

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021822\
 Data File : VN071007.D
 Acq On : 18 Feb 2022 17:51
 Operator : JC\MD
 Sample : N1587-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PRPH-1-GW

Quant Time: Feb 19 00:42:24 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	8.086	168	818944	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1284840	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1124998	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	371122	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	490302	52.411	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 104.820%	
35) Dibromofluoromethane	8.022	113	375219	49.098	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 98.200%	
50) Toluene-d8	10.439	98	1543414	50.289	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 100.580%	
62) 4-Bromofluorobenzene	12.727	95	503046	45.856	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 91.720%	
Target Compounds						
					Qvalue	
16) Acetone	4.298	43	13619	2.643	ug/l	99
68) m/p-Xylenes	11.957	106	6814	0.416	ug/l	90
69) o-Xylene	12.274	106	4202	0.262	ug/l	89
84) 1,2,4-Trimethylbenzene	13.362	105	12204	0.512	ug/l	100

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

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