Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-022
Sample ID: SS-10
Matrix: Solid Percent Dry: 79.6 %
Sampled: 10/26/10 14:30

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
styrene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
naphthalene	< 0.3	0.3	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
Surrogate Recovery		**Limits**								
dibromofluoromethane SUR	**94**	78-114	%	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
toluene-D8 SUR	**101**	88-110	%	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
4-bromofluorobenzene SUR	**100**	86-115	%	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B
a,a,a-trifluorotoluene SUR	**85**	70-130	%	1	LMM	10/29/10	3742	11/2/10	8:14	SW5035A8260B

VIL_RESP01947

RL Resource Laboratories, LLC

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-023
Sample ID: Trip Blank
Matrix: Solid Results expressed on a weight as received basis
Sampled: 10/26/10

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
bromomethane	< 0.2	0.2	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
diethyl ether	< 0.5	0.5	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
methylene chloride	< 0.2	0.2	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
tetrahydrofuran (THF)	< 0.5	0.5	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.4	0.4	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
2-hexanone	< 0.5	0.5	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B

VIL_RESP01948

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-023
Sample ID: Trip Blank
Matrix: Solid Results expressed on a weight as received basis
Sampled: 10/26/10

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
styrene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
naphthalene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
Surrogate Recovery		**Limits**								
dibromofluoromethane SUR	**93**	78-114	%	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
toluene-D8 SUR	**100**	88-110	%	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
4-bromofluorobenzene SUR	**103**	86-115	%	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B
a,a,a-trifluorotoluene SUR	**90**	70-130	%	1	LMM	10/28/10	3742	11/2/10	2:45	SW5035A8260B

VIL_RESP01949

RL Resource Laboratories, LLC

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-024
Sample ID: Trip Blank
Matrix: Water
Sampled: 10/26/10

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
dichlorodifluoromethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
m&p-xylenes	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
o-xylene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B

VIL_RESP01950

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-024
Sample ID: Trip Blank
Matrix: Water
Sampled: 10/26/10

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
styrene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
bromoform	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
isopropylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
n-propylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
bromobenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
2-chlorotoluene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
4-chlorotoluene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
tert-butylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
sec-butylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
4-isopropyltoluene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
n-butylbenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
naphthalene	< 5	5	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
Surrogate Recovery		**Limits**								
dibromofluoromethane SUR	**96**	78-114	%	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
toluene-D8 SUR	**102**	88-110	%	1	LMM		1002510	11/2/10	17:34	SW5030B8260B
4-bromofluorobenzene SUR	**103**	86-115	%	1	LMM		1002510	11/2/10	17:34	SW5030B8260B

VIL_RESP01951

RL Resource Laboratories, LLC

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-025
Sample ID: MW-09
Matrix: Water
Sampled: 10/26/10 10:15

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
dichlorodifluoromethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
m&p-xylenes	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
o-xylene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B

VIL_RESP01952

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-025
Sample ID: MW-09
Matrix: Water
Sampled: 10/26/10 10:15

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
styrene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
bromoform	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
isopropylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
n-propylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
bromobenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
2-chlorotoluene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
4-chlorotoluene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
tert-butylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
sec-butylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
4-isopropyltoluene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
n-butylbenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
naphthalene	< 5	5	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
Surrogate Recovery		**Limits**								
dibromofluoromethane SUR	**100**	78-114	%	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
toluene-D8 SUR	**106**	88-110	%	1	LMM		1002523	11/5/10	8:31	SW5030B8260B
4-bromofluorobenzene SUR	**101**	86-115	%	1	LMM		1002523	11/5/10	8:31	SW5030B8260B

VIL_RESP01953

RL Resource Laboratories, LLC

Project ID: 13 Depot Street 16989.4
Job ID: 20304

Sample#: 20304-001
Sample ID: B-11
Matrix: Solid Percent Dry: 77.3 %
Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 14:06

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
Unadjusted C5-C8 Aliphatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
Unadjusted C9-C12 Aliphatics	**36**	7	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
benzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
toluene	**0.2**	0.1	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
naphthalene	< 0.4	0.4	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
C5-C8 Aliphatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
C9-C12 Aliphatics	**34**	7	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
C9-C10 Aromatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**108**	70-130	%	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**105**	70-130	%	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH
a,a,a-trifluorotoluene SUR	**101**	70-130	%	1	LMM	10/28/10	3743	11/12/10	15:24	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01954

