

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW011019\
 Data File : VW008183.D
 Acq On : 11 Jan 2019 08:13
 Operator : SY/AP
 Sample : K1102-02
 Misc : 5.04G/5ML/MSVOA W/SOIL
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 CPSB-18(20-25)

Quant Time: Jan 11 11:38:21 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
 Quant Title : SW846 8260
 QLast Update : Wed Jan 09 04:40:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.95	168	109730	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.84	114	188517	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.63	117	170479	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.56	152	78423	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.31	65	61791	50.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.20%	
35) Dibromofluoromethane	7.88	113	57533	52.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.66%	
50) Toluene-d8	10.32	98	243369	52.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.50%	
62) 4-Bromofluorobenzene	12.62	95	78235	47.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.60%	
Target Compounds						
16) Acetone	4.14	43	10484	28.560	ug/l	98
17) Carbon Disulfide	4.38	76	10009	2.666	ug/l #	92
39) Methylcyclohexane	9.34	83	240566	97.699	ug/l	94
52) Toluene	10.38	92	6856	2.038	ug/l	94

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW011019\
Data File : VW008183.D
Acq On : 11 Jan 2019 08:13
Operator : SY/AP
Sample : K1102-02
Misc : 5.04G/5ML/MSVOA W/SOIL
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CPSB-18(20-25)

Quant Time: Jan 11 11:38:21 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260
QLast Update : Wed Jan 09 04:40:48 2019
Response via : Initial Calibration

