

Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR081125\
 Data File : VR026738.D
 Acq On : 11 Aug 2025 14:42
 Operator : SY/MD
 Sample : VIBLK
 Misc : 5mL/MSVOA_R/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_R
 ClientSampleId :
 VIBLK

Quant Time: Aug 12 02:48:58 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_R\methods\82R080425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 05 02:27:49 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.301	168	289861	50.000	ug/l	0.00
33) 1,4-Difluorobenzene	8.328	114	528523	50.000	ug/l	0.00
62) Chlorobenzene-d5	11.544	117	416063	50.000	ug/l	0.00
71) 1,4-Dichlorobenzene-d4	13.745	152	193086	50.000	ug/l	0.00

System Monitoring Compounds						
32) 1,2-Dichloroethane-d4	7.693	65	178692	48.707	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.420%
34) Dibromofluoromethane	7.199	113	146781	44.901	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.800%
49) Toluene-d8	10.035	98	598149	51.074	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	102.140%
61) 4-Bromofluorobenzene	12.673	95	153357	46.038	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	92.080%

Target Compounds	Qvalue

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

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