

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD121119\
 Data File : VD064502.D
 Acq On : 11 Dec 2019 16:13
 Operator : VA/SY
 Sample : K6265-02
 Misc : 7.31G/5.00ml/MSVOA D/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP-1A-S2

Quant Time: Dec 12 07:08:19 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D112719S.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 29 07:18:03 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.98	168	219694	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.86	114	347431	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.65	117	462180	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.57	152	1476043	50.00	ug/l	-0.02
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.33	65	120540	60.62	ug/l	0.00
Spiked Amount	50.000		Recovery	=	121.24%	
35) Dibromofluoromethane	7.91	113	115227	52.64	ug/l	0.00
Spiked Amount	50.000		Recovery	=	105.28%	
50) Toluene-d8	10.34	98	1936766	234.88	ug/l	0.00
Spiked Amount	50.000		Recovery	=	469.76%	
62) 4-Bromofluorobenzene	12.64	95	3811938	1438.50	ug/l	0.00
Spiked Amount	50.000		Recovery	=	2877.00%	
Target Compounds						
16) Acetone	4.16	43	364023	905.876	ug/l	100
25) 2-Butanone	7.22	43	632482	1112.948	ug/l	98
29) Tetrahydrofuran	7.57	42	347063	959.898	ug/l	99
40) Benzene	8.35	78	26235	2.874	ug/l	96

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD121119\
Data File : VD064502.D
Acq On : 11 Dec 2019 16:13
Operator : VA/SY
Sample : K6265-02
Misc : 7.31G/5.00ml/MSVOA D/SOIL
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TP-1A-S2

Quant Time: Dec 12 07:08:19 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D112719S.M
Quant Title : SW846 8260
QLast Update : Fri Nov 29 07:18:03 2019
Response via : Initial Calibration

