

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100125\
 Data File : VX047934.D
 Acq On : 01 Oct 2025 13:04
 Operator : JC/MD
 Sample : Q3193-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GM38-RW3-MW1-20250922

Quant Time: Oct 03 01:20:21 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	137828	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	279360	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	288741	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	149951	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	114553	52.215	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.420%
35) Dibromofluoromethane	5.373	113	93231	49.762	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.520%
50) Toluene-d8	8.635	98	315092	48.604	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.200%
62) 4-Bromofluorobenzene	11.061	95	137437	55.861	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	111.720%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	1583	1.659	ug/l	# 89
30) Chloroform	5.080	83	1515	0.431	ug/l	88
44) Trichloroethene	7.117	130	35808	17.812	ug/l	96
64) Tetrachloroethene	9.263	164	2804	1.490	ug/l	88

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

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