

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082020\
 Data File : VN063066.D
 Acq On : 21 Aug 2020 4:30
 Operator : JC/MD
 Sample : L3502-15
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OK-02-081920

Quant Time: Aug 21 07:50:15 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 19 06:09:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	138756	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	237478	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	264069	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	90924	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	96899	58.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	117.84%	
35) Dibromofluoromethane	7.55	113	84400	54.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.08%	
50) Toluene-d8	10.06	98	324212	54.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.10%	
62) 4-Bromofluorobenzene	12.38	95	106281	50.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.34%	
Target Compounds						
16) Acetone	3.78	43	8612	13.076	ug/l	95
18) Methyl Acetate	4.30	43	2457	1.503	ug/l #	66
20) Methylene Chloride	4.50	84	13690	8.329	ug/l	98
43) Isopropyl Acetate	8.13	43	33470	10.303	ug/l #	89

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN082020\
Data File : VN063066.D
Acq On : 21 Aug 2020 4:30
Operator : JC/MD
Sample : L3502-15
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
OK-02-081920

Quant Time: Aug 21 07:50:15 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N081820W.M
Quant Title : SW846 8260
QLast Update : Wed Aug 19 06:09:27 2020
Response via : Initial Calibration

