

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092525\
 Data File : VX047803.D
 Acq On : 25 Sep 2025 12:07
 Operator : JC/MD
 Sample : Q3165-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-184-GW-260-262

Quant Time: Sep 26 01:25:46 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	129350	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	276238	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	289210	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	151799	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	112618	54.697	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 109.400%	
35) Dibromofluoromethane	5.379	113	92276	49.809	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.620%	
50) Toluene-d8	8.634	98	312248	48.710	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 97.420%	
62) 4-Bromofluorobenzene	11.061	95	137883	56.675	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 113.360%	
Target Compounds						
						Qvalue
3) Chloromethane	1.307	50	515	0.293	ug/l	89
16) Acetone	2.379	43	7471	8.345	ug/l	94
17) Carbon Disulfide	2.514	76	1366	0.325	ug/l #	89

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

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