

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100925\
 Data File : VX048087.D
 Acq On : 09 Oct 2025 15:22
 Operator : JC/MD
 Sample : Q3311-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-5BA

Quant Time: Oct 10 01:26:04 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	148485	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	308267	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	312468	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	154528	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	125754	53.207	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 106.420%	
35) Dibromofluoromethane	5.373	113	103224	49.929	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.860%	
50) Toluene-d8	8.635	98	365877	51.146	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 102.300%	
62) 4-Bromofluorobenzene	11.061	95	154182	56.790	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 113.580%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	56479	54.959	ug/l	100
17) Carbon Disulfide	2.520	76	1798	0.372	ug/l	99
18) Methyl Acetate	2.709	43	1127	0.493	ug/l #	74
25) 2-Butanone	4.556	43	10397	8.111	ug/l	97
27) cis-1,2-Dichloroethene	4.483	96	1236	0.531	ug/l	89
44) Trichloroethene	7.110	130	2451	1.105	ug/l	87
59) 2-Hexanone	9.427	43	1278	0.605	ug/l	98
64) Tetrachloroethene	9.256	164	495261	243.137	ug/l	99

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

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