

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060925\
 Data File : VN086903.D
 Acq On : 09 Jun 2025 14:25
 Operator : JC\MD
 Sample : Q2192-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-1

Quant Time: Jun 10 03:33:17 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	404204	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	709385	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	610140	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	297993	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	228526	42.227	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	84.460%
35) Dibromofluoromethane	8.177	113	199463	47.446	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.900%
50) Toluene-d8	10.565	98	843736	50.698	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.400%
62) 4-Bromofluorobenzene	12.847	95	302575	48.935	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	97.880%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	5.530	59	10650	7.253	ug/l	95
16) Acetone	4.447	43	111039	38.690	ug/l	96
17) Carbon Disulfide	4.736	76	22072	1.774	ug/l	97
20) Methylene Chloride	5.294	84	19653	3.658	ug/l	92
40) Benzene	8.612	78	273368	13.345	ug/l	100
43) Isopropyl Acetate	8.694	43	31404	2.454	ug/l	93
52) Toluene	10.629	92	160408	12.813	ug/l	100
67) Ethyl Benzene	11.959	91	26103	1.126	ug/l	100
68) m/p-Xylenes	12.065	106	44211	4.983	ug/l	96
69) o-Xylene	12.394	106	34276	4.033	ug/l	97
84) 1,2,4-Trimethylbenzene	13.482	105	38218	2.126	ug/l	99
95) Naphthalene	15.635	128	2153008	96.254	ug/l	100

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

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