



Filename = JXS20_PROTON_FT-1-1.jdf
Author = chemteaching2
Experiment = proton.jxp
Sample_Id = JXS20
Solvent = CHLOROFORM-D
Actual_Start_Time = 4-FEB-2022 17:54:52
Revision_Time = 4-FEB-2022 17:56:02

Data_Format = 1D COMPLEX
Dim_Size = 26214
Dim_Title = Proton
Dim_Units = [ppm]
Dimensions = X
Site = CFNMR JEOL400
Spectrometer = JNM-ECZ400S/L1

Field_Strength = 9.389966[T] (400[MHz])
X_Acq_Duration = 4.37256192[s]
X_Domain = Proton
X_Freq = 399.79071365[MHz]
X_Offset = 5[ppm]
X_Points = 32768
X_Prescans = 0
X_Resolution = 0.22869888[Hz]
X_Sweep = 7.4940048[kHz]
X_Sweep_Clipped = 5.99520384[kHz]
Irr_Domain = Proton
Irr_Freq = 399.79071365[MHz]
Irr_Offset = 5[ppm]
Tri_Domain = Proton
Tri_Freq = 399.79071365[MHz]
Tri_Offset = 5[ppm]
Blanking = 2[us]
Clipped = FALSE
Scans = 8
Total_Scans = 8

Relaxation_Delay = 4[s]
Recvr_Gain = 32
Temp_Get = 25[dC]
X_90_Width = 6.39[us]
X_Acq_Time = 4.37256192[s]
X_Angle = 45[deg]
X_Atn = 5[dB]
X_Pulse = 3.195[us]
Irr_Mode = Off
Tri_Mode = Off
Dante_Loop = 400
Dante_Presat = FALSE
Decimation_Rate = 0
Experiment_Path = c:\Program Files\JEOL\C
Initial_Wait = 1[s]
Phase = {0, 90, 270, 180, 180,
Presat_Time = 4[s]
Presat_Time_Flag = FALSE
Relaxation_Delay_Calc = 0[s]
Relaxation_Delay_Temp = 4[s]
Repetition_Time = 8.37256192[s]