

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV110525\
 Data File : VV039336.D
 Acq On : 05 Nov 2025 15:18
 Operator : SY/MD
 Sample : VIBLK
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VIBLK

Quant Time: Nov 06 00:16:10 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V100725W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 08 05:01:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	4.596	168	827153	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.529	114	1540457	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.773	117	1418179	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	626525	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.937	65	639120	54.377	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	108.760%
35) Dibromofluoromethane	4.497	113	570030	46.964	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	93.920%
50) Toluene-d8	7.233	98	1846757	44.724	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	89.440%
62) 4-Bromofluorobenzene	9.998	95	570095	38.767	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	77.540%#

Target Compounds	Qvalue

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

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