

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100925\
 Data File : VX048082.D
 Acq On : 09 Oct 2025 13:39
 Operator : JC/MD
 Sample : Q3311-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1S

Quant Time: Oct 10 01:23:24 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091625W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 17 06:39:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	131146	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	272486	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	257011	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	132444	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	112939	54.102	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 108.200%	
35) Dibromofluoromethane	5.373	113	91039	49.818	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.640%	
50) Toluene-d8	8.635	98	309824	48.997	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 98.000%	
62) 4-Bromofluorobenzene	11.067	95	123862	51.613	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 103.220%	
Target Compounds						
					Qvalue	
5) Bromomethane	1.630	94	535	0.483	ug/l #	72
16) Acetone	2.380	43	1874	2.065	ug/l #	84
25) 2-Butanone	4.562	43	10546	9.315	ug/l #	87
27) cis-1,2-Dichloroethene	4.483	96	13327	6.487	ug/l	87
30) Chloroform	5.080	83	908	0.271	ug/l	90
44) Trichloroethene	7.117	130	26022	13.271	ug/l	95
64) Tetrachloroethene	9.263	164	2927961	1747.576	ug/l	100

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

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