

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081018\
 Data File : VX003975.D
 Acq On : 10 Aug 2018 14:30
 Operator : JC/MD
 Sample : J4379-20
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-102I

Quant Time: Aug 13 10:47:34 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	320361	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	499969	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	465007	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	255033	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	221695	52.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.40%	
35) Dibromofluoromethane	5.51	113	192562	50.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.68%	
50) Toluene-d8	8.72	98	702346	47.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.36%	
62) 4-Bromofluorobenzene	11.14	95	242591	50.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.80%	
Target Compounds						
16) Acetone	2.45	43	15815	6.26	ug/l	97
27) cis-1,2-Dichloroethene	4.60	96	3284	0.78	ug/l	78
30) Chloroform	5.22	83	1837	0.27	ug/l	85
44) Trichloroethene	7.21	130	10899	2.37	ug/l	99
64) Tetrachloroethene	9.34	164	20784	4.54	ug/l #	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081018\
Data File : VX003975.D
Acq On : 10 Aug 2018 14:30
Operator : JC/MD
Sample : J4379-20
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-102I

Quant Time: Aug 13 10:47:34 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

