

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091721\
 Data File : VX024310.D
 Acq On : 17 Sep 2021 15:34
 Operator : JC/MD
 Sample : M3791-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SW-CB-04-(1.5)

Quant Time: Sep 17 16:06:27 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	141050	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	247261	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	199190	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	104553	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	105917	52.660	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 105.320%		
35) Dibromofluoromethane	5.391	113	83941	50.621	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.240%		
50) Toluene-d8	8.653	98	306374	50.717	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 101.440%		
62) 4-Bromofluorobenzene	11.085	95	102024	43.398	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 86.800%		
Target Compounds						
						Qvalue
16) Acetone	2.386	43	62738	73.154	ug/l #	88
25) 2-Butanone	4.574	43	7801	6.176	ug/l	92
40) Benzene	6.050	78	9289	1.425	ug/l	94
52) Toluene	8.720	92	12820	3.000	ug/l	99
65) Chlorobenzene	10.086	112	5861	1.430	ug/l	99
67) Ethyl Benzene	10.201	91	1480	0.206	ug/l	94
86) p-Isopropyltoluene	12.012	119	14667	2.202	ug/l	96
95) Naphthalene	13.780	128	57017	8.721	ug/l	99

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

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