

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031622\
 Data File : VD072215.D
 Acq On : 16 Mar 2022 13:43
 Operator : VA/SY
 Sample : N1918-12
 Misc : 5.11G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 JDSH-TB-STOCKPILE-B

Quant Time: Mar 16 15:01:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
 Quant Title : SW846 8260
 QLast Update : Mon Mar 14 05:18:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.973	168	470833	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	774305	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	829047	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	389211	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	231447	44.109	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	88.220%
35) Dibromofluoromethane	7.914	113	238690	43.806	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	87.620%
50) Toluene-d8	10.332	98	971961	48.784	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	97.560%
62) 4-Bromofluorobenzene	12.620	95	402882	52.042	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	104.080%
Target Compounds						
					Qvalue	
16) Acetone	4.156	43	27030	23.726	ug/l #	85
17) Carbon Disulfide	4.409	76	29445	4.330	ug/l	98
18) Methyl Acetate	4.715	43	102189	36.602	ug/l	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031622\
Data File : VD072215.D
Acq On : 16 Mar 2022 13:43
Operator : VA/SY
Sample : N1918-12
Misc : 5.11G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
JDSH-TB-STOCKPILE-B

Quant Time: Mar 16 15:01:03 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
Quant Title : SW846 8260
QLast Update : Mon Mar 14 05:18:06 2022
Response via : Initial Calibration

