

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090120\
 Data File : VN063167.D
 Acq On : 1 Sep 2020 20:05
 Operator : JC/MD
 Sample : L3851-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 11995

Quant Time: Sep 02 05:01:42 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N083120W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 01 04:21:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	354337	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	709678	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	711170	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	288107	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	288205	49.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.52%	
35) Dibromofluoromethane	7.55	113	221045	48.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.10%	
50) Toluene-d8	10.06	98	897797	51.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.46%	
62) 4-Bromofluorobenzene	12.38	95	327030	53.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.70%	
Target Compounds						
8) Diethyl Ether	3.38	74	3112	1.159	ug/l	56
16) Acetone	3.78	43	39553	24.021	ug/l	98
18) Methyl Acetate	4.30	43	9650	2.286	ug/l #	55
20) Methylene Chloride	4.51	84	18986	3.645	ug/l	92
43) Isopropyl Acetate	8.13	43	128021	11.506	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN090120\
Data File : VN063167.D
Acq On : 1 Sep 2020 20:05
Operator : JC/MD
Sample : L3851-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
11995

Quant Time: Sep 02 05:01:42 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N083120W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 01 04:21:28 2020
Response via : Initial Calibration

