

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX040121\
 Data File : VX021709.D
 Acq On : 01 Apr 2021 15:37
 Operator : JC/MD
 Sample : M1913-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWBR1-149-174-040121

Quant Time: Apr 02 03:28:28 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	5.635	168	54117	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.835	114	97553	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	101907	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	46257	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.035	65	46584	51.74	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	103.48%
35) Dibromofluoromethane	5.470	113	36370	52.49	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.98%
50) Toluene-d8	8.694	98	123766	50.12	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.24%
62) 4-Bromofluorobenzene	11.124	95	49388	51.44	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	102.88%

Target Compounds						Qvalue
16) Acetone	2.423	43	4574	13.43	ug/l #	84
20) Methylene Chloride	2.834	84	2061	3.10	ug/l #	87
25) 2-Butanone	4.646	43	1063	2.06	ug/l	94
27) cis-1,2-Dichloroethene	4.570	96	6229	8.95	ug/l	99
30) Chloroform	5.182	83	7549	6.17	ug/l	98
40) Benzene	6.111	78	1102	0.41	ug/l #	90
44) Trichloroethene	7.188	130	23711	33.10	ug/l	94
51) 4-Methyl-2-Pentanone	8.623	43	2448	2.22	ug/l	87
52) Toluene	8.765	92	190644	113.71	ug/l	99

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

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