

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
 Data File : VN088109.D
 Acq On : 27 Oct 2025 13:31
 Operator : JC\MD
 Sample : Q3464-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW3

Quant Time: Oct 28 01:26:18 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 25 00:15:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	223573	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	506770	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	486491	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.771	152	231113	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	228087	58.997	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	118.000%
35) Dibromofluoromethane	8.147	113	173646	51.005	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.000%
50) Toluene-d8	10.547	98	622675	47.467	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.940%
62) 4-Bromofluorobenzene	12.829	95	233072	50.309	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	100.620%
Target Compounds						
						Qvalue
3) Chloromethane	2.383	50	16871	5.686	ug/l	99
16) Acetone	4.430	43	25969	13.371	ug/l	92
30) Chloroform	7.947	83	48967	8.542	ug/l	99
47) Bromodichloromethane	9.865	83	7823	1.433	ug/l	95

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102725\
Data File : VN088109.D
Acq On : 27 Oct 2025 13:31
Operator : JC\MD
Sample : Q3464-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW3

Quant Time: Oct 28 01:26:18 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102325W.M
Quant Title : SW846 8260
QLast Update : Sat Oct 25 00:15:22 2025
Response via : Initial Calibration

