

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
 Data File : VX021606.D
 Acq On : 30 Mar 2021 03:24
 Operator : JC/MD
 Sample : M1843-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE117D1-20210325

Quant Time: Mar 30 05:24:04 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 24 12:25:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	59109	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	109981	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	114852	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	52268	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	51488	52.35	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 104.70%		
35) Dibromofluoromethane	5.458	113	39593	50.68	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.36%		
50) Toluene-d8	8.694	98	141743	50.92	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 101.84%		
62) 4-Bromofluorobenzene	11.118	95	57922	53.52	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 107.04%		
Target Compounds						
						Qvalue
12) 1,1-Dichloroethene	2.352	96	450	0.76	ug/l	97
16) Acetone	2.423	43	2396	6.44	ug/l #	85
27) cis-1,2-Dichloroethene	4.558	96	737	0.97	ug/l	90
30) Chloroform	5.164	83	550	0.41	ug/l	100
44) Trichloroethene	7.188	130	82650	102.35	ug/l	99
55) 1,1,2-Trichloroethane	9.194	97	435	0.55	ug/l #	89

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

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