

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD092821\
 Data File : VD070509.D
 Acq On : 28 Sep 2021 18:20
 Operator : VA/SY
 Sample : M3938-03
 Misc : 6.80G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SUB-SLAB-05(6-2.6)

Quant Time: Sep 28 23:44:40 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D090821S.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 09 03:07:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.979	168	334772	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.861	114	671591	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.638	117	582884	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.567	152	198631	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	232706	54.136	ug/l	0.00
Spiked Amount 50.000	Range 50 - 163		Recovery = 108.280%			
35) Dibromofluoromethane	7.909	113	249135	52.115	ug/l	0.00
Spiked Amount 50.000	Range 54 - 147		Recovery = 104.240%			
50) Toluene-d8	10.332	98	839692	48.378	ug/l	0.00
Spiked Amount 50.000	Range 49 - 140		Recovery = 96.760%			
62) 4-Bromofluorobenzene	12.626	95	248093	40.899	ug/l	0.00
Spiked Amount 50.000	Range 25 - 144		Recovery = 81.800%			
Target Compounds					Qvalue	
16) Acetone	4.150	43	727409	947.836	ug/l	100
25) 2-Butanone	7.209	43	35786	35.030	ug/l	98
44) Trichloroethene	9.108	130	8350	1.661	ug/l	90
51) 4-Methyl-2-Pentanone	10.220	43	37433	14.575	ug/l	91
52) Toluene	10.397	92	33871	2.665	ug/l	94
64) Tetrachloroethene	10.873	164	7341	2.069	ug/l	90
68) m/p-Xylenes	11.844	106	15959	1.984	ug/l	88

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

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