

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101725\
 Data File : VX048214.D
 Acq On : 17 Oct 2025 15:52
 Operator : JC/MD
 Sample : Q3348-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WCS-TP1

Quant Time: Oct 18 00:19:56 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101725W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 18 00:06:31 2025
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 10/20/2025
 Supervised By : Mahesh Dadoda 10/20/2025

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

Internal Standards						
1) Pentafluorobenzene	5.543	168	100604	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.750	114	201041	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.036	117	218558	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	111919	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	85214	54.675	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 109.340%			
35) Dibromofluoromethane	5.379	113	70869	50.792	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 101.580%			
50) Toluene-d8	8.634	98	245094	48.759	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 97.520%			
62) 4-Bromofluorobenzene	11.061	95	107101	58.499	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 117.000%			
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	11828	22.247	ug/l	98
18) Methyl Acetate	2.702	43	4603	4.251	ug/l #	88
20) Methylene Chloride	2.794	84	5179	3.999	ug/l	86
25) 2-Butanone	4.562	43	4757m	7.680	ug/l	
37) Ethyl Acetate	4.714	43	2888	1.839	ug/l #	91
43) Isopropyl Acetate	6.330	43	6843	2.595	ug/l	99

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

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