

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111723\
 Data File : VN080166.D
 Acq On : 17 Nov 2023 16:29
 Operator : JC\MD
 Sample : 05462-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 S-28

Quant Time: Nov 18 02:32:48 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	8.230	168	167824	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	324857	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	349216	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	155094	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	152202	54.926	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 109.860%			
35) Dibromofluoromethane	8.171	113	119485	49.847	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 99.700%			
50) Toluene-d8	10.571	98	442018	51.155	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 102.300%			
62) 4-Bromofluorobenzene	12.853	95	173043	53.020	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 106.040%			
Target Compounds					Qvalue	
16) Acetone	4.430	43	43490	46.727	ug/l	92
20) Methylene Chloride	5.277	84	84604	43.407	ug/l	89
25) 2-Butanone	7.494	43	5876	5.314	ug/l	91
37) Ethyl Acetate	7.571	43	3037	1.262	ug/l #	73
40) Benzene	8.612	78	335824	34.048	ug/l	99
43) Isopropyl Acetate	8.694	43	131610	28.096	ug/l #	84
51) 4-Methyl-2-Pentanone	10.447	43	15801	6.515	ug/l #	80
52) Toluene	10.635	92	213629	34.609	ug/l	96
67) Ethyl Benzene	11.970	91	39642	2.815	ug/l	92
68) m/p-Xylenes	12.070	106	59199	11.531	ug/l	90
69) o-Xylene	12.406	106	33083	6.838	ug/l	87
84) 1,2,4-Trimethylbenzene	13.488	105	16730	1.619	ug/l	89
95) Naphthalene	15.647	128	150795	17.067	ug/l	99

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

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