

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD060922\
 Data File : VD073513.D
 Acq On : 09 Jun 2022 16:50
 Operator : VA/SY
 Sample : VIBLK
 Misc : 5.00G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VIBLK

Quant Time: Jun 10 06:10:57 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D060922S.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 10 05:32:19 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.973	168	374229	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.855	114	656152	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.632	117	728401	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	374969	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	163973	40.830	ug/l	0.00
Spiked Amount 50.000	Range 50 - 163		Recovery =	81.660%		
35) Dibromofluoromethane	7.908	113	195803	46.406	ug/l	0.00
Spiked Amount 50.000	Range 54 - 147		Recovery =	92.820%		
50) Toluene-d8	10.332	98	1024124	64.197	ug/l	0.00
Spiked Amount 50.000	Range 49 - 140		Recovery =	128.400%		
62) 4-Bromofluorobenzene	12.620	95	380952	65.781	ug/l	0.00
Spiked Amount 50.000	Range 25 - 144		Recovery =	131.560%		
Target Compounds					Qvalue	
16) Acetone	4.132	43	36749	51.970	ug/l #	65
20) Methylene Chloride	4.961	84	29817	3.704	ug/l #	69
25) 2-Butanone	7.202	43	34263	32.931	ug/l #	56
26) 2,2-Dichloropropane	7.185	77	11004	1.632	ug/l #	54
31) Cyclohexane	7.973	56	251061	37.111	ug/l	97
39) Methylcyclohexane	9.349	83	4377149	592.721	ug/l	97
73) Isopropylbenzene	12.467	105	47007	1.809	ug/l	100
78) n-propylbenzene	12.808	91	56980	1.761	ug/l	98

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

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