

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042922\
 Data File : VN072264.D
 Acq On : 29 Apr 2022 12:58
 Operator : JC\MD
 Sample : N2617-16DL 4X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-17IRDL

Quant Time: Apr 30 03:12:21 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042622W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 26 18:50:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	189728	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.963	114	324480	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	279299	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	87510	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	162397	55.754	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	111.500%
35) Dibromofluoromethane	8.022	113	106481	51.270	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	102.540%
50) Toluene-d8	10.439	98	385254	47.208	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	94.420%
62) 4-Bromofluorobenzene	12.727	95	113428	41.875	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	83.760%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	5370	4.276	ug/l	# 74
19) Methyl tert-butyl Ether	5.616	73	235024	33.088	ug/l	97
22) Diisopropyl ether	6.492	45	30893	3.986	ug/l	92
42) 1,2-Dichloroethane	8.528	62	54437	15.164	ug/l	96

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

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