

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092525\
 Data File : VV039173.D
 Acq On : 25 Sep 2025 12:10
 Operator : SY/MD
 Sample : Q3164-02
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 DA-2

Quant Time: Sep 26 01:24:03 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V092225W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 24 08:08:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	4.603	168	514758	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.535	114	932817	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.776	117	926008	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.175	152	438853	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.944	65	419629	56.325	ug/l	0.00
Spiked Amount 50.000	Range 81 - 118		Recovery = 112.660%			
35) Dibromofluoromethane	4.503	113	386984	51.604	ug/l	0.00
Spiked Amount 50.000	Range 80 - 119		Recovery = 103.200%			
50) Toluene-d8	7.240	98	1048304	47.596	ug/l	0.00
Spiked Amount 50.000	Range 89 - 112		Recovery = 95.200%			
62) 4-Bromofluorobenzene	10.001	95	406703	44.855	ug/l	0.00
Spiked Amount 50.000	Range 85 - 114		Recovery = 89.720%			
Target Compounds						
					Qvalue	
16) Acetone	2.124	43	108214	18.172	ug/l	99
17) Carbon Disulfide	2.259	76	6286	0.265	ug/l	97
51) 4-Methyl-2-Pentanone	7.169	43	16756	1.358	ug/l	91
65) Chlorobenzene	8.805	112	5564	0.216	ug/l	96

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

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