

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102125\
 Data File : VX048274.D
 Acq On : 21 Oct 2025 16:46
 Operator : JC/MD
 Sample : Q3400-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FRAC-TANK-0760450

Quant Time: Oct 24 04:28:49 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102025W.M
 Quant Title : SW846 8260
 QLast Update : Mon Oct 20 14:13:59 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.538	168	117685	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	236203	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	265928	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	136752	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	91897	55.694	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	111.380%
35) Dibromofluoromethane	5.373	113	76415	47.388	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.780%
50) Toluene-d8	8.635	98	281369	49.781	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.560%
62) 4-Bromofluorobenzene	11.061	95	124629	57.961	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	115.920%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	2910	5.141	ug/l #	84
27) cis-1,2-Dichloroethene	4.483	96	15375	9.707	ug/l	88
44) Trichloroethene	7.111	130	152383	86.725	ug/l	96
45) 1,2-Dichloropropane	7.415	63	27134	17.151	ug/l	99
64) Tetrachloroethene	9.256	164	12958	6.527	ug/l	89

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

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