

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101525\
 Data File : VX048192.D
 Acq On : 15 Oct 2025 15:38
 Operator : JC/MD
 Sample : Q3348-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WCS-TP2

Quant Time: Oct 16 04:07:27 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091625W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 17 06:39:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	117617	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	250641	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	278464	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	139994	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	108626	58.022	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 116.040%			
35) Dibromofluoromethane	5.379	113	86553	51.491	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 102.980%			
50) Toluene-d8	8.635	98	294220	50.585	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 101.160%			
62) 4-Bromofluorobenzene	11.061	95	127925	57.952	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 115.900%			
Target Compounds					Qvalue	
16) Acetone	2.380	43	32914	40.434	ug/l	95
18) Methyl Acetate	2.703	43	6885	3.800	ug/l	93
20) Methylene Chloride	2.794	84	9822	5.700	ug/l	96
25) 2-Butanone	4.550	43	11536	11.361	ug/l #	88
37) Ethyl Acetate	4.727	43	4108	1.562	ug/l #	87
43) Isopropyl Acetate	6.342	43	8787	2.037	ug/l #	92

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

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