

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092925\
 Data File : VX047890.D
 Acq On : 29 Sep 2025 17:26
 Operator : JC/MD
 Sample : Q3213-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-184-GW-420-422

Quant Time: Sep 30 01:24:35 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	128773	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	265004	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	273028	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	140306	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	105554	51.496	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	103.000%
35) Dibromofluoromethane	5.373	113	87688	49.339	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.680%
50) Toluene-d8	8.635	98	298703	48.572	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.140%
62) 4-Bromofluorobenzene	11.067	95	129436	55.459	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.920%
Target Compounds						
						Qvalue
3) Chloromethane	1.300	50	1095	0.625	ug/l	97
16) Acetone	2.380	43	13227	14.841	ug/l	99
19) Methyl tert-butyl Ether	3.123	73	1552	0.278	ug/l	100
25) 2-Butanone	4.574	43	3508	3.156	ug/l #	87

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

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