

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101325\
 Data File : VX048136.D
 Acq On : 13 Oct 2025 13:06
 Operator : JC/MD
 Sample : Q3311-12DL 5X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-214DDL

Quant Time: Oct 14 01:53:45 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	143231	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	299760	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	312119	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	155771	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	123182	54.030	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 108.060%			
35) Dibromofluoromethane	5.379	113	100199	49.841	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 99.680%			
50) Toluene-d8	8.635	98	343716	49.411	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 98.820%			
62) 4-Bromofluorobenzene	11.061	95	149702	56.705	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 113.420%			
Target Compounds						Qvalue
6) Chloroethane	1.703	64	1541	1.208	ug/l #	75
16) Acetone	2.380	43	4854	4.897	ug/l	90
17) Carbon Disulfide	2.526	76	2590	0.556	ug/l #	95
27) cis-1,2-Dichloroethene	4.483	96	1545	0.689	ug/l	95
44) Trichloroethene	7.123	130	594	0.275	ug/l	92
64) Tetrachloroethene	9.256	164	135347	66.520	ug/l	98

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

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