

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
 Data File : VN071010.D
 Acq On : 18 Feb 2022 19:03
 Operator : JC\MD
 Sample : N1587-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PRPH-DUP-GW

Quant Time: Feb 19 00:43:29 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	8.081	168	751635	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1184391	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1038989	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	342556	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	451185	52.549	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	105.100%
35) Dibromofluoromethane	8.016	113	354012	50.252	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.500%
50) Toluene-d8	10.439	98	1425153	50.374	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	100.740%
62) 4-Bromofluorobenzene	12.727	95	467304	46.211	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	92.420%
Target Compounds						
						Qvalue
3) Chloromethane	2.305	50	2472	0.252	ug/l	# 84
16) Acetone	4.298	43	13044	2.759	ug/l	90
84) 1,2,4-Trimethylbenzene	13.368	105	8670	0.394	ug/l	99

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

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