

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN093022\
 Data File : VN074695.D
 Acq On : 01 Oct 2022 07:14
 Operator : JC\MD
 Sample : N4882-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-108

Quant Time: Oct 03 02:12:41 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	7.939	168	486514	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.828	114	916996	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.616	117	900546	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.551	152	401039	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.292	65	636888	52.233	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	104.460%
35) Dibromofluoromethane	7.869	113	503440	51.869	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.740%
50) Toluene-d8	10.310	98	1948296	47.722	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.440%
62) 4-Bromofluorobenzene	12.604	95	618585	47.784	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.560%
Target Compounds						Qvalue
16) Acetone	4.122	43	179191	25.520	ug/l	96
19) Methyl tert-butyl Ether	5.428	73	475150	13.701	ug/l	97
40) Benzene	8.310	78	44478	0.928	ug/l	96
42) 1,2-Dichloroethane	8.386	62	27649	1.718	ug/l	92

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

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