

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092822\
 Data File : VX031750.D
 Acq On : 29 Sep 2022 01:23
 Operator : JC/MD
 Sample : N4863-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-8

Quant Time: Sep 29 04:36:39 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 27 01:44:18 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	126284	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	208679	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	219168	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	122688	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	136223	50.825	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	101.640%	
35) Dibromofluoromethane	5.385	113	114160	45.913	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	91.820%	
50) Toluene-d8	8.646	98	434290	48.177	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	96.360%	
62) 4-Bromofluorobenzene	11.079	95	151559	45.872	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	91.740%	
Target Compounds						
17) Carbon Disulfide	2.513	76	7200	1.073	ug/l	Qvalue 96
19) Methyl tert-butyl Ether	3.111	73	38389	6.440	ug/l	98
67) Ethyl Benzene	10.195	91	2949	0.284	ug/l #	87
70) Styrene	10.658	104	1627	0.245	ug/l	91

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

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