

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031919\
 Data File : VX008195.D
 Acq On : 19 Mar 2019 17:15
 Operator : JC/SP
 Sample : K1860-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE105D1-20190311

Quant Time: Mar 20 07:59:02 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031919W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 20 06:51:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	350863	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	564405	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	479630	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	189089	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	240139	45.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.88%	
35) Dibromofluoromethane	5.50	113	176504	46.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.42%	
50) Toluene-d8	8.71	98	700975	48.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.86%	
62) 4-Bromofluorobenzene	11.13	95	224136	45.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.88%	
Target Compounds						
9) 1,1,2-Trichlorotrifluoroet	2.39	101	13524	2.999	ug/l	89
16) Acetone	2.45	43	6661	2.223	ug/l	99
27) cis-1,2-Dichloroethene	4.60	96	4609	0.879	ug/l	89
44) Trichloroethene	7.21	130	401835	84.498	ug/l	89
49) 1,4-Dioxane	7.74	88	1686	14.432	ug/l #	91
64) Tetrachloroethene	9.33	164	2491	0.576	ug/l #	84

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX031919\
Data File : VX008195.D
Acq On : 19 Mar 2019 17:15
Operator : JC/SP
Sample : K1860-07
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RE105D1-20190311

Quant Time: Mar 20 07:59:02 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X031919W.M
Quant Title : SW846 8260
QLast Update : Wed Mar 20 06:51:58 2019
Response via : Initial Calibration

