

ONE WEEK SUMMER SCHOOL 2024

Hands-on Training in Computational and Animal Modeling Tools for Design & Discovery of Drugs for Neurodegenerative Disorders

2nd to 6th July 2024

Organized by
Department of Pharmaceutical Chemistry
and
Department of Pharmacology

Theme

Critical Concepts & Hands-on-Training:

- Molecular Docking
- ADME Screening
- Pharmacophore screening
- Network Pharmacology
- In-Vitro and In-Vivo pharmacological screening techniques

Eligibility

Any passionate learner is eligible to participate, preferably

- Faculty, Industry Personal
- Post-Doctoral Fellows
- Ph.D. Research Scholars
- M. Pharm. Students

Committee

Convener

Dr. Jagannath Sahoo, Dean, SPPSPTM

Workshop Coordinators

Dr. RajaSekhar Reddy Alavala

Dr. Pratigya Tripathi

Venue: SPPSPTM, SVKM's NMIMS Deemed to be University, V.L. Mehta Road, Vile Parle (W), Mumbai 400056.

Objectives

One week intensive summer school is aimed at meeting the objectives of

- Making the attendees able to understand the concept of in silico drug design for neurodegenerative disorders
- Getting trained on handling the applications of various computational techniques
- Demonstrate animal models and techniques for evaluation of drugs for treatment of neurodegenerative disorders

Registration

Online Registration Process

- Registration Fee: Rs. 1500/- (for students) & Rs. 2000 /- (for faculty)
- Working lunch included on all the days
- Fill the below registration form <https://forms.gle/4RNjsj9PcS2Cf9KL7>
- Limited to 30 seats (*On a first come, first serve basis*)
- Last Date to Register: 20th June 2024



For any queries, please contact: Workshop Coordinators

alavala.rajasekharreddy@nmims.edu (or) pratigya.tripathi@nmims.edu

Contact: 98852 71223, 89488 96950

Session No	Topic	Session No	Topic
Day-1: 02-07-2024 (Tuesday)		Day-2: 03-07-2024 (Wednesday)	
	Inauguration	5	Introduction to Network Pharmacology & its applications
1	Introduction to Molecular Modelling		Tea Break
	Tea Break	6	Building of Network Pharmacology Model:Protein-protein
2	Applications of Computational tools in Pharmaceutical Research		Lunch Break
	Lunch Break	7	Building of Network Pharmacology Model:Protein-gene
3	Drawing of the structures with various tools	8	Building of Network Pharmacology Model:Protein-drug
4	Conversion of 2D structures to 3D & Energy minimization		
Session No	Topic	Session No	Topic
Day-3: 04-07-2024 (Thursday)		Day-4: 05-07-2024 (Friday)	
9	Introduction to Molecular Docking & its process	13	Analysis of the Molecular Docking Results
	Tea Break		Tea Break
10	Protein Selection & preparation	14	Introduction to Pharmacophore modeling based screening
	Lunch Break		Lunch Break
11	Molecular Docking Hands-on for the protein-ligands	15	Hands-on for the Pharmacophore generation by LBDD
12	Molecular Docking Hands-on for the protein-protein	16	Hands-on for the Pharmacophore generation by SBDD
Session No	Topic	Session No	Topic
Day-5: 06-07-2024 (Saturday)			
17	Introduction to In-Vitro and In-Vivo pharmacological screening techniques		
	Tea Break		
18	Demonstration of Stereotaxic Surgery- Batch A & Demo on Fruit Fly Animal Model- Batch B		
	Lunch Break		
19	Demonstration of Stereotaxic Surgery- Batch B & Demo on Fruit Fly Animal Model- Batch A		
20	Demonstration of Behavioral Studies		
	Valedictory		

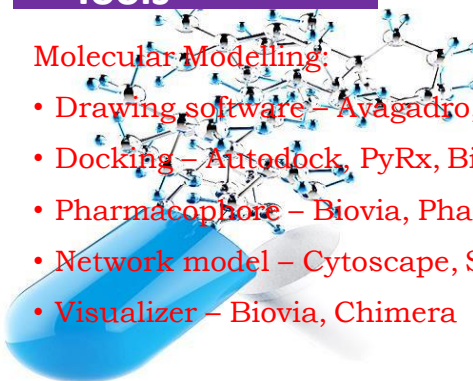
Registration

Name of the A/C Holder:	SVKM'S NMIMS
NEFT IFSC Code:	KKBK0000661
Bank Name:	Kotak Mahindra Bank
	JUHU - Vile Parle West
Account No:	2311578254
Type of Account:	Current
MICR Code:	400485022
SWIFT Code:	KKBKINBB

Tools

Molecular Modelling:

- Drawing software – Avagadro, ChemSketch
- Docking – Autodock, PyRx, Biovia
- Pharmacophore – Biovia, Pharmmaper
- Network model – Cytoscape, STRING
- Visualizer – Biovia, Chimera



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