

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072621\
 Data File : VX023449.D
 Acq On : 26 Jul 2021 14:27
 Operator : JC/MD
 Sample : M3145-20DL 5X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 152140-DDC-1-PDBDL

Quant Time: Jul 27 03:27:33 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 23 06:05:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	174414	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	294940	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	259374	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	101360	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	112799	51.846	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 103.700%	
35) Dibromofluoromethane	5.397	113	93363	50.628	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.260%	
50) Toluene-d8	8.653	98	349358	51.564	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 103.120%	
62) 4-Bromofluorobenzene	11.085	95	112114	48.663	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 97.320%	
Target Compounds						Qvalue
64) Tetrachloroethene	9.281	164	128248	62.127	ug/l	99
90) Hexachloroethane	12.543	117	1014	2.750	ug/l	86

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072621\
Data File : VX023449.D
Acq On : 26 Jul 2021 14:27
Operator : JC/MD
Sample : M3145-20DL 5X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
152140-DDC-1-PDBDL

Quant Time: Jul 27 03:27:33 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
Quant Title : SW846 8260
QLast Update : Fri Jul 23 06:05:38 2021
Response via : Initial Calibration

