

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120925\
 Data File : VX048774.D
 Acq On : 09 Dec 2025 11:45
 Operator : JC/MD
 Sample : Q3788-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OWBR-05-135-120525

Quant Time: Dec 10 00:16:56 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.538	168	135193	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	279762	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	295744	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	153431	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	94731	52.207	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 104.420%		
35) Dibromofluoromethane	5.373	113	84981	44.117	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 88.240%		
50) Toluene-d8	8.635	98	304625	45.115	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 90.240%#		
62) 4-Bromofluorobenzene	11.061	95	131405	56.437	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 112.880%		
Target Compounds						
					Qvalue	
3) Chloromethane	1.301	50	2432	1.399	ug/l	99
12) 1,1-Dichloroethene	2.325	96	187750	119.596	ug/l	99
13) Acrolein	2.252	56	302	41.073	ug/l	98
16) Acetone	2.374	43	8653	12.457	ug/l	98
17) Carbon Disulfide	2.520	76	19382	4.251	ug/l	99
21) trans-1,2-Dichloroethene	3.093	96	18284	11.411	ug/l	98
24) 1,1-Dichloroethane	3.605	63	25666	9.344	ug/l	99
27) cis-1,2-Dichloroethene	4.477	96	167466	92.623	ug/l	94
42) 1,2-Dichloroethane	6.074	62	14210	5.133	ug/l	99
44) Trichloroethene	7.111	130	45972	22.527	ug/l	98

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

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