

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010322\
 Data File : VX026225.D
 Acq On : 03 Jan 2022 17:10
 Operator : JC/MD
 Sample : M5335-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 20966

Quant Time: Jan 04 03:52:28 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	195805	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	345953	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	301097	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	135325	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	151055	53.325	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 106.660%			
35) Dibromofluoromethane	5.391	113	111948	49.773	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 99.540%			
50) Toluene-d8	8.653	98	401746	48.021	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 96.040%			
62) 4-Bromofluorobenzene	11.079	95	145380	48.815	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 97.620%			
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.953	59	12981	22.943	ug/l	98
16) Acetone	2.380	43	56811	37.015	ug/l	94
18) Methyl Acetate	2.703	43	3905	0.784	ug/l	100
25) 2-Butanone	4.568	43	14998	7.015	ug/l	98
40) Benzene	6.044	78	113695	10.981	ug/l	98
51) 4-Methyl-2-Pentanone	8.574	43	11302	2.649	ug/l	95
52) Toluene	8.720	92	185549	29.111	ug/l	98
67) Ethyl Benzene	10.195	91	22872	1.934	ug/l	99
68) m/p-Xylenes	10.299	106	36485	8.262	ug/l	94
69) o-Xylene	10.647	106	11896	2.753	ug/l	87
95) Naphthalene	13.780	128	2336	0.223	ug/l	98

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

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