

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD060722\
 Data File : VD073496.D
 Acq On : 07 Jun 2022 13:09
 Operator : VA/SY
 Sample : VIBLK
 Misc : 5.00G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VIBLK

Quant Time: Jun 08 06:45:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D051922S.M
 Quant Title : SW846 8260
 QLast Update : Thu May 19 15:13:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.973	168	350487	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.855	114	669106	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.632	117	750805	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.561	152	389749	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.326	65	175772	45.380	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	90.760%
35) Dibromofluoromethane	7.908	113	202289	45.816	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	91.640%
50) Toluene-d8	10.332	98	916528	56.313	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	112.620%
62) 4-Bromofluorobenzene	12.620	95	379479	66.438	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	132.880%
Target Compounds						
					Qvalue	
16) Acetone	4.144	43	8486	12.649	ug/l #	63
31) Cyclohexane	7.979	56	13683	2.021	ug/l #	31
39) Methylcyclohexane	9.350	83	405984	54.860	ug/l	100
73) Isopropylbenzene	12.467	105	38165	1.374	ug/l	96
78) n-propylbenzene	12.808	91	46676	1.338	ug/l	98

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

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