

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030721\
 Data File : VX021029.D
 Acq On : 05 Mar 2021 13:18
 Operator : JC/MD
 Sample : M1574-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWBR3-160-170-030421

Quant Time: Mar 05 23:57:30 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	5.629	168	234029	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.835	114	440004	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	405583	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	172646	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.035	65	156770	49.59	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	99.18%
35) Dibromofluoromethane	5.464	113	137702	51.51	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	103.02%
50) Toluene-d8	8.694	98	535412	51.36	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	102.72%
62) 4-Bromofluorobenzene	11.118	95	194170	50.38	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	100.76%
Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.364	96	42024	16.04	ug/l	95
16) Acetone	2.423	43	2513	1.93	ug/l	99
17) Carbon Disulfide	2.558	76	13113	1.76	ug/l	100
20) Methylene Chloride	2.840	84	1889	0.60	ug/l	89
21) trans-1,2-Dichloroethene	3.146	96	1278	0.43	ug/l	95
24) 1,1-Dichloroethane	3.670	63	9921	1.96	ug/l	99
27) cis-1,2-Dichloroethene	4.570	96	45871	13.53	ug/l	90
30) Chloroform	5.176	83	18940	3.71	ug/l	93
42) 1,2-Dichloroethane	6.170	62	6346	1.52	ug/l	98
44) Trichloroethene	7.194	130	43332	13.15	ug/l	95
49) 1,4-Dioxane	7.718	88	1502	18.95	ug/l	99
52) Toluene	8.765	92	115500	14.79	ug/l	99

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

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