

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD022222\
 Data File : VD072054.D
 Acq On : 22 Feb 2022 13:42
 Operator : VA/SY
 Sample : N1626-02
 Misc : 4.41G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 CC-B1

Quant Time: Feb 23 05:30:19 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D021822S.M
 Quant Title : SW846 8260
 QLast Update : Sat Feb 19 04:05:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.979	168	394633	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	697215	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	649718	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.567	152	254604	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	217724	53.188	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	= 106.380%	
35) Dibromofluoromethane	7.908	113	116750	27.638	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	= 55.280%	
50) Toluene-d8	10.338	98	834206	53.252	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	= 106.500%	
62) 4-Bromofluorobenzene	12.626	95	289989	50.396	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	= 100.800%	
Target Compounds						Qvalue
16) Acetone	4.156	43	149468	168.951	ug/l	96
17) Carbon Disulfide	4.409	76	4676	1.121	ug/l #	84
25) 2-Butanone	7.214	43	54592	44.647	ug/l	98
59) 2-Hexanone	10.979	43	19602	9.814	ug/l	92
80) 1,3,5-Trimethylbenzene	12.949	105	17565	1.203	ug/l	96
84) 1,2,4-Trimethylbenzene	13.261	105	53648	3.739	ug/l	85

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

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