

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
 Data File : VX022476.D
 Acq On : 05 Jun 2021 16:55
 Operator : JC/MD
 Sample : M2589-23
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 72 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 18-B7(4-6)

Quant Time: Jun 07 03:26:05 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 19 14:23:10 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	61479	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	125986	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	111965	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	41103	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	52285	55.132	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	110.260%	
35) Dibromofluoromethane	5.397	113	38281	47.915	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.820%	
50) Toluene-d8	8.653	98	155964	50.677	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	101.360%	
62) 4-Bromofluorobenzene	11.085	95	53658	47.170	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	94.340%	
Target Compounds						
						Qvalue
16) Acetone	2.392	43	6742	16.327	ug/l	94
20) Methylene Chloride	2.794	84	4811	5.247	ug/l	88
43) Isopropyl Acetate	6.348	43	12356	5.290	ug/l	93

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022476.D
Acq On : 05 Jun 2021 16:55
Operator : JC/MD
Sample : M2589-23
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 72 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
18-B7(4-6)

Quant Time: Jun 07 03:26:05 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260
QLast Update : Wed May 19 14:23:10 2021
Response via : Initial Calibration

