

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120722\
 Data File : VN075612.D
 Acq On : 07 Dec 2022 17:31
 Operator : JC\MD
 Sample : N5876-02DL 200X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 LIN-VRG-12-02DL

Quant Time: Dec 07 17:58:00 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 23 00:18:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	225261	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	357239	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	309283	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	138425	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	59805	49.460	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	98.920%
35) Dibromofluoromethane	8.171	113	78863	56.196	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	112.400%
50) Toluene-d8	10.571	98	163932	51.865	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.740%
62) 4-Bromofluorobenzene	12.853	95	117345	50.299	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.600%
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	52790	46.805	ug/l	95
25) 2-Butanone	7.489	43	261888	166.317	ug/l	98
29) Tetrahydrofuran	7.847	42	36732	38.350	ug/l	98
40) Benzene	8.612	78	53569	5.717	ug/l	98
52) Toluene	10.635	92	140688	23.300	ug/l	99
67) Ethyl Benzene	11.965	91	34343	3.152	ug/l	99
68) m/p-Xylenes	12.071	106	53904	12.181	ug/l	96
69) o-Xylene	12.400	106	28271	6.430	ug/l	93
84) 1,2,4-Trimethylbenzene	13.482	105	10040	1.191	ug/l	94

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

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