

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031822\
 Data File : VD072299.D
 Acq On : 18 Mar 2022 19:37
 Operator : VA/SY
 Sample : N1962-03
 Misc : 5.03G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 C-18-17-18

Quant Time: Mar 19 05:27:59 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031122S.M
 Quant Title : SW846 8260
 QLast Update : Mon Mar 14 05:18:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.973	168	347626	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	574354	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	628355	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	327190	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	170126	43.914	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	87.820%
35) Dibromofluoromethane	7.908	113	177844	44.001	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	88.000%
50) Toluene-d8	10.332	98	732745	49.581	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	99.160%
62) 4-Bromofluorobenzene	12.626	95	308151	53.663	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	107.320%
Target Compounds						
					Qvalue	
16) Acetone	4.162	43	14589	17.344	ug/l #	79
17) Carbon Disulfide	4.403	76	19568	4.048	ug/l	100
40) Benzene	8.344	78	56849	3.514	ug/l	93
67) Ethyl Benzene	11.738	91	35798	1.439	ug/l #	82
68) m/p-Xylenes	11.844	106	21336	2.258	ug/l	100
84) 1,2,4-Trimethylbenzene	13.255	105	29988	1.366	ug/l	98

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

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