

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082420\
 Data File : VN063096.D
 Acq On : 24 Aug 2020 15:28
 Operator : JC/MD
 Sample : L3780-23
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 JC-03-082120-C

Quant Time: Aug 25 01:29:38 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	7.62	168	146240	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	243996	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	250212	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	93300	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	96160	55.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.96%	
35) Dibromofluoromethane	7.54	113	81606	51.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.66%	
50) Toluene-d8	10.06	98	314894	51.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.12%	
62) 4-Bromofluorobenzene	12.38	95	103950	48.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.48%	
Target Compounds						
16) Acetone	3.80	43	10681	16.092	ug/l	89
18) Methyl Acetate	4.28	43	4991	2.898	ug/l #	72
20) Methylene Chloride	4.47	84	10717	5.833	ug/l	94
43) Isopropyl Acetate	8.13	43	32976	9.879	ug/l #	87

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN082420\
Data File : VN063096.D
Acq On : 24 Aug 2020 15:28
Operator : JC/MD
Sample : L3780-23
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
JC-03-082120-C

Quant Time: Aug 25 01:29:38 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

