

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092525\
 Data File : VX047806.D
 Acq On : 25 Sep 2025 13:10
 Operator : JC/MD
 Sample : Q3165-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-184-GW-300-302

Quant Time: Sep 26 01:27:01 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	119390	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	254220	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	261705	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	127218	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	100498	52.883	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	105.760%
35) Dibromofluoromethane	5.373	113	84663	49.658	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.320%
50) Toluene-d8	8.634	98	292394	49.563	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.120%
62) 4-Bromofluorobenzene	11.067	95	121659	54.338	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	108.680%
Target Compounds						
					Qvalue	
3) Chloromethane	1.300	50	1199	0.738	ug/l #	87
16) Acetone	2.380	43	13378	16.190	ug/l #	89
17) Carbon Disulfide	2.520	76	2759	0.710	ug/l	99
25) 2-Butanone	4.580	43	3619	3.511	ug/l #	88

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

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