

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030721\
 Data File : VX021028.D
 Acq On : 05 Mar 2021 12:57
 Operator : JC/MD
 Sample : M1574-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWBR3-150-160-030421

Quant Time: Mar 05 23:57:10 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	221668	50.00	ug/l	0.01
34) 1,4-Difluorobenzene	6.829	114	416801	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	413128	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	184561	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	146071	48.79	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	97.58%
35) Dibromofluoromethane	5.464	113	128397	50.70	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	101.40%
50) Toluene-d8	8.694	98	518946	52.56	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	105.12%
62) 4-Bromofluorobenzene	11.118	95	199783	54.72	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	109.44%
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.358	96	13944	5.62	ug/l	99
16) Acetone	2.417	43	5086	4.12	ug/l	100
17) Carbon Disulfide	2.552	76	13409	1.90	ug/l	99
20) Methylene Chloride	2.834	84	2150	0.73	ug/l	93
21) trans-1,2-Dichloroethene	3.146	96	769	0.28	ug/l #	77
24) 1,1-Dichloroethane	3.676	63	5149	1.07	ug/l	99
27) cis-1,2-Dichloroethene	4.564	96	19758	6.15	ug/l	89
30) Chloroform	5.170	83	29605	6.12	ug/l	99
42) 1,2-Dichloroethane	6.164	62	3364	0.85	ug/l	98
44) Trichloroethene	7.188	130	28454	9.11	ug/l	97
52) Toluene	8.765	92	107885	14.59	ug/l	99

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

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