

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111423\
 Data File : VN080095.D
 Acq On : 14 Nov 2023 17:38
 Operator : JC\MD
 Sample : 05382-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 S-17

Quant Time: Nov 15 04:44:09 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110123W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 02 05:49:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	185559	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	349865	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	386561	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	176209	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	161623	57.170	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 114.340%	
35) Dibromofluoromethane	8.171	113	128097	51.731	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 103.460%	
50) Toluene-d8	10.571	98	465573	51.073	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.140%	
62) 4-Bromofluorobenzene	12.853	95	188659	62.403	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 124.800%#	
Target Compounds						
						Qvalue
16) Acetone	4.430	43	43051	43.353	ug/l	93
18) Methyl Acetate	5.030	43	4596	0.967	ug/l #	42
20) Methylene Chloride	5.289	84	7323	2.923	ug/l #	79
40) Benzene	8.612	78	137002	12.937	ug/l	98
43) Isopropyl Acetate	8.689	43	152487	26.816	ug/l #	83
52) Toluene	10.630	92	80587	12.836	ug/l	98
67) Ethyl Benzene	11.965	91	16866	1.154	ug/l	97
68) m/p-Xylenes	12.071	106	24966	4.587	ug/l	89
69) o-Xylene	12.400	106	14632	2.792	ug/l	83
80) 1,3,5-Trimethylbenzene	13.177	105	3924	0.377	ug/l	83
84) 1,2,4-Trimethylbenzene	13.482	105	10458	1.022	ug/l	91
95) Naphthalene	15.641	128	79078	8.529	ug/l	99

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

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