

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062521\
 Data File : VX022920.D
 Acq On : 25 Jun 2021 14:05
 Operator : JC/MD
 Sample : M2838-19
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE104D2-20210622

Quant Time: Jun 26 03:53:23 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 23 19:12:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	52390	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	107538	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	100517	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	41448	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	50151	52.748	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	105.500%
35) Dibromofluoromethane	5.397	113	35898	52.568	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	105.140%
50) Toluene-d8	8.653	98	144878	54.132	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	108.260%
62) 4-Bromofluorobenzene	11.085	95	51316	51.375	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	102.760%
Target Compounds						
						Qvalue
12) 1,1-Dichloroethene	2.325	96	775	1.140	ug/l	# 63
16) Acetone	2.392	43	1354	3.212	ug/l	# 71
27) cis-1,2-Dichloroethene	4.501	96	27624	33.010	ug/l	97
30) Chloroform	5.111	83	2986	2.069	ug/l	98
44) Trichloroethene	7.135	130	162383	207.893	ug/l	92

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

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