

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061822\
 Data File : VX029575.D
 Acq On : 18 Jun 2022 17:23
 Operator : JC/MD
 Sample : N3325-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-24

Quant Time: Jun 20 01:35:11 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	273557	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	455228	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	395924	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	175066	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	156815	43.305	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	86.620%
35) Dibromofluoromethane	5.391	113	135879	45.802	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.600%
50) Toluene-d8	8.653	98	519028	47.891	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.780%
62) 4-Bromofluorobenzene	11.085	95	190252	46.550	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.100%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.971	59	60888	70.515	ug/l	99
16) Acetone	2.380	43	7291	5.051	ug/l #	84
17) Carbon Disulfide	2.514	76	3462	0.482	ug/l #	94
19) Methyl tert-butyl Ether	3.111	73	25966	2.699	ug/l #	1
22) Diisopropyl ether	3.763	45	11015	1.244	ug/l #	86
31) Cyclohexane	5.470	56	58469	12.027	ug/l	93
39) Methylcyclohexane	7.385	83	2837	0.567	ug/l	90
40) Benzene	6.044	78	3165	0.266	ug/l	91
52) Toluene	8.720	92	3086	0.405	ug/l	94
68) m/p-Xylenes	10.305	106	9619	1.765	ug/l	100
69) o-Xylene	10.646	106	2276	0.423	ug/l	91
73) Isopropylbenzene	10.963	105	11910	0.907	ug/l	100
78) n-propylbenzene	11.305	91	4702	0.318	ug/l	98
95) Naphthalene	13.780	128	4770	0.388	ug/l	98

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

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