

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD111621\
 Data File : VD071045.D
 Acq On : 16 Nov 2021 16:06
 Operator : VA/SY
 Sample : M4659-02
 Misc : 6.11G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 HB-1-(7.5-8)-11-11-21

Quant Time: Nov 17 03:24:30 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D110921S.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 10 03:55:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.979	168	68373	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	165320	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	142672	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	60322	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	80768	110.949	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	221.900%#
35) Dibromofluoromethane	7.914	113	60321	64.662	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	129.320%
50) Toluene-d8	10.338	98	161418	44.184	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	88.360%
62) 4-Bromofluorobenzene	12.620	95	62700	47.665	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	95.340%
Target Compounds						
					Qvalue	
16) Acetone	4.156	43	27897	150.463	ug/l #	77
17) Carbon Disulfide	4.409	76	4625	2.206	ug/l #	73
18) Methyl Acetate	4.726	43	4354	6.972	ug/l #	85
20) Methylene Chloride	4.967	84	4930	5.579	ug/l #	86
25) 2-Butanone	7.209	43	10406	57.990	ug/l #	88

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

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