

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD080321\
 Data File : VD069969.D
 Acq On : 03 Aug 2021 16:39
 Operator : VA/SY
 Sample : M3268-10
 Misc : 5.40G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 20210507-93

Quant Time: Aug 04 01:33:13 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D072221S.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 23 04:51:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.979	168	524952	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	1083505	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	958784	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.567	152	330422	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	358114	59.316	ug/l	0.00
Spiked Amount 50.000	Range 50 - 163		Recovery = 118.640%			
35) Dibromofluoromethane	7.914	113	361534	52.930	ug/l	0.00
Spiked Amount 50.000	Range 54 - 147		Recovery = 105.860%			
50) Toluene-d8	10.338	98	1361390	51.690	ug/l	0.00
Spiked Amount 50.000	Range 49 - 140		Recovery = 103.380%			
62) 4-Bromofluorobenzene	12.626	95	425070	48.292	ug/l	0.00
Spiked Amount 50.000	Range 25 - 144		Recovery = 96.580%			
Target Compounds					Qvalue	
16) Acetone	4.156	43	120413	110.406	ug/l	93
17) Carbon Disulfide	4.409	76	77263	3.598	ug/l #	93
25) 2-Butanone	7.209	43	43576	29.462	ug/l #	89
73) Isopropylbenzene	12.473	105	31573	1.330	ug/l	98
84) 1,2,4-Trimethylbenzene	13.261	105	28393	1.417	ug/l	95

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

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