

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080918\
 Data File : VX003956.D
 Acq On : 09 Aug 2018 17:43
 Operator : JC/MD
 Sample : J4379-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-105I

Quant Time: Aug 10 16:04:10 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X080618W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 07 07:45:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	323825	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	505695	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	479464	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	248101	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	220286	51.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.64%	
35) Dibromofluoromethane	5.51	113	189852	49.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.12%	
50) Toluene-d8	8.72	98	713239	47.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.74%	
62) 4-Bromofluorobenzene	11.14	95	255232	52.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.90%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.60	96	3020	0.71	ug/l	70
44) Trichloroethene	7.22	130	2508	0.54	ug/l	95
64) Tetrachloroethene	9.34	164	1695	0.36	ug/l	92
68) m/p-Xylenes	10.36	106	3458	0.43	ug/l	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX080918\
Data File : VX003956.D
Acq On : 09 Aug 2018 17:43
Operator : JC/MD
Sample : J4379-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-105I

Quant Time: Aug 10 16:04:10 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X080618W.M
Quant Title : SW846 8260
QLast Update : Tue Aug 07 07:45:48 2018
Response via : Initial Calibration

