

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092121\
 Data File : VX024390.D
 Acq On : 21 Sep 2021 16:49
 Operator : JC/MD
 Sample : M3835-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

MMDadoda
 9/23/2021 11:06:46 AM

Quant Time: Sep 22 04:51:20 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 21 10:52:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	154336	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	242838	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	243554	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	95456	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	112222	50.992	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 101.980%			
35) Dibromofluoromethane	5.397	113	93831	57.615	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 115.240%			
50) Toluene-d8	8.653	98	316241	53.304	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 106.600%			
62) 4-Bromofluorobenzene	11.085	95	112973	48.930	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 97.860%			
Target Compounds					Qvalue	
8) Diethyl Ether	2.142	74	1618	1.808	ug/l	85
16) Acetone	2.392	43	2198	2.342	ug/l	92
40) Benzene	6.050	78	109114	17.040	ug/l	98
44) Trichloroethene	7.135	130	1096	0.609	ug/l	65
65) Chlorobenzene	10.085	112	74620	14.892	ug/l	99
87) 1,3-Dichlorobenzene	11.969	146	1555	0.486	ug/l	99
88) 1,4-Dichlorobenzene	12.042	146	3910m	1.197	ug/l	
91) 1,2-Dichlorobenzene	12.341	146	3863	1.230	ug/l	93

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

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