

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100125\
 Data File : VX047941.D
 Acq On : 01 Oct 2025 15:30
 Operator : JC/MD
 Sample : Q3193-18
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW203D-20250923

Quant Time: Oct 03 01:24:15 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	142268	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	297268	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	304101	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	151044	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	122524	54.105	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 108.220%		
35) Dibromofluoromethane	5.373	113	99694	50.006	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.020%		
50) Toluene-d8	8.635	98	336034	48.712	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 97.420%		
62) 4-Bromofluorobenzene	11.067	95	144029	55.014	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 110.020%		
Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.331	96	897	0.531	ug/l	79
16) Acetone	2.386	43	1883	1.912	ug/l #	82
19) Methyl tert-butyl Ether	3.123	73	1953	0.317	ug/l	92
24) 1,1-Dichloroethane	3.611	63	4216	1.173	ug/l #	85
27) cis-1,2-Dichloroethene	4.483	96	4851	2.177	ug/l	88
30) Chloroform	5.093	83	1166	0.321	ug/l	97
44) Trichloroethene	7.110	130	26386	12.335	ug/l	93
64) Tetrachloroethene	9.263	164	1731	0.873	ug/l	91

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

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