

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD111121\
 Data File : VD070996.D
 Acq On : 11 Nov 2021 17:24
 Operator : VA/SY
 Sample : M4583-04
 Misc : 5.54G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VWE-B7-110621-0-1

Quant Time: Nov 12 08:03:06 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D110921S.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 10 03:55:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.973	168	348776	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	607531	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	572275	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	239465	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	198083	53.342	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 106.680%	
35) Dibromofluoromethane	7.914	113	191294	55.800	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 111.600%	
50) Toluene-d8	10.332	98	784642	58.444	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	= 116.880%	
62) 4-Bromofluorobenzene	12.620	95	269068	55.661	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	= 111.320%	
Target Compounds						
					Qvalue	
16) Acetone	4.150	43	1042362	976.911	ug/l	91
17) Carbon Disulfide	4.409	76	38247	3.577	ug/l #	87
18) Methyl Acetate	4.715	43	8315	2.610	ug/l	92
25) 2-Butanone	7.209	43	41368	49.687	ug/l	92
39) Methylcyclohexane	9.350	83	75688	10.804	ug/l	98
68) m/p-Xylenes	11.844	106	11194	1.549	ug/l	93
84) 1,2,4-Trimethylbenzene	13.255	105	19069	1.371	ug/l	96

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

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