

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092521\
 Data File : VN068675.D
 Acq On : 25 Sep 2021 18:25
 Operator : JC/MD
 Sample : M3870-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-C

Quant Time: Sep 27 04:46:52 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N092321W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 23 13:52:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	479754	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.971	114	812647	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	731052	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	257232	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	390912	52.273	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 104.540%	
35) Dibromofluoromethane	8.027	113	286716	52.752	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 105.500%	
50) Toluene-d8	10.443	98	1158029	50.246	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 100.500%	
62) 4-Bromofluorobenzene	12.733	95	329772	41.837	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 83.680%	
Target Compounds						
					Qvalue	
16) Acetone	4.291	43	120722	34.555	ug/l	94
18) Methyl Acetate	4.867	43	13355	1.212	ug/l	# 80
20) Methylene Chloride	5.106	84	26085	2.717	ug/l	88
25) 2-Butanone	7.340	43	26343	5.770	ug/l	98
43) Isopropyl Acetate	8.563	43	73939	4.639	ug/l	# 81

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

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