



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/-**

Venue

DR Bioscience & Scientific Bio-Minds
#17, KK Chambers, 1st Floor, Bettahalasur,
Bengaluru - 562157
Karnataka, India
Email: scientificbiominds@gmail.com
Mobile: +91-9844158444

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.

(Registration Fees including Certificate)

For Registration and queries contact
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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>