

Sr. No	Lecture Title	Description	Category
Segment 6: Molecular Dynamics & Simulation			
1	Molecular Dynamics of Protein Using GROMACS	<ul style="list-style-type: none"> • Introduction to GROMACS and its installation. • Preparation of the protein for simulation. • Energy minimization • Complete simulation of the protein. 	Protein Simulation
2	Molecular Dynamics of Protein-Ligand Complex Using GROMACS	<ul style="list-style-type: none"> • Preparation of the protein and ligand for simulation. • Energy minimization. • Complete simulation of the complex. 	Protein-Ligand Complex Simulation
3	Molecular Dynamics of Protein-Protein Complex Using GROMACS	<ul style="list-style-type: none"> • Preparation of the protein-protein-complex for simulation. • Energy minimization. • Complete simulation of the complex. 	Protein-Protein Complex Simulation
4	Molecular Dynamics of DNA-Ligand Complex Using GROMACS	<ul style="list-style-type: none"> • Preparation of the DNA and ligand for simulation. • Energy minimization. • Complete simulation of the complex. 	DNA-Ligand Complex Simulation
5	Molecular Dynamics of RNA-Ligand Complex Using GROMACS	<ul style="list-style-type: none"> • Preparation of the RNA and ligand for simulation. • Energy minimization. • Complete simulation of the complex. 	RNA-Ligand Complex Simulation