

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101425\
 Data File : VX048159.D
 Acq On : 14 Oct 2025 12:49
 Operator : JC/MD
 Sample : Q3311-14
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-227B

Quant Time: Oct 15 02:42:49 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	127304	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	263250	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	264360	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	131679	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	105919	52.271	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 104.540%	
35) Dibromofluoromethane	5.373	113	90563	51.296	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.600%	
50) Toluene-d8	8.635	98	291356	47.693	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 95.380%	
62) 4-Bromofluorobenzene	11.061	95	126098	54.389	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 108.780%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	1754	1.991	ug/l #	78
25) 2-Butanone	4.556	43	9203	8.374	ug/l	93
27) cis-1,2-Dichloroethene	4.489	96	17007	8.528	ug/l	86
30) Chloroform	5.093	83	14541	4.476	ug/l	95
44) Trichloroethene	7.111	130	29470	15.557	ug/l	91
64) Tetrachloroethene	9.256	164	364388	211.442	ug/l	98

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

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