



# Hands on Training and Workshop on Computational Drug Discovery

## **Important Dates**

Registration open: 10 Jan 2024 Last date for Registration: 25 Feb 2024 Batch-1: 28–29 Feb, 2024 Batch-2: 01–02 Mar, 2024

# **Registration fees**

UG Students: 500/-PG Students: 600/-Research Scholars: 1000/-Faculty & Industry : 1500/-

## **Course Content**

## Day-1

- Introduction to Drug Discovery and Development
- Disease Target identification
- Bioinformatics Methods to identify disease targets
- Sequence Analysis
- Protein Structure Prediction
- Molecular Modelling
- Active Site prediction
- Hands on Session

# Day-2

- Lead discovery and Analog Based Drug Design (Rational approaches to lead discovery based on traditional medicine, Random screening, Non-random screening, lead discovery based on drug metabolism, lead discovery based on clinical observation.)
- Quantitative Structure Activity Relationship (QSAR)
- Virtual Screening techniques: Drug likeness screening, Concept of pharmacophore mapping and pharmacophore based Screening,
- Molecular docking: Rigid docking, flexible docking, manual docking, Docking based screening.

#### Venue

DR Bioscience & Scientific Bio-Minds

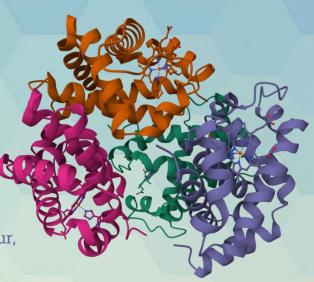
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# (Registration Fees including Certificate)

For Registration and queries contact Prashantha C N (+91-9844158444) Email: scientificbiominds@gmail.com

# Learning Outcomes

- Hands on Training in essential computing skills in drug discovery
- Working with bioinformatics software's, Tools and databases
- Explore Chemoinformatics and Drug Discovery concepts
- Mechanism of Structure and Ligand based drug discovery process



# **Registration fees**



# **Registration Form**

Participants are requested to fill the registration form along with course fees before: 25 Feb 2024. https://forms.gle/dmnqmbMM3M3SZ34w7