

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV110425\
 Data File : VV039313.D
 Acq On : 04 Nov 2025 14:34
 Operator : SY/MD
 Sample : Q3434-02RE
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 PR132-WC2-20251022RE

Quant Time: Nov 05 00:37:35 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

Manual Integrations APPROVED

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

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

Internal Standards						
1) Pentafluorobenzene	4.600	168	632494	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.532	114	1126953	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.773	117	1080628	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	492663	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.940	65	588764	65.510	ug/l	0.00
Spiked Amount 50.000	Range 81 - 118		Recovery = 131.020%#			
35) Dibromofluoromethane	4.500	113	512576	57.725	ug/l	0.00
Spiked Amount 50.000	Range 80 - 119		Recovery = 115.460%			
50) Toluene-d8	7.236	98	1626339	53.838	ug/l	0.00
Spiked Amount 50.000	Range 89 - 112		Recovery = 107.680%			
62) 4-Bromofluorobenzene	10.001	95	627116	58.292	ug/l	0.00
Spiked Amount 50.000	Range 85 - 114		Recovery = 116.580%#			
Target Compounds					Qvalue	
16) Acetone	2.124	43	79561	9.137	ug/l	96
17) Carbon Disulfide	2.256	76	8462	0.351	ug/l #	95
25) 2-Butanone	3.915	43	23215m	4.224	ug/l	

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

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