

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092021\
 Data File : VX024355.D
 Acq On : 20 Sep 2021 16:52
 Operator : JC/MD
 Sample : M3817-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

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

Quant Time: Sep 21 08:42:09 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	144263	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	251719	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	239056	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	106898	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	104594	50.844	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.680%
35) Dibromofluoromethane	5.397	113	85720	50.778	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.560%
50) Toluene-d8	8.653	98	308136	50.105	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.200%
62) 4-Bromofluorobenzene	11.085	95	114540	47.859	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.720%
Target Compounds						
						Qvalue
16) Acetone	2.386	43	67498	76.951	ug/l	89
25) 2-Butanone	4.580	43	12544	9.709	ug/l #	87
40) Benzene	6.050	78	18728	2.821	ug/l	100
52) Toluene	8.720	92	83099	19.101	ug/l	97
65) Chlorobenzene	10.079	112	8799	1.789	ug/l	99
67) Ethyl Benzene	10.195	91	2324	0.270	ug/l	93
86) p-Isopropyltoluene	12.012	119	13112	1.925	ug/l	93
95) Naphthalene	13.780	128	34790	5.724	ug/l	98

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

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