

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091725\
 Data File : VX047609.D
 Acq On : 17 Sep 2025 12:00
 Operator : JC/MD
 Sample : Q3095-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW201D-20250908

Quant Time: Sep 18 02:02:17 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	5.543	168	164440	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	329692	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	318034	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	152611	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	120762	46.137	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	92.280%
35) Dibromofluoromethane	5.373	113	99168	44.850	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.700%
50) Toluene-d8	8.634	98	346603	45.303	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	90.600%#
62) 4-Bromofluorobenzene	11.067	95	145287	50.036	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.080%
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.178	85	18261	10.724	ug/l	95
7) Trichlorofluoromethane	1.898	101	463624	142.533	ug/l	100
9) 1,1,2-Trichlorotrifluo...	2.337	101	4819	2.534	ug/l	97
12) 1,1-Dichloroethene	2.325	96	88304	45.198	ug/l	94
19) Methyl tert-butyl Ether	3.111	73	5994	0.840	ug/l	98
24) 1,1-Dichloroethane	3.611	63	18699	4.503	ug/l	96
27) cis-1,2-Dichloroethene	4.489	96	5493	2.132	ug/l	92
30) Chloroform	5.086	83	31914	7.605	ug/l	94
32) 1,1,1-Trichloroethane	5.373	97	15234	4.280	ug/l	98
44) Trichloroethene	7.110	130	197550	83.268	ug/l	96
55) 1,1,2-Trichloroethane	9.140	97	907	0.389	ug/l #	85
64) Tetrachloroethene	9.256	164	179402	86.532	ug/l	99

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

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