample in the sample lid 1)
- Pour sample from Lid gently over wide range pH paper 2)
- .3)
- Do Not dip the pH paper in the sample bottle or lid

  If sample is not preserved properly list its extension and receiving pH in the appropriate column above 4)
- Flag COC, notify client if requested Place client conversation on COC 5)
- 6)
- Samples may be adjusted 7)

Frequency:

All samples requiring preservation