

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD010220\
 Data File : VD064718.D
 Acq On : 02 Jan 2020 15:32
 Operator : VA/SY
 Sample : K6465-06
 Misc : 8.33G/5.00ml/MSVOA D/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 GP-9-(8-10)

Quant Time: Jan 03 01:20:12 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D122419S.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 24 13:19:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.98	168	507019	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.87	114	782002	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.65	117	665534	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.58	152	279637	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.33	65	230015	51.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.90%	
35) Dibromofluoromethane	7.92	113	242985	50.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.30%	
50) Toluene-d8	10.34	98	921203	50.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.26%	
62) 4-Bromofluorobenzene	12.64	95	260683	44.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.74%	
Target Compounds						
16) Acetone	4.16	43	26640	29.297	ug/l #	81
17) Carbon Disulfide	4.40	76	98876	6.855	ug/l	99
25) 2-Butanone	7.23	43	8535	6.945	ug/l	95
30) Chloroform	7.71	83	10973	1.256	ug/l	96

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD010220\
Data File : VD064718.D
Acq On : 02 Jan 2020 15:32
Operator : VA/SY
Sample : K6465-06
Misc : 8.33G/5.00ml/MSVOA D/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
GP-9-(8-10)

Quant Time: Jan 03 01:20:12 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D122419S.M
Quant Title : SW846 8260
QLast Update : Tue Dec 24 13:19:06 2019
Response via : Initial Calibration

