

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN093022\
 Data File : VN074689.D
 Acq On : 01 Oct 2022 04:48
 Operator : JC\MD
 Sample : N4882-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-101

Quant Time: Oct 03 02:11:39 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N092922W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 30 07:44:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.934	168	520005	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.828	114	1010451	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.622	117	1026557	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.551	152	469259	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.292	65	668613	51.304	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	102.600%
35) Dibromofluoromethane	7.869	113	549069	51.338	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.680%
50) Toluene-d8	10.310	98	2122135	47.172	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.340%
62) 4-Bromofluorobenzene	12.604	95	697451	48.893	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.780%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	5.163	59	64401	18.503	ug/l	# 72
19) Methyl tert-butyl Ether	5.428	73	186032	5.019	ug/l	99
30) Chloroform	7.657	83	14240	0.710	ug/l	86
40) Benzene	8.310	78	64602	1.224	ug/l	95
42) 1,2-Dichloroethane	8.387	62	32359	1.825	ug/l	95

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

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