

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042021\
 Data File : VX021779.D
 Acq On : 21 Apr 2021 01:19
 Operator : JC/MD
 Sample : M2079-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-13

Quant Time: Apr 21 11:29:38 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042021W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 20 13:46:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	165554	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.830	114	277601	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.092	117	239977	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.061	152	100088	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.025	65	94040	50.07	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.14%	
35) Dibromofluoromethane	5.464	113	84647	47.68	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 95.36%	
50) Toluene-d8	8.696	98	329071	48.21	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 96.42%	
62) 4-Bromofluorobenzene	11.122	95	110337	43.28	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 86.56%	
Target Compounds						
						Qvalue
16) Acetone	2.416	43	28393	43.97	ug/l	97
25) 2-Butanone	4.635	43	8638	7.90	ug/l	98
30) Chloroform	5.172	83	3464	1.09	ug/l	98
40) Benzene	6.105	78	3513	0.48	ug/l	99
67) Ethyl Benzene	10.232	91	25603	2.99	ug/l	99
68) m/p-Xylenes	10.342	106	22345	6.73	ug/l	97
69) o-Xylene	10.683	106	12506	3.91	ug/l	97
73) Isopropylbenzene	11.000	105	3158	0.45	ug/l	95
78) n-propylbenzene	11.341	91	5974	0.77	ug/l	97
80) 1,3,5-Trimethylbenzene	11.488	105	13754	2.31	ug/l	100
84) 1,2,4-Trimethylbenzene	11.786	105	42516	7.08	ug/l	95
95) Naphthalene	13.817	128	11788	1.75	ug/l	99

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

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