

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070722\
 Data File : VD073817.D
 Acq On : 07 Jul 2022 19:28
 Operator : VA/SY
 Sample : N3593-10
 Misc : 4.77G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 20220486A-4A

Quant Time: Jul 08 06:35:37 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	94758	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.855	114	212109	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.632	117	194257	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	84819	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.320	65	78345	67.197	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 134.400%	
35) Dibromofluoromethane	7.908	113	72258	55.520	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 111.040%	
50) Toluene-d8	10.332	98	270008	61.242	ug/l	0.00
Spiked Amount	50.000	Range	78 - 12	Recovery	= 122.480%#	
62) 4-Bromofluorobenzene	12.620	95	97416	53.754	ug/l	0.00
Spiked Amount	50.000	Range	50 - 14	Recovery	= 107.500%#	
Target Compounds						
					Qvalue	
16) Acetone	4.156	43	47156	222.991	ug/l	95
17) Carbon Disulfide	4.397	76	58048	56.548	ug/l	97
25) 2-Butanone	7.208	43	14582	38.072	ug/l #	90
39) Methylcyclohexane	9.344	83	7133	4.149	ug/l #	90
73) Isopropylbenzene	12.467	105	9133	1.399	ug/l	100
84) 1,2,4-Trimethylbenzene	13.255	105	17651	3.268	ug/l	96

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

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