

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090420\
 Data File : VN063253.D
 Acq On : 4 Sep 2020 17:25
 Operator : JC/MD
 Sample : L3864-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CL-01-090320

Manual Integrations
 APPROVED

MMDadoda
 9/8/2020 1:33:28 PM

Quant Time: Sep 07 04:06:12 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	289243	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	571714	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	586417	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	224887	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	225979	47.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.58%	
35) Dibromofluoromethane	7.55	113	180616	49.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.48%	
50) Toluene-d8	10.06	98	740407	52.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.90%	
62) 4-Bromofluorobenzene	12.38	95	253349	51.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.60%	
Target Compounds						
16) Acetone	3.78	43	42410	31.552	ug/l	98
20) Methylene Chloride	4.51	84	14142m	3.282	ug/l	
43) Isopropyl Acetate	8.13	43	82754	9.232	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN090420\
Data File : VN063253.D
Acq On : 4 Sep 2020 17:25
Operator : JC/MD
Sample : L3864-07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
CL-01-090320

Manual Integrations
APPROVED

MMDadoda
9/8/2020 1:33:28 PM

Quant Time: Sep 07 04:06:12 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

