

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD041122\
 Data File : VD072586.D
 Acq On : 11 Apr 2022 14:26
 Operator : VA/SY
 Sample : N2331-04
 Misc : 7.03G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SR-EX1B-BASE-02

Quant Time: Apr 12 01:03:09 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D032922S.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 30 02:44:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.973	168	544391	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	994538	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	949802	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	459362	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	292456	44.409	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	88.820%
35) Dibromofluoromethane	7.909	113	309568	40.053	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	80.100%
50) Toluene-d8	10.332	98	1585039	58.997	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	118.000%
62) 4-Bromofluorobenzene	12.620	95	529350	50.450	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	100.900%
Target Compounds						
					Qvalue	
17) Carbon Disulfide	4.409	76	14218	4.871	ug/l	96
31) Cyclohexane	7.979	56	47010	5.459	ug/l #	56
39) Methylcyclohexane	9.350	83	308811	28.179	ug/l	90
69) o-Xylene	12.173	106	37369	2.755	ug/l	95
73) Isopropylbenzene	12.467	105	78153	2.107	ug/l	98
80) 1,3,5-Trimethylbenzene	12.949	105	216630	6.818	ug/l	100
86) p-Isopropyltoluene	13.502	119	42012	1.241	ug/l #	85

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

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