

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100925\
 Data File : VX048085.D
 Acq On : 09 Oct 2025 14:41
 Operator : JC/MD
 Sample : Q3311-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-101S

Quant Time: Oct 10 01:25:07 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	123020	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	255257	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	258974	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	132248	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	108754	55.539	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 111.080%	
35) Dibromofluoromethane	5.379	113	86859	50.739	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.480%	
50) Toluene-d8	8.635	98	294044	49.640	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.280%	
62) 4-Bromofluorobenzene	11.061	95	124713	55.476	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 110.960%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.307	50	651	0.389	ug/l	90
16) Acetone	2.380	43	1763	2.071	ug/l	90
25) 2-Butanone	4.568	43	5416	5.100	ug/l	90
27) cis-1,2-Dichloroethene	4.483	96	20366	10.568	ug/l	88
44) Trichloroethene	7.117	130	16310	8.879	ug/l	96
64) Tetrachloroethene	9.257	164	902554	534.613	ug/l	99

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

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