

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX111218\
 Data File : VX005913.D
 Acq On : 12 Nov 2018 20:54
 Operator : JC/MD
 Sample : J5912-09
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS003EW296-181109

Quant Time: Nov 13 06:22:25 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 07 14:40:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	108110	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	157747	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	142483	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	70830	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	65692	52.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.72%	
35) Dibromofluoromethane	5.50	113	53239	49.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.24%	
50) Toluene-d8	8.71	98	190403	53.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.76%	
62) 4-Bromofluorobenzene	11.14	95	64727	50.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.94%	
Target Compounds						
65) Chlorobenzene	10.14	112	35278	12.359	ug/l	99
87) 1,3-Dichlorobenzene	12.02	146	14434	6.515	ug/l	99
88) 1,4-Dichlorobenzene	12.10	146	39294	17.324	ug/l	98
91) 1,2-Dichlorobenzene	12.39	146	201185	90.640	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX111218\
Data File : VX005913.D
Acq On : 12 Nov 2018 20:54
Operator : JC/MD
Sample : J5912-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SS003EW296-181109

Quant Time: Nov 13 06:22:25 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 07 14:40:53 2018
Response via : Initial Calibration

