

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111523\
 Data File : VN080117.D
 Acq On : 15 Nov 2023 17:52
 Operator : JC\MD
 Sample : 05417-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1118

Quant Time: Nov 15 21:56:42 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	8.230	168	213652	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	403504	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	407836	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	190252	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	168336	51.715	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 103.440%	
35) Dibromofluoromethane	8.171	113	134112	46.961	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 93.920%	
50) Toluene-d8	10.571	98	502078	47.756	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 95.520%	
62) 4-Bromofluorobenzene	12.847	95	198465	56.920	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 113.840%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	39833	34.838	ug/l	100
18) Methyl Acetate	5.036	43	6852	1.253	ug/l #	88
20) Methylene Chloride	5.289	84	9939	3.445	ug/l #	95
37) Ethyl Acetate	7.571	43	4006	1.071	ug/l #	89
40) Benzene	8.606	78	12988	1.063	ug/l	94
43) Isopropyl Acetate	8.694	43	230956	35.216	ug/l #	77
51) 4-Methyl-2-Pentanone	10.447	43	149228	40.416	ug/l	95
52) Toluene	10.635	92	14386	1.987	ug/l	97
67) Ethyl Benzene	11.965	91	16660	1.081	ug/l	93
68) m/p-Xylenes	12.071	106	41011	7.143	ug/l	86
69) o-Xylene	12.400	106	54500	9.857	ug/l	85
95) Naphthalene	15.641	128	129685	12.955	ug/l	99

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

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