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-002
Sample ID: B-12
Matrix: Solid Percent Dry: 66.5 %
Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 14:17

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
Unadjusted C5-C8 Aliphatics	< 10	10	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
Unadjusted C9-C12 Aliphatics	**15**	10	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
methyl t-butyl ether (MTBE)	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
benzene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
toluene	**0.5**	0.2	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
ethylbenzene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
m&p-xylenes	**0.4**	0.2	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
o-xylene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
naphthalene	< 0.5	0.5	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
C5-C8 Aliphatics	< 10	10	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
C9-C12 Aliphatics	**13**	10	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
C9-C10 Aromatics	< 10	10	ug/g	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**112**	70-130	%	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**106**	70-130	%	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH
a,a,a-trifluorotoluene SUR	**111**	70-130	%	1	LMM	10/28/10	3743	11/12/10	14:12	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01955

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-003
Sample ID: B-03
Matrix: Solid Percent Dry: 81.7 %
Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 13:45

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
Unadjusted C5-C8 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
Unadjusted C9-C12 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
benzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
toluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
naphthalene	< 0.3	0.3	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
C5-C8 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
C9-C12 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
C9-C10 Aromatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**104**	70-130	%	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**105**	70-130	%	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH
a,a,a-trifluorotoluene SUR	**100**	70-130	%	1	LMM	10/28/10	3743	11/8/10	17:59	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01956

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-008
Sample ID: B-06
Matrix: Solid Percent Dry: 86.4 %
Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 12:45

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
Unadjusted C5-C8 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
Unadjusted C9-C12 Aliphatics	**6**	6	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
benzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
toluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
naphthalene	< 0.3	0.3	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
C5-C8 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
C9-C12 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
C9-C10 Aromatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**84**	70-130	%	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**83**	70-130	%	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH
a,a,a-trifluorotoluene SUR	**99**	70-130	%	1	LMM	10/28/10	3743	11/8/10	20:20	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01957

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-009
Sample ID: B-02
Matrix: Solid Percent Dry: 80.5 %
Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 10:45

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
Unadjusted C5-C8 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
Unadjusted C9-C12 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
benzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
toluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
naphthalene	< 0.3	0.3	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
C5-C8 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
C9-C12 Aliphatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
C9-C10 Aromatics	< 6	6	ug/g	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**102**	70-130	%	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**100**	70-130	%	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH
a,a,a-trifluorotoluene SUR	**101**	70-130	%	1	LMM	10/28/10	3743	11/8/10	19:44	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01958

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-012
Sample ID: BK-B-01
Matrix: Solid Percent Dry: 77 %
Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 8:35

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
Unadjusted C5-C8 Aliphatics	< 9	9	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
Unadjusted C9-C12 Aliphatics	< 9	9	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
methyl t-butyl ether (MTBE)	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
benzene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
toluene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
ethylbenzene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
m&p-xylenes	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
o-xylene	< 0.2	0.2	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
naphthalene	< 0.4	0.4	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
C5-C8 Aliphatics	< 9	9	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
C9-C12 Aliphatics	< 9	9	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
C9-C10 Aromatics	< 9	9	ug/g	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**103**	70-130	%	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**102**	70-130	%	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH
a,a,a-trifluorotoluene SUR	**82**	70-130	%	1	LMM	10/28/10	3743	11/8/10	18:34	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01959

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-013
Sample ID: B-10
Matrix: Solid Percent Dry: 76.8 %
Samples prepared in methanol within a 1:1 ratio +/- 25% mL MeOH/g soil
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 11:17

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
Unadjusted C5-C8 Aliphatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
Unadjusted C9-C12 Aliphatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
benzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
toluene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
naphthalene	< 0.4	0.4	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
C5-C8 Aliphatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
C9-C12 Aliphatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
C9-C10 Aromatics	< 7	7	ug/g	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**104**	70-130	%	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**102**	70-130	%	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH
a,a,a-trifluorotoluene SUR	**82**	70-130	%	1	LMM	10/28/10	3743	11/8/10	19:09	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01960

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-018
Sample ID: MW-10
Matrix: Water
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 12:45

