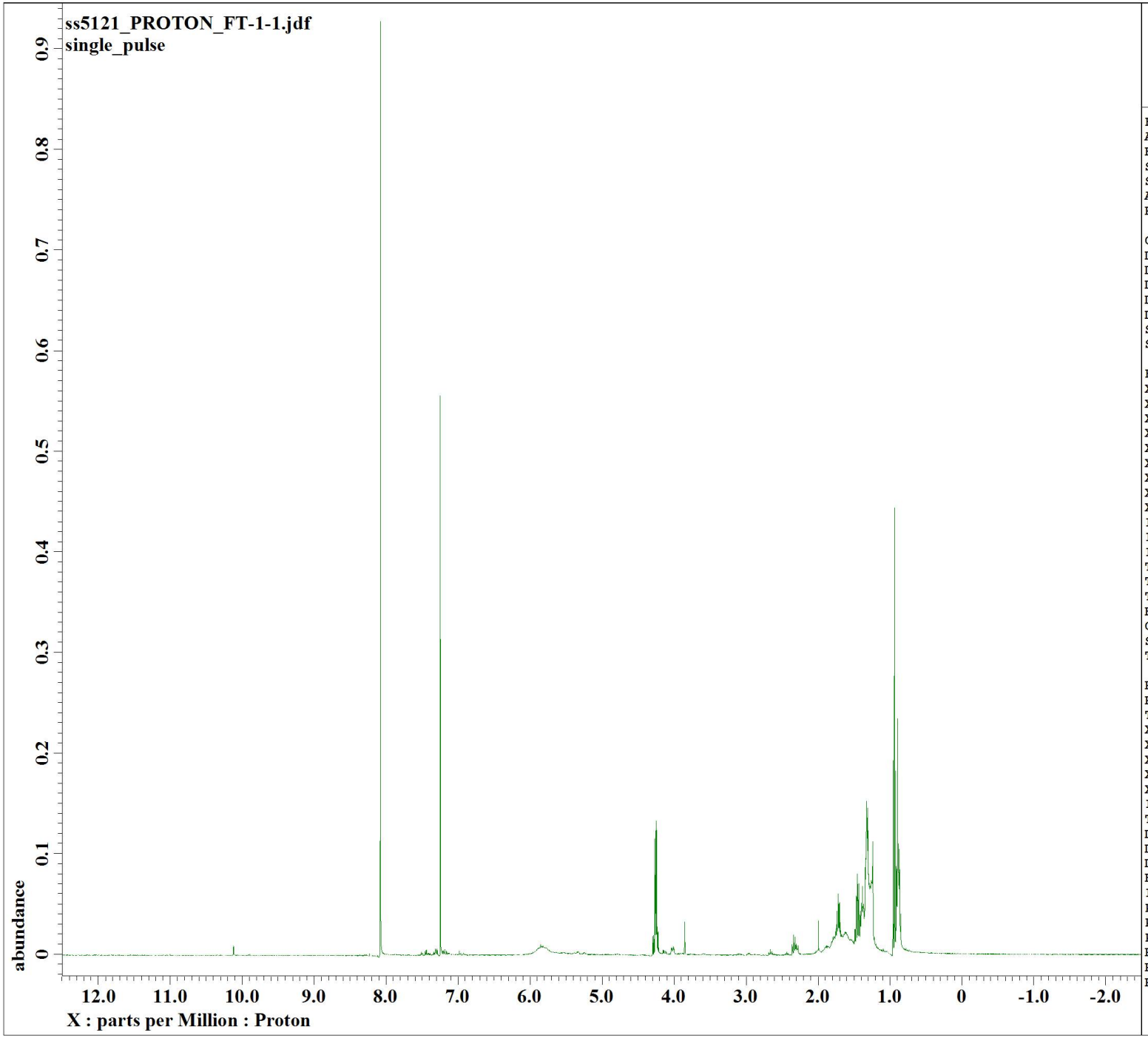


```
Filename                = ss5121_PROTON_FT-1-1.jd
Author                  = chemteaching2
Experiment              = proton.jxp
Sample_Id               = ss5121
Solvent                 = CHLOROFORM-D
Actual_Start_Time       =  4-MAR-2022 20:39:16
Revision_Time           =  4-MAR-2022 20:40:28

Comment                 = single_pulse
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              = 52
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]
```