

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD031023\
 Data File : VD075493.D
 Acq On : 10 Mar 2023 14:28
 Operator : KP/SY
 Sample : VIBLK
 Misc : 5.00G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 VIBLK

Quant Time: Mar 10 23:06:25 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D031023S.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 10 16:02:39 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.875	168	127454	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.781	114	269337	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.587	117	282738	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.522	152	127886	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.234	65	54422	47.058	ug/l	0.00
Spiked Amount	50.000	Range	50 - 163	Recovery	=	94.120%
35) Dibromofluoromethane	7.805	113	83984	45.528	ug/l	0.00
Spiked Amount	50.000	Range	54 - 147	Recovery	=	91.060%
50) Toluene-d8	10.269	98	339941	46.648	ug/l	0.00
Spiked Amount	50.000	Range	49 - 140	Recovery	=	93.300%
62) 4-Bromofluorobenzene	12.575	95	109701	52.074	ug/l	0.00
Spiked Amount	50.000	Range	25 - 144	Recovery	=	104.140%
Target Compounds						
3) Chloromethane	2.152	50	746	Below Cal	Qvalue	89
16) Acetone	4.028	43	252	Below Cal	#	44
20) Methylene Chloride	4.805	84	5998	Below Cal	#	60

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

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