

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX081418\
 Data File : VX004056.D
 Acq On : 15 Aug 2018 04:21
 Operator : JC/MD
 Sample : J4439-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-289-SRI3

Quant Time: Aug 15 15:54:53 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X081418W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 14 13:27:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	280627	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	444412	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	417008	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	228309	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	207987	52.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.20%	
35) Dibromofluoromethane	5.51	113	172103	51.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.48%	
50) Toluene-d8	8.72	98	632374	46.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.78%	
62) 4-Bromofluorobenzene	11.14	95	215241	44.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.02%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.39	101	60612	19.10	ug/l	Qvalue 100
30) Chloroform	5.22	83	7259	1.19	ug/l	100
44) Trichloroethene	7.21	130	5403	1.31	ug/l	87
64) Tetrachloroethene	9.34	164	9130	2.31	ug/l	93

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX081418\
Data File : VX004056.D
Acq On : 15 Aug 2018 04:21
Operator : JC/MD
Sample : J4439-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-289-SRI3

Quant Time: Aug 15 15:54:53 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X081418W.M
Quant Title : SW846 8260
QLast Update : Tue Aug 14 13:27:25 2018
Response via : Initial Calibration

