

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
 Data File : VD073812.D
 Acq On : 07 Jul 2022 16:58
 Operator : VA/SY
 Sample : N3593-03
 Misc : 4.38G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 20220483C-1C

Quant Time: Jul 08 06:33:44 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D070622S.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 08 05:39:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.973	168	99799	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.855	114	224767	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	204436	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	89633	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.320	65	80328	65.418	ug/l	0.00
Spiked Amount	50.000	Range 50 - 163	Recovery	= 130.840%		
35) Dibromofluoromethane	7.908	113	74526	54.038	ug/l	0.00
Spiked Amount	50.000	Range 54 - 147	Recovery	= 108.080%		
50) Toluene-d8	10.332	98	281335	60.217	ug/l	0.00
Spiked Amount	50.000	Range 78 - 12	Recovery	= 120.440%#		
62) 4-Bromofluorobenzene	12.620	95	99286	51.701	ug/l	0.00
Spiked Amount	50.000	Range 50 - 14	Recovery	= 103.400%#		
Target Compounds						
					Qvalue	
16) Acetone	4.156	43	200566	974.012	ug/l	96
17) Carbon Disulfide	4.403	76	66292	61.903	ug/l	98
25) 2-Butanone	7.208	43	73614	182.489	ug/l	93
39) Methylcyclohexane	9.349	83	14925	8.192	ug/l	91
40) Benzene	8.344	78	5841	1.120	ug/l #	88
84) 1,2,4-Trimethylbenzene	13.255	105	6640	1.163	ug/l	96

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

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