

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061818\
 Data File : VX002707.D
 Acq On : 19 Jun 2018 04:33
 Operator : JC/MD
 Sample : J3537-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS037MW208-180613

Quant Time: Jun 19 07:11:03 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 16 01:37:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	221791	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	315455	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	292731	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.09	152	166563	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	165807	49.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.40%	
35) Dibromofluoromethane	5.51	113	130426	51.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.60%	
50) Toluene-d8	8.72	98	491953	51.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.92%	
62) 4-Bromofluorobenzene	11.14	95	171778	47.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.84%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.60	96	4160	1.521	ug/l	56
30) Chloroform	5.23	83	1066	0.229	ug/l #	82
44) Trichloroethene	7.21	130	3067	1.058	ug/l	90
64) Tetrachloroethene	9.34	164	10328	3.178	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061818\
Data File : VX002707.D
Acq On : 19 Jun 2018 04:33
Operator : JC/MD
Sample : J3537-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SS037MW208-180613

Quant Time: Jun 19 07:11:03 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061518W.M
Quant Title : SW846 8260
QLast Update : Sat Jun 16 01:37:33 2018
Response via : Initial Calibration

