

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062022\
 Data File : VX029654.D
 Acq On : 21 Jun 2022 10:24
 Operator : JC/MD
 Sample : N3354-17
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-61

Quant Time: Jun 21 17:09:53 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	242168	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	411056	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	352987	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	155991	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	141403	44.111	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	88.220%
35) Dibromofluoromethane	5.391	113	124614	46.519	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.040%
50) Toluene-d8	8.653	98	461418	47.150	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.300%
62) 4-Bromofluorobenzene	11.085	95	167468	45.378	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.760%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.959	59	19077	18.749	ug/l	100
16) Acetone	2.380	43	16869	13.202	ug/l	99
17) Carbon Disulfide	2.514	76	1916	0.301	ug/l	99
19) Methyl tert-butyl Ether	3.117	73	30107	3.535	ug/l	98
40) Benzene	6.044	78	2457	0.229	ug/l	99
73) Isopropylbenzene	10.963	105	8082	0.691	ug/l	99
78) n-propylbenzene	11.305	91	2829	0.215	ug/l	98

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

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