

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091520\
 Data File : VX018419.D
 Acq On : 15 Sep 2020 19:55
 Operator : JC/SP
 Sample : L3871-10
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TR-04-091420

Quant Time: Sep 16 04:28:38 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082120W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 21 03:03:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	439671	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	702861	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	704232	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	336474	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	319994	52.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.86%	
35) Dibromofluoromethane	5.47	113	249883	52.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.40%	
50) Toluene-d8	8.70	98	898179	50.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.76%	
62) 4-Bromofluorobenzene	11.12	95	332485	47.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.80%	
Target Compounds						
16) Acetone	2.42	43	24922	10.478	ug/l	98
18) Methyl Acetate	2.75	43	8181	1.425	ug/l #	87
20) Methylene Chloride	2.84	84	13589	1.665	ug/l #	81
43) Isopropyl Acetate	6.42	43	128534	10.569	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX091520\
Data File : VX018419.D
Acq On : 15 Sep 2020 19:55
Operator : JC/SP
Sample : L3871-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TR-04-091420

Quant Time: Sep 16 04:28:38 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X082120W.M
Quant Title : SW846 8260
QLast Update : Fri Aug 21 03:03:56 2020
Response via : Initial Calibration

