

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN073120\
 Data File : VN062697.D
 Acq On : 31 Jul 2020 19:56
 Operator : JC/MD
 Sample : L3511-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1

Quant Time: Aug 01 07:21:26 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 21 18:48:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	124627	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	237295	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	235903	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	86253	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	111154	57.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	114.82%	
35) Dibromofluoromethane	7.55	113	76319	48.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.92%	
50) Toluene-d8	10.06	98	308921	49.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.34%	
62) 4-Bromofluorobenzene	12.38	95	112465	49.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.92%	
Target Compounds						
16) Acetone	3.78	43	14585	20.020	ug/l	97
18) Methyl Acetate	4.28	43	6922	3.753	ug/l	95
20) Methylene Chloride	4.50	84	8280	4.739	ug/l	93
43) Isopropyl Acetate	8.13	43	47471	10.801	ug/l #	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN073120\
Data File : VN062697.D
Acq On : 31 Jul 2020 19:56
Operator : JC/MD
Sample : L3511-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-1

Quant Time: Aug 01 07:21:26 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N072120W.M
Quant Title : SW846 8260
QLast Update : Tue Jul 21 18:48:59 2020
Response via : Initial Calibration

