

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061122\
 Data File : VN072788.D
 Acq On : 11 Jun 2022 14:29
 Operator : JC\MD
 Sample : N3322-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 002-35TH-AVE(JUNE)

Quant Time: Jun 13 01:19:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N060622W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Mon Jun 06 14:10:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	7.657	128	32098	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	8.963	114	193431	30.000	ug/l	# 0.00
57) Chlorobenzene-d5	11.745	117	156881	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.434	65	116377	31.791	ug/l	0.00
Spiked Amount	30.000	Range 91 - 110		Recovery	= 105.967%	
60) 4-Bromofluorobenzene	12.727	95	81659	25.307	ug/l	0.00
Spiked Amount	30.000	Range 63 - 112		Recovery	= 84.367%	
63) Toluene-d8	10.439	98	282674	32.527	ug/l	0.00
Spiked Amount	30.000	Range 91 - 112		Recovery	= 108.433%	
Target Compounds						
					Qvalue	
15) Acetone	4.281	58	164953	341.591	ug/l	85
16) Carbon Disulfide	4.534	76	7187	1.344	ug/l	# 90
30) 2-Butanone	7.334	43	34253	12.836	ug/l	# 89
58) 4-Methyl-2-Pentanone	10.328	43	15566	3.308	ug/l	98
62) Toluene	10.504	91	52025	4.891	ug/l	99

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

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