

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061725\
 Data File : VN087060.D
 Acq On : 17 Jun 2025 15:26
 Operator : JC\MD
 Sample : Q2319-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-B

Quant Time: Jun 18 01:49:32 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	8.230	168	213420	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	420742	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	382712	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	183788	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	146641	51.319	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 102.640%		
35) Dibromofluoromethane	8.177	113	126756	50.836	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.680%		
50) Toluene-d8	10.565	98	518462	52.525	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 105.040%		
62) 4-Bromofluorobenzene	12.847	95	183759	50.108	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 100.220%		
Target Compounds						
						Qvalue
16) Acetone	4.459	43	37760	24.919	ug/l	93
18) Methyl Acetate	5.048	43	25058	5.674	ug/l	98
20) Methylene Chloride	5.300	84	17542	6.183	ug/l	91
30) Chloroform	7.977	83	6454	1.352	ug/l	97
37) Ethyl Acetate	7.571	43	17959	3.797	ug/l	99
43) Isopropyl Acetate	8.700	43	11228	1.479	ug/l #	93

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

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