week1

August 23, 2023

Notebook file for Week 1

Import libraries. The nmrbase folder needs to be located within the folder containing this Jupyter notebook

```
[]: import numpy as np
import matplotlib.pyplot as plt
import nmrbase.expbase as expbase
import nmrbase.expdta as expdta
```

Tuning response experiment

```
[]: filename = r"../DIRECTORY/FILENAME" #Defines the path to the data file. Here \rightarrow the path is relative to the current folder.
```

```
[]: a = expbase.expbase()  # create an experiment base object (a)
a.load(filename)  # load data into a

f1 = plt.figure()
ax1 = f1.subplots()  # create axes for a figure

a.plottm(ax1,1)  # plot the time domain data of the

→ first scan from the data set
```

```
[]: ## TASKS:

## change parameter "ftmin" and "ftmax" to select the time interval

corresponding to the ringdown for the Fourier transform

## change parameter "ffmin" and "ffmax" to select the frequency range of the

cutuing response signal

## include statements to label axes

## use statements to change appearance such as font size, etc.
```

```
## Adjust horizontal axis limits to display the spectrum of the ring-down signal
     f2=plt.figure()
     ax2=f2.subplots()
     a.pproc['ftmin']=VALUE
                                                    # time interval for Fourier_
      \rightarrow transform (0 = all data)
     a.pproc['ftmax']=VALUE
     a.pproc['ffmin']=VALUE
                                                    # frequency interval for spectrum_
      \rightarrow display (0 = all data)
     a.pproc['ffmax']=VALUE
     a.proc()
                                            # calculate Fourier transform of the data
      \rightarrow in a
     a.plotfrq(ax2,1)
                                            # plot the frequency domain data of the
      \rightarrow first scan from the data set
                                     # find data point index of largest peak in spectrum
[]: i=np.argmax(a.frq[0].dta)
     f=a.frq[0].ind_to_x(i)
                                     # find frequency corresponding to largest peak in_
     \hookrightarrow spectrum
     print("Tuning frequency: {:.2f} Hz".format(f)) # this statement prints the
      →identified frequency in a nice format
[]: # plot figure again with line indicating maximum peak position. Notice this,
      →works only if previous parameters have been correctly selected.
     ax2.axvline(x = f, color = 'r', label = 'Larmor Frequency')
     f2
          # show the figure again with the line included
    Pulse length calibration. NOTE: To analyze and report multiple datasets, simply copy the necessary
    blocks of code and change the name of the veriables.
[]: filename2 = r"../DIRECTORY/FILENAME"
                                                     # path of data file
[]: a2 = expbase.expbase()
     a2.load(filename2)
                                                 # load data
     f3=plt.figure()
     ax3=f3.subplots()
     a2.plottm(ax3,5)
[ ]: ## TASKS:
     ## fine tune with ax3.set_xlim and ax3.set_ylim parameters to zoom in on the echo
     ax3.set_ylim([VALUE, VALUE])
```

ax3.set_xlim([VALUE, VALUE])

```
f3
          # display adapted figrue
                                   # set appropriate digital filter parameters
[]: a2.pproc['digfmin']=1500
     a2.pproc['digfmax']=3500
     a2.digfilt()
                                    # perform digital filter operation
[]: f4=plt.figure()
     ax4=f4.subplots()
     a2.plottm(ax4,5)
                        # plot the 5th scan (the 1st scan doesn't have signal because,
      → the pulse is too short)
[ ]: ## TASKS:
     ## fine tune with ax4.set_xlim and ax4.set_ylim parameters to zoom in on the echo
     ax4.set_ylim([VALUE, VALUE])
     ax4.set_xlim([VALUE, VALUE])
          # display adapted figrue
[]: f5=plt.figure()
     ax5 = f5.subplots()
     a2.pproc['ftmin']=VALUE
                                                    # time interval for Fourier_
      \rightarrow transform (0 = all data)
     a2.pproc['ftmax']=VALUE
     a2.pproc['ffmin']=VALUE
                                                    # frequency interval for spectrum_
      \rightarrow display (0 = all data)
     a2.pproc['ffmax']=VALUE
     a2.proc()
                                             # calculate Fourier transform of the data
      \rightarrow in a
     ## this will use the digitally filtered data from before. Instead, the original \Box
      ⇒ data can be processed by loading it again.
     a2.plotfrq(ax5,5)
                                             # plot the frequency domain data of the
      \hookrightarrow first scan from the data set
     ## TASKS:
     ## change parameter "ftmin" and "ftmax" to select the time interval_{\sqcup}
     →corresponding to the ringdown for the Fourier transform
     ## change parameter "ffmin" and "ffmax" to select the frequency range of the
     → tuning response signal
     ## include statements to label axes
     ## use statements to change appearance such as font size, etc.
     ## Adjust horizontal axis limits to display the spectrum of the ring-down signal
```

```
[]: # This section is to integrate the peaks in the NMR spectra to determine pulse.
      \rightarrow length
     ## TASKS:
     ## change the intmin and intmax parameters to select the correct frequency range_
     \rightarrow for integration
     ## use set_xlabel and set_ylabel to set the labels to qet a publication quality.
     \hookrightarrow figure
     ## adjust appearance of figure as needed
     ## change and report the "x" parameter to indicate the pulse length for 90° ___
     →pulse with a vertical line
     f5=plt.figure()
     ax5=f5.subplots()
                                  # set correct frequency range for integration
     a2.pproc['intmin']=VALUE
     a2.pproc['intmax']=VALUE
     a2.integrate()
                                  # perform integration
     #find the starting pulselength and increment, then set the correct x-axis
     dx=a2.pinc["inc"][0]
     x0=a2.p["p1"]
     print('x0 = ',x0,'s , dx = ',dx,'s') # x0 is the starting pulse length,
     \rightarrow and dx is the increment
     a2.idt.x0=x0
     a2.idt.dx=dx
     a2.idt.plot(ax5,disp=[0])
                                                 # disp=0 plots only the NMR signal
     \rightarrow in trace 0
     ax5.axvline(x = VALUE, color = 'r', label = '90° pulse length') # display
      →vertical line at 90 degree pulse length
```