

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081721\
 Data File : VX023772.D
 Acq On : 17 Aug 2021 18:51
 Operator : JC/MD
 Sample : M3389-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CRT-18

Quant Time: Aug 18 06:12:03 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 17 11:57:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	142304	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	246287	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	228340	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94258	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	100153	50.427	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 100.860%		
35) Dibromofluoromethane	5.397	113	78772	48.323	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 96.640%		
50) Toluene-d8	8.653	98	299437	48.263	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 96.520%		
62) 4-Bromofluorobenzene	11.085	95	102047	44.680	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 89.360%		
Target Compounds						
						Qvalue
16) Acetone	2.392	43	5061	5.699	ug/l #	78
20) Methylene Chloride	2.794	84	17047	10.050	ug/l	87
30) Chloroform	5.111	83	3094	1.047	ug/l #	82

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081721\
Data File : VX023772.D
Acq On : 17 Aug 2021 18:51
Operator : JC/MD
Sample : M3389-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
CRT-18

Quant Time: Aug 18 06:12:03 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260
QLast Update : Tue Aug 17 11:57:51 2021
Response via : Initial Calibration

