

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX093025\
 Data File : VX047916.D
 Acq On : 30 Sep 2025 17:03
 Operator : JC/MD
 Sample : Q3193-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP12-20250923

Quant Time: Oct 01 01:14:20 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	132206	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	271875	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	282976	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	144063	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	111786	53.120	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 106.240%	
35) Dibromofluoromethane	5.379	113	90914	49.861	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.720%	
50) Toluene-d8	8.635	98	313666	49.716	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.440%	
62) 4-Bromofluorobenzene	11.067	95	134480	56.164	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 112.320%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	1558	1.703	ug/l	100
19) Methyl tert-butyl Ether	3.105	73	1807	0.315	ug/l #	81
24) 1,1-Dichloroethane	3.605	63	4380	1.312	ug/l	97
27) cis-1,2-Dichloroethene	4.495	96	4539	2.192	ug/l	93
30) Chloroform	5.093	83	1488	0.441	ug/l #	56
44) Trichloroethene	7.117	130	27403	14.007	ug/l	95
64) Tetrachloroethene	9.256	164	1764	0.956	ug/l	96

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

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