

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
 Data File : VN072777.D
 Acq On : 10 Jun 2022 20:14
 Operator : JC\MD
 Sample : N3293-09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 52322-B

Quant Time: Jun 11 01:56:53 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 09 05:17:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	164134	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.963	114	324880	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	280051	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	105986	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	242939	55.148	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	110.300%
35) Dibromofluoromethane	8.016	113	179586	55.865	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	111.740%
50) Toluene-d8	10.439	98	562279	43.840	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	87.680%
62) 4-Bromofluorobenzene	12.727	95	211686	48.098	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	96.200%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	5.363	59	131921	125.536	ug/l	100
16) Acetone	4.287	43	91143	43.178	ug/l	# 88
18) Methyl Acetate	4.851	43	9017	1.642	ug/l	# 87
25) 2-Butanone	7.334	43	68302	20.115	ug/l	91
51) 4-Methyl-2-Pentanone	10.333	43	8977	1.277	ug/l	88
52) Toluene	10.504	92	14007	1.657	ug/l	98
95) Naphthalene	15.510	128	10871	1.561	ug/l	97

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

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