



Filename = 171003_3_13C-1-4.jdf
Author = delta
Experiment = carbon.jxp
Sample_Id = 171003_3
Solvent = METHANOL-D4
Creation_Time = 4-OCT-2017 01:10:16
Revision_Time = 4-OCT-2017 09:43:07
Current_Time = 4-OCT-2017 09:43:43

Comment = 3/13C
Data_Format = 1D COMPLEX
Dim_Size = 26214
Dim_Title = Carbon13
Dim_Units = [ppm]
Dimensions = X
Spectrometer = DELTA2_NMR

Field_Strength = 14.09636928[T] (600[MHz])
X_Acq_Duration = 0.69206016[s]
X_Domain = 13C
X_Freq = 150.91343039[MHz]
X_Offset = 100[ppm]
X_Points = 32768
X_Prescans = 4
X_Resolution = 1.44496109[Hz]
X_Sweep = 47.34848485[kHz]
X_Sweep_Clipped = 37.87878788[kHz]
Irr_Domain = Proton
Irr_Freq = 600.1723046[MHz]
Irr_Offset = 5[ppm]
Clipped = FALSE
Scans = 4000
Total_Scans = 4000

Relaxation_Delay = 2[s]
Recvr_Gain = 60
Temp_Get = 24.7[dC]
X_90_Width = 10.2[us]
X_Acq_Time = 0.69206016[s]
X_Angle = 30[deg]
X_Atn = 11.1[dB]
X_Pulse = 3.4[us]
Irr_Atn_Dec = 24.81[dB]
Irr_Atn_Noe = 24.81[dB]
Irr_Noise = WALTZ
Irr_Pwidth = 76[us]
Decoupling = TRUE
Initial_Wait = 1[s]
Noe = TRUE
Noe_Time = 2[s]
Repetition_Time = 2.69206016[s]