

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061021\
 Data File : VX022653.D
 Acq On : 10 Jun 2021 19:16
 Operator : JC/MD
 Sample : M2645-03 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BUR-1041

Quant Time: Jun 11 06:47:34 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 12:20:00 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	58541	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	118322	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	105380	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	42081	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	49490	47.006	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	94.020%
35) Dibromofluoromethane	5.391	113	35551	47.092	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.180%
50) Toluene-d8	8.653	98	147001	49.396	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.800%
62) 4-Bromofluorobenzene	11.085	95	52475	46.623	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.240%
Target Compounds						
						Qvalue
13) Acrolein	2.239	56	8234	80.601	ug/l	98
16) Acetone	2.386	43	12139	25.749	ug/l	96
17) Carbon Disulfide	2.508	76	1070	0.669	ug/l	# 75
25) 2-Butanone	4.568	43	13204	17.270	ug/l	91
37) Ethyl Acetate	4.727	43	3770	2.516	ug/l	# 87
84) 1,2,4-Trimethylbenzene	11.756	105	734	0.251	ug/l	89
89) n-Butylbenzene	12.341	91	1364	0.489	ug/l	87
93) 1,2,4-Trichlorobenzene	13.597	180	492	0.601	ug/l	93
95) Naphthalene	13.780	128	1410	1.857	ug/l	# 94
96) 1,2,3-Trichlorobenzene	13.969	180	393	0.492	ug/l	93

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

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