

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092021\
 Data File : VX024352.D
 Acq On : 20 Sep 2021 15:42
 Operator : JC/MD
 Sample : M3817-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SW-CB-03-(1.0)

Manual Integrations
 APPROVED

MMDadoda
 9/21/2021 7:00:39 PM

Quant Time: Sep 21 08:41:39 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 14 12:18:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	124591	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	211904	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	233081	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94816	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	91187	51.326	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 102.660%			
35) Dibromofluoromethane	5.397	113	71302	50.173	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.340%			
50) Toluene-d8	8.653	98	308844	59.656	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 119.320%#			
62) 4-Bromofluorobenzene	11.085	95	107701	53.456	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 106.920%			
Target Compounds						Qvalue
8) Diethyl Ether	2.142	74	1716	2.375	ug/l	84
16) Acetone	2.386	43	3224	4.256	ug/l	97
40) Benzene	6.043	78	48245	8.634	ug/l	96
65) Chlorobenzene	10.085	112	40363	8.417	ug/l	97
87) 1,3-Dichlorobenzene	11.969	146	1556m	0.490	ug/l	
88) 1,4-Dichlorobenzene	12.048	146	3139m	0.967	ug/l	
91) 1,2-Dichlorobenzene	12.341	146	3497	1.121	ug/l	94

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

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