

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092325\
 Data File : VX047730.D
 Acq On : 23 Sep 2025 14:07
 Operator : JC/MD
 Sample : Q3141-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-184-GW-60-62

Quant Time: Sep 24 01:32:45 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	141797	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	298818	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	301046	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	155521	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	118593	52.543	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 105.080%		
35) Dibromofluoromethane	5.379	113	97478	48.641	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 97.280%		
50) Toluene-d8	8.635	98	334878	48.293	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 96.580%		
62) 4-Bromofluorobenzene	11.067	95	146404	55.631	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 111.260%		
Target Compounds						
						Qvalue
3) Chloromethane	1.307	50	678	0.351	ug/l	# 85
16) Acetone	2.380	43	10429	10.627	ug/l	97
25) 2-Butanone	4.556	43	3588	2.931	ug/l	94
30) Chloroform	5.080	83	5853	1.617	ug/l	89
40) Benzene	6.031	78	6765	0.761	ug/l	92
52) Toluene	8.708	92	2151	0.393	ug/l	96
60) Dibromochloromethane	9.506	129	881	0.362	ug/l	87
65) Chlorobenzene	10.061	112	1894	0.277	ug/l	97

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

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