

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072121\
 Data File : VN067840.D
 Acq On : 21 Jul 2021 15:50
 Operator : JC/MD
 Sample : M3089-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-N2-C-BOTTOM-(4-11)

Quant Time: Jul 22 02:12:21 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N070721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jul 07 16:18:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	323869	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	618811	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	627629	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	227489	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	253101	51.710	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	103.420%
35) Dibromofluoromethane	8.024	113	185559	49.632	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.260%
50) Toluene-d8	10.443	98	797782	50.979	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.960%
62) 4-Bromofluorobenzene	12.733	95	270796	48.984	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	97.960%
Target Compounds						
						Qvalue
16) Acetone	4.296	43	37141	24.115	ug/l	95
18) Methyl Acetate	4.864	43	5946	1.374	ug/l	94
20) Methylene Chloride	5.098	84	32834	8.555	ug/l #	87
43) Isopropyl Acetate	8.566	43	37798	3.858	ug/l #	91

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

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