

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101325\
 Data File : VX048134.D
 Acq On : 13 Oct 2025 12:25
 Operator : JC/MD
 Sample : Q3311-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-5BA

Quant Time: Oct 14 01:53:01 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	158387	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	323274	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	331796	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	163977	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	129189	51.243	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 102.480%	
35) Dibromofluoromethane	5.379	113	108295	49.950	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.900%	
50) Toluene-d8	8.635	98	373516	49.790	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.580%	
62) 4-Bromofluorobenzene	11.061	95	159213	55.921	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 111.840%	
Target Compounds						
						Qvalue
16) Acetone	2.380	43	68533	62.519	ug/l	100
17) Carbon Disulfide	2.520	76	3341	0.649	ug/l	96
25) 2-Butanone	4.556	43	14658	10.720	ug/l	92
44) Trichloroethene	7.123	130	2166	0.931	ug/l	79
51) 4-Methyl-2-Pentanone	8.561	43	957	0.310	ug/l #	65
59) 2-Hexanone	9.427	43	1597	0.721	ug/l	87
64) Tetrachloroethene	9.256	164	130759	60.454	ug/l	99

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

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