



Webinar on “Introduction to Molecular Dynamics (MD) Simulation for Bioinformatics Application”



Call-to-action: Tired of tutorials showing step-by-step commands for running MD simulation??? Dive into the theory behind MD simulation, force fields and applications through our **free webinar!**

What to expect?

- **Introduction to Molecular Dynamics**
- **Role of Force fields in MD simulation of proteins**
- **Types of MD simulation along with different tools**
- **Applications of MD simulation in Bioinformatics**
- **Challenges and limitations of MD simulation**

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Date : 25.05.2024 (Saturday)

Time : 10:00 AM to 12:30 PM

Platform : Google Meet

(Credentials will be sent later)

Speaker

Dr. Kamalesh Damodaran

Incubatee (Autosimulate - Startup), VITTBI



[Scan here for Free Registration](#)



[For Further Details Contact](#)

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