

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011719\
 Data File : VX007118.D
 Acq On : 18 Jan 2019 05:08
 Operator : JC/SP
 Sample : K1157-22
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-04-011519

Quant Time: Jan 18 07:38:54 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	577519	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	884990	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	860744	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	423819	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	309962	51.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.72%	
35) Dibromofluoromethane	5.51	113	272497	48.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.16%	
50) Toluene-d8	8.71	98	1082375	52.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.36%	
62) 4-Bromofluorobenzene	11.13	95	394775	54.59	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.18%	
Target Compounds						
16) Acetone	2.45	43	31522	10.564	ug/l	97
18) Methyl Acetate	2.78	43	21346	2.470	ug/l	98
20) Methylene Chloride	2.86	84	38466	6.283	ug/l	98
43) Isopropyl Acetate	6.46	43	20594	1.583	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX011719\
Data File : VX007118.D
Acq On : 18 Jan 2019 05:08
Operator : JC/SP
Sample : K1157-22
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
WC-04-011519

Quant Time: Jan 18 07:38:54 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

