

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100225\
 Data File : VX047960.D
 Acq On : 02 Oct 2025 12:26
 Operator : JC/MD
 Sample : Q3246-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-184-GW-480-482

Quant Time: Oct 03 03:28:59 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	162247	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	333521	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	336858	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	168922	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	131258	50.825	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.640%
35) Dibromofluoromethane	5.379	113	108404	48.464	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.920%
50) Toluene-d8	8.635	98	373781	48.294	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.580%
62) 4-Bromofluorobenzene	11.061	95	162094	55.184	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.360%
Target Compounds						
						Qvalue
3) Chloromethane	1.307	50	705	0.319	ug/l	95
16) Acetone	2.380	43	4441	3.955	ug/l #	89
17) Carbon Disulfide	2.520	76	3029	0.574	ug/l	98

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

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