

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092021\
 Data File : VX024357.D
 Acq On : 20 Sep 2021 17:39
 Operator : JC/MD
 Sample : M3817-07RE
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

MMDadoda
 9/21/2021 7:01:00 PM

Quant Time: Sep 21 08:42:31 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	151903	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	265819	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	216666	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94088	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	111438	51.447	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 102.900%			
35) Dibromofluoromethane	5.397	113	88672	49.740	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 99.480%			
50) Toluene-d8	8.653	98	328075	50.517	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 101.040%			
62) 4-Bromofluorobenzene	11.085	95	99401	39.330	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 78.660%#			
Target Compounds						Qvalue
8) Diethyl Ether	2.142	74	1566	1.778	ug/l	72
16) Acetone	2.386	43	3387	3.667	ug/l #	83
40) Benzene	6.044	78	54637	7.795	ug/l	99
65) Chlorobenzene	10.086	112	32178	7.219	ug/l	99
87) 1,3-Dichlorobenzene	11.975	146	1262	0.400	ug/l	96
88) 1,4-Dichlorobenzene	12.042	146	2538m	0.788	ug/l	
91) 1,2-Dichlorobenzene	12.341	146	3033	0.980	ug/l	92

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

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