

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111723\
 Data File : VN080175.D
 Acq On : 17 Nov 2023 20:03
 Operator : JC\MD
 Sample : 05462-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 S-27

Quant Time: Nov 18 02:38:20 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N111623W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 17 01:19:32 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	162609	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	315914	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	347550	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	162766	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	152528	56.809	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 113.620%	
35) Dibromofluoromethane	8.171	113	117099	50.235	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.460%	
50) Toluene-d8	10.570	98	434647	51.726	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 103.460%	
62) 4-Bromofluorobenzene	12.853	95	177890	56.048	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 112.100%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	30944	34.313	ug/l	92
18) Methyl Acetate	5.041	43	3680	1.270	ug/l #	60
20) Methylene Chloride	5.277	84	90229	47.778	ug/l	95
37) Ethyl Acetate	7.565	43	3096	1.323	ug/l #	74
40) Benzene	8.612	78	138119	14.400	ug/l	100
43) Isopropyl Acetate	8.694	43	137348	30.151	ug/l #	84
51) 4-Methyl-2-Pentanone	10.447	43	24690	10.469	ug/l	90
52) Toluene	10.635	92	88677	14.773	ug/l	97
67) Ethyl Benzene	11.970	91	16543	1.181	ug/l	91
68) m/p-Xylenes	12.070	106	26400	5.167	ug/l	87
69) o-Xylene	12.406	106	18140	3.767	ug/l	87
84) 1,2,4-Trimethylbenzene	13.488	105	11591	1.069	ug/l	81
95) Naphthalene	15.647	128	125280	13.511	ug/l	98

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

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