

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112425\
 Data File : VX048668.D
 Acq On : 24 Nov 2025 16:50
 Operator : JC/MD
 Sample : Q3711-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VNJ-222

Quant Time: Nov 26 00:27:42 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 08 05:08:11 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.537	168	147321	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.750	114	302792	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	329711	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	177161	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	116400	56.703	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	113.400%
35) Dibromofluoromethane	5.373	113	95232	41.555	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	83.100%
50) Toluene-d8	8.634	98	354888	49.652	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.300%
62) 4-Bromofluorobenzene	11.061	95	154133	57.866	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	115.740%
Target Compounds						
					Qvalue	
16) Acetone	2.373	43	15029	15.467	ug/l #	85
18) Methyl Acetate	2.702	43	3072	1.150	ug/l	98
20) Methylene Chloride	2.794	84	8327	4.168	ug/l #	96
25) 2-Butanone	4.556	43	11655	8.108	ug/l	98
43) Isopropyl Acetate	6.330	43	12285	2.023	ug/l	96

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

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