

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060925\
 Data File : VX046578.D
 Acq On : 09 Jun 2025 14:44
 Operator : JC/MD
 Sample : Q2260-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP10-MHG-WC

Quant Time: Jun 10 01:44:06 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060625W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 06 16:56:12 2025
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/10/2025
 Supervised By : Mahesh Dadoda 06/10/2025

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	84479	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	156035	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	142402	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	70692	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	74964	49.877	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	99.760%		
35) Dibromofluoromethane	5.403	113	54142	47.184	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.360%		
50) Toluene-d8	8.653	98	189785	50.167	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	100.340%		
62) 4-Bromofluorobenzene	11.079	95	74611	47.668	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	95.340%		
Target Compounds					Qvalue	
16) Acetone	2.392	43	33483	60.283	ug/l	96
20) Methylene Chloride	2.800	84	3971	3.293	ug/l	92
25) 2-Butanone	4.586	43	4864	5.814	ug/l	94
43) Isopropyl Acetate	6.354	43	10323m	3.701	ug/l	

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

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