

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
 Data File : VN071013.D
 Acq On : 18 Feb 2022 20:15
 Operator : JC\MD
 Sample : N1599-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SPN-4

Quant Time: Feb 19 00:44:38 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 16 07:52:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.087	168	847276	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1336396	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1184356	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	387616	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	496141	51.262	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	102.520%
35) Dibromofluoromethane	8.022	113	395120	49.708	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.420%
50) Toluene-d8	10.439	98	1601032	50.154	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	100.300%
62) 4-Bromofluorobenzene	12.727	95	522136	45.760	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	91.520%
Target Compounds						
						Qvalue
3) Chloromethane	2.299	50	2721	0.246	ug/l	# 80
16) Acetone	4.287	43	67006	12.571	ug/l	100
37) Ethyl Acetate	7.410	43	273215	19.562	ug/l	99

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

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