

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030821\
 Data File : VX021057.D
 Acq On : 08 Mar 2021 13:53
 Operator : JC/MD
 Sample : M1594-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWBR3-187-200-030521

Quant Time: Mar 09 01:05:26 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022421W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 24 13:50:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.635	168	208513	50.00	ug/l	0.01
34) 1,4-Difluorobenzene	6.829	114	393313	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	382738	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	178264	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	140328	49.82	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	99.64%		
35) Dibromofluoromethane	5.470	113	121116	50.68	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	101.36%		
50) Toluene-d8	8.694	98	479582	51.47	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	102.94%		
62) 4-Bromofluorobenzene	11.118	95	189716	55.06	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	110.12%		
Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.358	96	7372	3.16	ug/l	93
16) Acetone	2.417	43	3940	3.39	ug/l	99
17) Carbon Disulfide	2.558	76	6228	0.94	ug/l #	92
20) Methylene Chloride	2.840	84	3253	1.17	ug/l	94
24) 1,1-Dichloroethane	3.676	63	3664	0.81	ug/l	94
27) cis-1,2-Dichloroethene	4.564	96	14064	4.66	ug/l	88
30) Chloroform	5.182	83	12125	2.66	ug/l	98
42) 1,2-Dichloroethane	6.164	62	2718	0.73	ug/l	99
44) Trichloroethene	7.194	130	9144	3.10	ug/l	97
52) Toluene	8.765	92	41186	5.90	ug/l	99

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

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