

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010622\
 Data File : VX026384.D
 Acq On : 06 Jan 2022 19:21
 Operator : JC/MD
 Sample : N1020-17
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-12

Quant Time: Jan 07 04:18:26 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	153425	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	257373	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	218012	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	92183	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	109916	49.520	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.040%
35) Dibromofluoromethane	5.391	113	82506	49.308	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.620%
50) Toluene-d8	8.653	98	294279	47.282	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	94.560%
62) 4-Bromofluorobenzene	11.079	95	107053	48.317	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.640%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	6625	5.509	ug/l	89
37) Ethyl Acetate	4.721	43	6167	1.971	ug/l #	92
40) Benzene	6.044	78	2766	0.359	ug/l	95
67) Ethyl Benzene	10.195	91	2030	0.237	ug/l	96
68) m/p-Xylenes	10.305	106	2538	0.794	ug/l	88

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010622\
Data File : VX026384.D
Acq On : 06 Jan 2022 19:21
Operator : JC/MD
Sample : N1020-17
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-12

Quant Time: Jan 07 04:18:26 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
Quant Title : SW846 8260
QLast Update : Thu Dec 30 08:02:21 2021
Response via : Initial Calibration

