

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102125\
 Data File : VX048263.D
 Acq On : 21 Oct 2025 12:58
 Operator : JC/MD
 Sample : Q3369-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-22

Quant Time: Oct 21 13:16:16 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102025W.M
 Quant Title : SW846 8260
 QLast Update : Mon Oct 20 14:13:59 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	122880	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	241629	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	255176	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	130902	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	92553	53.720	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 107.440%	
35) Dibromofluoromethane	5.379	113	78710	47.715	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 95.420%	
50) Toluene-d8	8.635	98	275827	47.705	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 95.400%	
62) 4-Bromofluorobenzene	11.061	95	120181	54.637	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 109.280%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	10779	18.237	ug/l	98
18) Methyl Acetate	2.703	43	4581	3.711	ug/l	99
20) Methylene Chloride	2.788	84	5297	3.678	ug/l	87
25) 2-Butanone	4.568	43	5163	6.901	ug/l #	85
43) Isopropyl Acetate	6.330	43	6162	1.930	ug/l #	87
51) 4-Methyl-2-Pentanone	8.531	43	40705	22.116	ug/l #	52

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

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