

Cloud-based Hands-on Workshop:

Computational Structure-based Screening and Explicit Molecular Dynamics

Date: 05th & 06th Nov 2020 ~ Time: 10:00 - 16:00 IST



SCHRODINGER.

**Cloud-based Hands-on Workshop:
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and Explicit Molecular Dynamics**

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Outline: Two-day cloud-based hands-on workshop targeting structure-based drug designing. Participants will get practical experience and in-person guidance in using the Maestro GUI, covering the organic molecule sketching, protein selection, preparation, and screening for hit identification of molecules against therapeutic targets. The workshop will also include a brief recap of background theory for Molecular mechanics, Molecular Docking, and Molecular Dynamics via case studies on the real-time industrial projects.

Day 1 - 05th Nov 2020

Time	Session
9:30	Technical set-up
9:50	Audio & Visual check
10:00	1. Opening: Agenda and split into breakout rooms
10:10	Breakout 1: Introductions and open software
10:30	2. Maestro GUI: Building Molecules and Enumeration
11:00	Breakout 2: Build Challenges from Schrodinger
11:20	3. Ligand Preparation and ADME
11:35	Breakout 3: Ligand, LigPrep
11:45	Intro to Molecular Docking, Protein, and Ligand Preparations
12:10	4. Protein Preparation and Grid Generations
12:30	Breakout 4: Prepare protein and grid generations and molecular docking
13:00	Break: Continue to use software during the break
14:00	Welcome Back
14:10	5. Molecular Docking
14:25	Breakout 5: Analyzing 3D molecular docking analysis
14:40	6. Molecular Docking Analysis
14:55	Breakout 6: 2D molecular docking analysis
15:30	7. Full Revision of Day 1
15:55	Breakout 7: Full session practice
16:00	Wrapup
16:00	Finish

Day 2 - 06th Nov 2020

Time	Session
10:00	1. Opening: Agenda and molecular dynamics theory
10:30	Breakout 1: Docking results and ligand docking
10:45	2. Desmond Introduction
11:00	Breakout 2: Protein-ligand complex formation, system building
11:10	3. Molecular Dynamics Demo
11:25	Breakout 3: MD submission
11:35	4. Desmond Trajectory Visualization
11:45	Breakout 4: Post-clash analysis
12:15	Break: Continue to use software during the break
14:00	Welcome Back
14:10	5. Desmond Molecular Simulation Analysis
14:20	Breakout 5: Analyzing molecular dynamics analysis
14:45	6. Organic Molecules Enumeration
15:00	Breakout 6: Reaction-based enumeration, library enumeration, AUMs
15:30	7. Full Revision of Day 2
15:55	Wrapup
16:00	Finish

Workshop Speakers:
Dr. Pritesh Bhat
Dr. Prajwal Nandekar
Dr. Koushik Kasavajhala
Ms. Shelvia Malik
Mr. Kishore V
Mr. Vinod D

Registration: Registration is open to 30 per. (Organis), please register here - <https://bit.ly/fy6da>

General Inquiries:
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The workshop was organized by Schrodinger and Central Sophisticated Instrumentation facility (CSIF), University of Calicut. Prof. (Dr.) M K Jayaraj, Hon'ble Vice Chancellor, University of Calicut inaugurated the workshop. Dr. Denoj Sebastian, Asst. Professor, Dept. of Life Science, Prof. (Dr.) Jos T Puthur, Director, CSIF, Dr. Raghu R, Vice president, Schrodinger attended the function. 31 participants from various universities/research institutes participated in the Workshop.

Inauguration:

Prof. (Dr.) M K Jayaraj

Hon'ble Vice Chancellor, University of Calicut

Workshop Speakers:

- Dr. Pritesh Bhat
- Dr. Prajwal Nandekar
- Dr. Koushik Kasavajhala
- Ms. Shelvia Malik
- Mr. Kishore V
- Mr. Vinod D



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Organized by
Schrodinger and Central Sophisticated Instrumentation Facility (CSIF), University of Calicut
on 05th & 06th Nov 2020

PROGRAMME

Date: 05-11-2020

9:00 am - Technical updation regarding the Workshop
(All Participants should Login)

9:30 am - Inauguration of the Workshop

Inaugural Message
Prof. (Dr.) M K Jayaraj
Hon. Vice Chancellor,
University of Calicut

Welcome
Dr. Denoj Sebastian
Asst. Professor
Dept. of Life science
University of Calicut

Introduction about CSIF
Prof. (Dr.) Jos T Puthur
Director
CSIF
University of Calicut

Felicitations
Dr. Raghu R
Vice President
Schrodinger

Technical sessions are led by Scientists and Technical experts from Schrodinger