

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
 Data File : VX017601.D
 Acq On : 25 Jul 2020 17:09
 Operator : JC/SP
 Sample : L3383-19
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 75 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CRT-008

Quant Time: Jul 27 04:02:31 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 24 05:48:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	442152	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	697970	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	820934	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	422109	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	298876	52.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.24%	
35) Dibromofluoromethane	5.46	113	256748	55.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.00%	
50) Toluene-d8	8.70	98	963571	57.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.44%	
62) 4-Bromofluorobenzene	11.12	95	415257	61.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	123.24%	
Target Compounds						
16) Acetone	2.42	43	29196	11.679	ug/l	98
18) Methyl Acetate	2.75	43	7204	1.224	ug/l #	78
20) Methylene Chloride	2.84	84	37492	6.855	ug/l	94
43) Isopropyl Acetate	6.42	43	142513	11.793	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX072420\
Data File : VX017601.D
Acq On : 25 Jul 2020 17:09
Operator : JC/SP
Sample : L3383-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 75 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
CRT-008

Quant Time: Jul 27 04:02:31 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
Quant Title : SW846 8260
QLast Update : Fri Jul 24 05:48:09 2020
Response via : Initial Calibration

