

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011719\
 Data File : VX007098.D
 Acq On : 17 Jan 2019 19:23
 Operator : JC/SP
 Sample : K1157-18
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-02-011519

Quant Time: Jan 18 06:55:53 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 08 14:06:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	572855	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	879970	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	847258	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	419587	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	304976	50.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.88%	
35) Dibromofluoromethane	5.50	113	268106	47.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.16%	
50) Toluene-d8	8.71	98	1066217	51.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.38%	
62) 4-Bromofluorobenzene	11.13	95	383193	53.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.58%	
Target Compounds						
16) Acetone	2.45	43	72136	24.373	ug/l	98
18) Methyl Acetate	2.79	43	14780	1.724	ug/l	93
20) Methylene Chloride	2.86	84	11878	1.562	ug/l	90
43) Isopropyl Acetate	6.46	43	13811	1.067	ug/l #	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX011719\
Data File : VX007098.D
Acq On : 17 Jan 2019 19:23
Operator : JC/SP
Sample : K1157-18
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-02-011519

Quant Time: Jan 18 06:55:53 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
Quant Title : SW846 8260
QLast Update : Tue Jan 08 14:06:59 2019
Response via : Initial Calibration

