

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062022\
 Data File : VX029646.D
 Acq On : 21 Jun 2022 07:18
 Operator : JC/MD
 Sample : N3354-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-53

Quant Time: Jun 21 08:48:25 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	260863	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	445562	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	384120	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	167349	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	152620	44.198	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.400%
35) Dibromofluoromethane	5.391	113	131797	45.390	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	90.780%
50) Toluene-d8	8.653	98	499579	47.096	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.200%
62) 4-Bromofluorobenzene	11.085	95	180353	45.085	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.180%
Target Compounds						
					Qvalue	
8) Diethyl Ether	2.136	74	1332	0.774	ug/l	83
11) Tert butyl alcohol	2.971	59	339135	458.388	ug/l	99
16) Acetone	2.386	43	5082	3.692	ug/l	93
17) Carbon Disulfide	2.520	76	5063	0.739	ug/l	99
19) Methyl tert-butyl Ether	3.117	73	216460	23.592	ug/l	98
22) Diisopropyl ether	3.757	45	26755	3.170	ug/l	88
40) Benzene	6.037	78	279250	23.971	ug/l	97
52) Toluene	8.726	92	3775	0.506	ug/l	100

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

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