

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092925\
 Data File : VV039239.D
 Acq On : 29 Sep 2025 18:00
 Operator : SY/MD
 Sample : VIBLK
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VIBLK

Quant Time: Sep 30 01:30:52 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V092225W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 24 08:08:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	4.603	168	449829	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.535	114	872909	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.776	117	853992	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.175	152	381061	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.944	65	397769	61.098	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	122.200%#
35) Dibromofluoromethane	4.503	113	353297	50.345	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	100.700%
50) Toluene-d8	7.236	98	1069842	51.908	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	103.820%
62) 4-Bromofluorobenzene	10.001	95	335542	39.546	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	79.100%#

Target Compounds	Qvalue

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

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