

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062419\
 Data File : VX010415.D
 Acq On : 24 Jun 2019 16:37
 Operator : JC/SP
 Sample : K3164-14
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BAT-B02

Quant Time: Jun 25 04:22:00 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 20 03:31:56 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	288108	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	440021	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	399555	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	185097	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	135174	50.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.08%	
35) Dibromofluoromethane	5.50	113	133969	49.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.52%	
50) Toluene-d8	8.71	98	516953	50.19	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.38%	
62) 4-Bromofluorobenzene	11.13	95	175941	47.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.62%	
Target Compounds						
3) Chloromethane	1.33	50	6893	2.955	ug/l	95
16) Acetone	2.45	43	6647	4.910	ug/l	88
44) Trichloroethene	7.22	130	2123	0.651	ug/l	83
64) Tetrachloroethene	9.34	164	24351	7.535	ug/l	92

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062419\
Data File : VX010415.D
Acq On : 24 Jun 2019 16:37
Operator : JC/SP
Sample : K3164-14
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BAT-B02

Quant Time: Jun 25 04:22:00 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260
QLast Update : Thu Jun 20 03:31:56 2019
Response via : Initial Calibration

