

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX040121\
 Data File : VX021717.D
 Acq On : 01 Apr 2021 18:37
 Operator : JC/MD
 Sample : M1895-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-8

Quant Time: Apr 02 03:31:06 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 24 12:25:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	56479	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	103430	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	107970	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.065	152	48184	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	49459	52.63	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	105.26%
35) Dibromofluoromethane	5.464	113	38439	52.32	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.64%
50) Toluene-d8	8.694	98	131942	50.40	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.80%
62) 4-Bromofluorobenzene	11.118	95	53031	52.10	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.20%
Target Compounds						Qvalue
16) Acetone	2.423	43	4973	14.00	ug/l	95
17) Carbon Disulfide	2.558	76	3217	2.01	ug/l	100
19) Methyl tert-butyl Ether	3.170	73	47309	22.26	ug/l	95
25) 2-Butanone	4.641	43	1126	2.09	ug/l #	87
29) Tetrahydrofuran	5.105	42	1077	3.08	ug/l #	73
40) Benzene	6.111	78	1122	0.39	ug/l	97
52) Toluene	8.765	92	807	0.45	ug/l	97
67) Ethyl Benzene	10.235	91	2033	0.51	ug/l	92
68) m/p-Xylenes	10.335	106	778	0.55	ug/l	85
69) o-Xylene	10.677	106	468	0.34	ug/l	90
70) Styrene	10.694	104	1069	0.46	ug/l	98

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

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