

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062022\
 Data File : VX029639.D
 Acq On : 21 Jun 2022 04:36
 Operator : JC/MD
 Sample : N3354-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-47

Quant Time: Jun 21 08:47:10 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	265965	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	454923	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	385325	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	169914	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	153439	43.583	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.160%
35) Dibromofluoromethane	5.391	113	136910	46.181	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.360%
50) Toluene-d8	8.653	98	510848	47.168	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.340%
62) 4-Bromofluorobenzene	11.079	95	184388	45.145	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.300%
Target Compounds						
					Qvalue	
8) Diethyl Ether	2.130	74	1301	0.742	ug/l	62
11) Tert butyl alcohol	2.977	59	49526	57.424	ug/l	99
16) Acetone	2.386	43	3969	2.828	ug/l	93
17) Carbon Disulfide	2.520	76	4812	0.689	ug/l	97
19) Methyl tert-butyl Ether	3.117	73	76553	8.184	ug/l	97
22) Diisopropyl ether	3.770	45	12839	1.492	ug/l	90
68) m/p-Xylenes	10.305	106	1844	0.348	ug/l	91
73) Isopropylbenzene	10.964	105	6355	0.499	ug/l	99

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

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