

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV110425\
 Data File : VV039311.D
 Acq On : 04 Nov 2025 13:49
 Operator : SY/MD
 Sample : Q3434-02
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 PR132-WC2-20251022

Manual Integrations
 APPROVED

Reviewed By :Amit Patel 11/05/2025
 Supervised By :Mahesh Dadoda 11/05/2025

Quant Time: Nov 05 00:36:50 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V100725W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 08 05:01:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	4.596	168	659851	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.529	114	1187277	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.773	117	1121779	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	499329	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.937	65	591835	63.122	ug/l	0.00
Spiked Amount 50.000	Range 81 - 118		Recovery = 126.240%#			
35) Dibromofluoromethane	4.497	113	519604	55.544	ug/l	0.00
Spiked Amount 50.000	Range 80 - 119		Recovery = 111.080%			
50) Toluene-d8	7.233	98	1642276	51.603	ug/l	0.00
Spiked Amount 50.000	Range 89 - 112		Recovery = 103.200%			
62) 4-Bromofluorobenzene	9.998	95	628109	55.418	ug/l	0.00
Spiked Amount 50.000	Range 85 - 114		Recovery = 110.840%			
Target Compounds					Qvalue	
16) Acetone	2.121	43	77273	7.765	ug/l	96
17) Carbon Disulfide	2.256	76	7768	0.309	ug/l	98
25) 2-Butanone	3.908	43	22325m	3.893	ug/l	

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

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