

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010622\
 Data File : VX026383.D
 Acq On : 06 Jan 2022 18:58
 Operator : JC/MD
 Sample : N1020-16
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-28

Quant Time: Jan 07 04:18:07 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	5.562	168	152993	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	257381	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	219063	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	93484	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	108992	49.243	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.480%
35) Dibromofluoromethane	5.391	113	82085	49.055	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.120%
50) Toluene-d8	8.653	98	295630	47.498	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.000%
62) 4-Bromofluorobenzene	11.079	95	107586	48.556	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.120%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.965	59	16734	37.853	ug/l	100
16) Acetone	2.386	43	4140	3.452	ug/l	92
19) Methyl tert-butyl Ether	3.117	73	8248	1.394	ug/l	94
22) Diisopropyl ether	3.764	45	19270	3.016	ug/l #	70
37) Ethyl Acetate	4.721	43	11321	3.618	ug/l	99
40) Benzene	6.050	78	3947	0.512	ug/l	91
67) Ethyl Benzene	10.195	91	2711	0.315	ug/l	94
68) m/p-Xylenes	10.305	106	3683	1.146	ug/l	84

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

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