

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061822\
 Data File : VX029585.D
 Acq On : 18 Jun 2022 21:15
 Operator : JC/MD
 Sample : N3290-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-19

Quant Time: Jun 20 01:37:26 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 20 01:31:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	272600	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	456598	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	393707	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	170347	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	156712	43.429	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.860%
35) Dibromofluoromethane	5.391	113	136536	45.885	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.780%
50) Toluene-d8	8.653	98	523550	48.163	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.320%
62) 4-Bromofluorobenzene	11.079	95	186015	45.377	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.760%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.977	59	249288	319.590	ug/l	99
16) Acetone	2.386	43	4702	3.269	ug/l	92
17) Carbon Disulfide	2.514	76	2563	0.358	ug/l #	91
19) Methyl tert-butyl Ether	3.117	73	2583113	269.416	ug/l	100
22) Diisopropyl ether	3.763	45	28437	3.224	ug/l #	89
40) Benzene	6.043	78	10054	0.842	ug/l	94
67) Ethyl Benzene	10.195	91	8534	0.614	ug/l	89
68) m/p-Xylenes	10.305	106	7313	1.349	ug/l	91
69) o-Xylene	10.646	106	1902	0.355	ug/l	98
78) n-propylbenzene	11.311	91	3350	0.233	ug/l	99
84) 1,2,4-Trimethylbenzene	11.756	105	4136	0.386	ug/l	95
95) Naphthalene	13.780	128	4274	0.358	ug/l #	94

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

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