PAYMENT

REGISTRATION FEES

Early-bird Registration Deadline 31st August 2018

Workshop Workshop + Conference

·Early Bird International **USD 300 USD 80 USD 350** International **USD 100** •Early Bird lecturers/ Researchers/others RM 250 RM 900 RM 1000 Lecturers/ Researchers/others RM 300 ·Early bird local students **RM 200 RM 700** RM 850 Local students RM 250

LIMITED TO 30 Participants!

Package

REGISTRATION FORM

Prefix	
First Name	
Last Name	
E-Mail	
Quantity Ticket Type	
Total Paid	
Cell Phone	
Gender	
Country	
Area of Research	
Food Preference	
Job Title	
Company /Institute	

3rd Announcement



PRE-CONFERENCE WORKSHOP ON MOLECULAR DOCKING IN DRUG DISCOVERY RESEARCH

In conjunction with



3rd DECEMBER 2018 8.30 AM - 5.00 PM **BILIK SAINTIS GEMILANG, FACULTY OF SCIENCE, UPM**

SPEAKERS



ZHANG (JAPAN) Leader of Structural Bioinformatics Team Center for Life Science Technologies, RIKEN, Japan.

PROF. KAM Y.



DR. LAM KOK WAI, (UKM) Senior Lecturer Drugs and Herbal Research Centre Faculty of Pharmacy, Univ<mark>ersiti Ke</mark>bangsa<mark>an</mark> Malaysia

Organized by the Department of Chemistry

www.icddtm18.macr-cancer.ora

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Payment Inquiries

Ms Nura (+60173059630)

For further information or query, please contact:

The ICDDTM '18 Secretariat **Department of Chemistry Faculty of Science** Universiti Putra Malaysia 43400 UPM, Serdang, Selangor Malaysia

REGISTRATION

Tel: Mrs. Nurhidayu (+60173157977)

Fax: 03-89432508

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INTRODUCTION

This one-day workshop is exclusively tailoredmade for beginners (students and researchers) with a diverse disciplines related to drug discovery or relevant applications. The main objective is to provide the participants an introduction to basic molecular docking with hands-on practice, thus improving understanding of molecular mechanisms in specific biological activity. This computational technique may also help to conceive and evaluate potential molecules to become actual drugs. Last but not least, it would also increase the productivity by offering time-saving and costeffective solutions in the development of new pharmaceutical products.

TOPICS TO BE COVERED

- •Theory on molecular docking and computational methods
- •Selection of target receptor (protein–protein interactions
- •(PPIs) and protein-small molecule interactions)
- •Scoring functions and protein-ligand interaction
- •Hands-on session on performing molecular docking and analysis of the docking results

WHO SHOULD ATTEND

The workshop is suitable for all personnel involved in natural product, organic synthesis, pharmaceutical, bio-medical, life sciences, structural biology, computational biology or those people who are intended to work with bioactive-small organic molecules.

DATE VENUE

: 3rd DECEMBER 2018 :BILIK SAINTIS GEMILANG, FACULTY OF SCIENCE, UPM

3rd DECEMBER 2018		
0830 - 0900	Registration	
0900 - 1000	Introduction of Computational Methods in Drug Discoveries	
	(Prof. Kam Y.J. Zhang, Japan)	
1000 - 1020	Tea Break & Photography Session	
1030 - 1230	Practical I : Get to know Your Drug Target! Finding the Right Protein and Principles Behind Docking (Dr. Lam Kok Wai, UKM)	
1230 - 1400	Lunch	
1400 - 1630	Practical II: Performing Molecular Docking and Data Analysis (Dr. Lam Kok Wai, UKM)	
1630 - 1700	Refreshment	

