

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW111125\
 Data File : VW032420.D
 Acq On : 11 Nov 2025 11:22
 Operator : SY/MD
 Sample : VIBLK
 Misc : 5.00g/5mL/MSVOA_W/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 VIBLK

Quant Time: Nov 12 03:45:26 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W111025S.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 11 03:48:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.965	168	264316	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.855	114	607073	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.635	117	629228	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.556	152	293574	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.319	65	227494	66.760	ug/l	0.00
Spiked Amount	50.000	Range 63 - 155	Recovery	=	133.520%	
35) Dibromofluoromethane	7.904	113	204824	56.353	ug/l	0.00
Spiked Amount	50.000	Range 70 - 134	Recovery	=	112.700%	
50) Toluene-d8	10.331	98	767506	55.042	ug/l	0.00
Spiked Amount	50.000	Range 74 - 123	Recovery	=	110.080%	
62) 4-Bromofluorobenzene	12.617	95	312486	59.324	ug/l	0.00
Spiked Amount	50.000	Range 17 - 146	Recovery	=	118.640%	
Target Compounds						
					Qvalue	
16) Acetone	4.167	43	8218	8.061	ug/l #	73
18) Methyl Acetate	4.710	43	8761	3.470	ug/l	98
20) Methylene Chloride	4.953	84	24764	6.792	ug/l	94

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW111125\
Data File : VW032420.D
Acq On : 11 Nov 2025 11:22
Operator : SY/MD
Sample : VIBLK
Misc : 5.00g/5mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK

Quant Time: Nov 12 03:45:26 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\82W111025S.M
Quant Title : SW846 8260
QLast Update : Tue Nov 11 03:48:21 2025
Response via : Initial Calibration

