

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX072320\
 Data File : VX017502.D
 Acq On : 23 Jul 2020 19:01
 Operator : JC/SP
 Sample : L3187-13
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SU-04-072120

Quant Time: Jul 24 06:44:11 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	5.63	168	470652	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	770116	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	846699	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	428079	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	320314	52.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.96%	
35) Dibromofluoromethane	5.47	113	269231	52.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.54%	
50) Toluene-d8	8.70	98	927523	49.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.84%	
62) 4-Bromofluorobenzene	11.12	95	433334	58.28	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.56%	
Target Compounds						
16) Acetone	2.42	43	79270	29.790	ug/l	99
18) Methyl Acetate	2.75	43	6385	1.019	ug/l #	87
20) Methylene Chloride	2.84	84	46179	8.139	ug/l	97
43) Isopropyl Acetate	6.42	43	152411	11.430	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX072320\
Data File : VX017502.D
Acq On : 23 Jul 2020 19:01
Operator : JC/SP
Sample : L3187-13
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SU-04-072120

Quant Time: Jul 24 06:44:11 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

