

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD121818\
 Data File : VD060623.D
 Acq On : 18 Dec 2018 16:10
 Operator : VA/AP
 Sample : J6397-03
 Misc : 6.99g/5ml/MSVOA D/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 EB11 0-2

Quant Time: Dec 19 03:09:10 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D120518S.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 06 01:49:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	6.40	168	1685736	50.00	ug/l	-0.02
34) 1,4-Difluorobenzene	7.44	114	2346451	50.00	ug/l	-0.02
63) Chlorobenzene-d5	11.27	117	1864116	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.37	152	951355	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.79	65	480638	40.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	80.46%	
35) Dibromofluoromethane	6.32	113	653647	37.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	75.86%	
50) Toluene-d8	9.45	98	1917162	38.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	76.88%	
62) 4-Bromofluorobenzene	12.43	95	672723	37.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	75.40%	
Target Compounds						
16) Acetone	3.16	43	39787	20.774	ug/l	97
20) Methylene Chloride	3.73	84	131453	3.297	ug/l	95
40) Benzene	6.80	78	80404	1.372	ug/l	94
52) Toluene	9.54	92	602652	16.449	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD121818\
Data File : VD060623.D
Acq On : 18 Dec 2018 16:10
Operator : VA/AP
Sample : J6397-03
Misc : 6.99g/5ml/MSVOA D/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EB11_0-2

Quant Time: Dec 19 03:09:10 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120518S.M
Quant Title : SW846 8260
QLast Update : Thu Dec 06 01:49:20 2018
Response via : Initial Calibration

