

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD061521\
 Data File : VD069447.D
 Acq On : 15 Jun 2021 18:36
 Operator : VA/SY
 Sample : M2709-10
 Misc : 5.00G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 0249-AMND-COMP-N(CL-22)

Quant Time: Jun 16 02:01:59 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D061121S.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 11 12:50:13 2021
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.979	168	286561	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.867	114	570109	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.644	117	624961	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.567	152	298789	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.332	65	171888	53.169	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 106.340%	
35) Dibromofluoromethane	7.914	113	116292	27.234	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 54.460%	
50) Toluene-d8	10.338	98	772844	49.485	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	= 98.960%	
62) 4-Bromofluorobenzene	12.626	95	297417	59.350	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	= 118.700%	
Target Compounds						
					Qvalue	
16) Acetone	4.156	43	43629	81.532	ug/l	100
17) Carbon Disulfide	4.414	76	207318	21.089	ug/l	99
39) Methylcyclohexane	9.355	83	62956	10.543	ug/l	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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Instrument :
MSVOA_D
ClientSampleId :
0249-AMND-COMP-N(CL-22)

Quant Time: Jun 16 02:01:59 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D061121S.M
Quant Title : SW846 8260
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