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
Unadjusted C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
benzene	< 1	1	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
toluene	< 2	2	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
ethylbenzene	< 2	2	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
m&p-xylenes	< 2	2	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
o-xylene	< 2	2	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
naphthalene	< 5	5	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
C9-C10 Aromatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	15:38	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**102**	70-130	%	1	LMM		1002541	11/8/10	15:38	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**104**	70-130	%	1	LMM		1002541	11/8/10	15:38	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01961

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-020
Sample ID: MW-08
Matrix: Water
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 12:15

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
Unadjusted C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
benzene	< 1	1	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
toluene	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
ethylbenzene	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
m&p-xylenes	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
o-xylene	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
naphthalene	< 5	5	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
C9-C10 Aromatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:13	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**106**	70-130	%	1	LMM		1002541	11/8/10	16:13	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**108**	70-130	%	1	LMM		1002541	11/8/10	16:13	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01962

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-021
Sample ID: MW-02
Matrix: Water
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 12:15

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
Unadjusted C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
benzene	< 1	1	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
toluene	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
ethylbenzene	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
m&p-xylenes	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
o-xylene	< 2	2	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
naphthalene	< 5	5	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
C9-C10 Aromatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	16:48	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**106**	70-130	%	1	LMM		1002541	11/8/10	16:48	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**107**	70-130	%	1	LMM		1002541	11/8/10	16:48	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01963

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-025
Sample ID: MW-09
Matrix: Water
Received on ice at 4°C, in satisfactory condition.
Sampled: 10/26/10 10:15

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
Unadjusted C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
benzene	< 1	1	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
toluene	< 2	2	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
ethylbenzene	< 2	2	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
m&p-xylenes	< 2	2	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
o-xylene	< 2	2	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
naphthalene	< 5	5	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
C9-C10 Aromatics	< 100	100	ug/L	1	LMM		1002541	11/8/10	17:23	MAVPH
Surrogate Recovery		**Limits**								
2,5-dibromotoluene as Aromatic SUR	**104**	70-130	%	1	LMM		1002541	11/8/10	17:23	MAVPH
2,5-dibromotoluene as Aliphatic SUR	**105**	70-130	%	1	LMM		1002541	11/8/10	17:23	MAVPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.
C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.
C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

VIL_RESP01964

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-004
Sample ID: SS-07
Matrix: Solid Percent Dry: 89.9 %
Sampled: 10/26/10 13:40

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
Surrogate Recovery		**Limits**								
2-fluorobiphenyl SUR	**85**	43-116	%	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D
o-terphenyl SUR	**113**	33-141	%	1	AJD	10/29/10	3737	11/4/10	10:50	SW3550B8270D

VIL_RESP01965

RL Resource Laboratories, LLC

Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-006
Sample ID: SS-04
Matrix: Solid Percent Dry: 78.5 %
Sampled: 10/26/10 12:45

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Analysis Time	Reference
naphthalene	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
2-methylnaphthalene	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
acenaphthylene	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
acenaphthene	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
dibenzofuran	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
fluorene	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
phenanthrene	**1.2**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
anthracene	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
fluoranthene	**3.5**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
pyrene	**3.0**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
benzo(a)anthracene	**1.5**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
chrysene	**2.0**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
benzo(b)fluoranthene	**2.0**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
benzo(k)fluoranthene	**1.3**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
benzo(a)pyrene	**1.6**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
indeno(1,2,3-cd)pyrene	**0.8**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
dibenzo(a,h)anthracene	< 0.6	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
benzo(g,h,i)perylene	**1.0**	0.6	ug/g	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
Surrogate Recovery		**Limits**								
2-fluorobiphenyl SUR	**69**	43-116	%	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D
o-terphenyl SUR	**89**	33-141	%	1	AJD	10/29/10	3737	11/4/10	14:34	SW3550B8270D

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Project ID: 13 Depot Street 16989.4
Job ID: 20304
Sample#: 20304-007
Sample ID: SS-03
Matrix: Solid Percent Dry: 94 %
Sampled: 10/26/10 13:30

Parameter	Result	Quant Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
naphthalene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
acenaphthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
fluorene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
phenanthrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
anthracene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
fluoranthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
pyrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
benzo(a)anthracene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
chrysene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
benzo(a)pyrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
Surrogate Recovery		**Limits**								
2-fluorobiphenyl SUR	**83**	43-116	%	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D
o-terphenyl SUR	**95**	33-141	%	1	AJD	10/29/10	3737	11/1/10	18:17	SW3550B8270D

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