

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101425\
 Data File : VX048176.D
 Acq On : 14 Oct 2025 18:38
 Operator : JC/MD
 Sample : Q3311-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-219

Quant Time: Oct 15 02:50:39 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	5.537	168	101378	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	188364	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	214540	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	104145	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	79441	49.230	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.460%
35) Dibromofluoromethane	5.379	113	68594	54.299	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	108.600%
50) Toluene-d8	8.635	98	232782	53.254	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	106.500%
62) 4-Bromofluorobenzene	11.061	95	95184	57.376	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	114.760%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	1061	1.512	ug/l #	80
25) 2-Butanone	4.568	43	3074	3.512	ug/l #	87
27) cis-1,2-Dichloroethene	4.483	96	1654	1.041	ug/l	82
30) Chloroform	5.086	83	2151	0.831	ug/l #	78
44) Trichloroethene	7.116	130	2167	1.599	ug/l	94
64) Tetrachloroethene	9.256	164	203167	145.267	ug/l	98

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

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