

# **Students' Annual Webinar**

## **Experimental time saving for assignment and restraints at fast MAS**

**Sahil R**

The chemical shift assignment is the first step for structure calculations by NMR, followed by obtaining sufficient distance restraints. The assignments and restraints usually require 3-6 different sets of spectra which makes the processing time intensive. In the talk, I will present the pulse programs to acquire multiple experiments in a single experiment (multiple acquisitions). The pulse programs for backbone, side-chain assignment and restraints measurement will be discussed. I will also discuss the different recoupling sequences which can be used with these experiments.

***Friday, Apr 22<sup>nd</sup> 2022***

***11:30 AM***