

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061825\
 Data File : VN087081.D
 Acq On : 18 Jun 2025 13:34
 Operator : JC\MD
 Sample : Q2339-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-C

Quant Time: Jun 19 01:27:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 07 02:12:50 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.236	168	194601	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	394387	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	359527	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	166703	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	141721	54.393	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 108.780%	
35) Dibromofluoromethane	8.177	113	120405	51.516	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 103.040%	
50) Toluene-d8	10.571	98	489109	52.862	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 105.720%	
62) 4-Bromofluorobenzene	12.847	95	169746	49.380	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 98.760%	
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	41726	30.199	ug/l	97
18) Methyl Acetate	5.053	43	24199	6.010	ug/l	96
20) Methylene Chloride	5.295	84	10072	3.894	ug/l	91
37) Ethyl Acetate	7.577	43	16981	3.830	ug/l	98
43) Isopropyl Acetate	8.694	43	10380	1.459	ug/l	95

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

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