

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100125\
 Data File : VX047942.D
 Acq On : 01 Oct 2025 15:50
 Operator : JC/MD
 Sample : Q3193-20
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT17411-20250924

Quant Time: Oct 03 01:24:39 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	118608	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	242382	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	250160	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	123424	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	94847	50.238	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.480%	
35) Dibromofluoromethane	5.379	113	79617	48.979	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 97.960%	
50) Toluene-d8	8.635	98	269894	47.984	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 95.960%	
62) 4-Bromofluorobenzene	11.067	95	113615	53.223	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 106.440%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.331	96	1496	1.062	ug/l	89
16) Acetone	2.380	43	1697	2.067	ug/l #	85
19) Methyl tert-butyl Ether	3.117	73	1972	0.383	ug/l	93
24) 1,1-Dichloroethane	3.611	63	10503	3.506	ug/l	98
30) Chloroform	5.086	83	1371	0.453	ug/l #	64
44) Trichloroethene	7.111	130	2666	1.529	ug/l	95
64) Tetrachloroethene	9.263	164	894	0.548	ug/l	87

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

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