

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091825\
 Data File : VX047649.D
 Acq On : 18 Sep 2025 18:08
 Operator : JC/MD
 Sample : Q3095-24
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW178I-20250910

Quant Time: Sep 19 01:35:05 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	5.538	168	110217	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	210734	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	190251	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	91199	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	99338	56.623	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 113.240%	
35) Dibromofluoromethane	5.373	113	72961	51.625	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 103.240%	
50) Toluene-d8	8.635	98	248250	50.764	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 101.520%	
62) 4-Bromofluorobenzene	11.067	95	96032	51.743	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 103.480%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	3230	4.234	ug/l #	82
19) Methyl tert-butyl Ether	3.111	73	5465	1.143	ug/l	100
29) Tetrahydrofuran	5.013	42	2712	4.738	ug/l #	73
30) Chloroform	5.086	83	1794	0.638	ug/l	90
44) Trichloroethene	7.110	130	3476	2.292	ug/l	98
64) Tetrachloroethene	9.263	164	898	0.724	ug/l	87

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

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