

244th OMICS Group Conference

Scientific Program



3rd International Conference on

Medicinal Chemistry & Computer Aided Drug Designing

December 08-10, 2014 DoubleTree by Hilton Hotel San Francisco Airport, USA



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OMICS Group Conferences

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Keynote Forum

09:55-10:00 Introduction

10:00-10:25 **Concepción González-Bello**

University of Santiago de Compostela, Spain

10:25-10:50 **Patrick Y S Lam**

Drexel University, USA

Coffee Break 10:50-11:05 @ Foyer

11:05-11:30 **Thorsten Nowak**

C4X Discovery Holdings PLC., UK

Track 1: Rational Drug Design

Track 2: Computer-Aided Drug Design and Structure Determination

Session Chair: Patrick Y S Lam, Drexel University, USA

Session Co-Chair: Concepción González-Bello, University of Santiago de Compostela, Spain

Session Introduction

Title: Predicting ligand binding affinity: A comparative study on the use of docking vs. Bayesian categorization and random forest recursive partitioning

11:30-11:50 **David C Kombo**, Proteostasis Therapeutics, Inc., USA

Title: In silico screening for anti-HPV agents using pharmacophore models

11:50-12:10 **Tatsuya Takagi**, Osaka University, Japan

Title: NMR driven conformational design - A powerful drug design tool

12:10-12:30 **Thorsten Nowak**, C4X Discovery Holdings PLC., UK

Title: Computer aided design and optimization of kinase and phosphatase inhibitors

12:30-12:50 **Kal Ramnarayan**, Sapient Discovery, LLC., USA

Title: The use of epitopes against schistosomiasis: The role of molecular modelling

12:50-13:10 **Moacyr Comar Junior**, Federal University of São João del Rei, Brazil

Title: Bacterial type II dehydroquinase enzyme: From the reaction mechanism to the structure-based design of inhibitors

13:10-13:30 **Concepción González-Bello**, University of Santiago de Compostela, Spain

Title: Structure-based design of covalent inhibitors or protein-protein interactions

13:30-13:50 **Maurizio Pellecchia**, Sanford-Burnham Institute for Medical Research, USA

Lunch Break 13:50-14:35 @ Tiburon-Sausalito

Track 3: Quantitative Structure-Activity Relationships

Track 4: Advanced Medicinal Chemistry

Track 8: Drug Interactions and Drug Metabolism

Track 9: Neurodegenerative Diseases

Session Chair: Thorsten Nowak, C4X Discovery Holdings PLC., UK

Session Co-Chair: Istvan J Enyedy, Biogen Idec, USA

Session Introduction

Title: Revisiting INH: QSAR-based design of new anti-tubercular compounds

14:35-14:55 **Filomena Martins**, Universidade de Lisboa, Portugal

Title: In vitro COX inhibitory activity of novel α -aminoarylpropionic acid derivatives

14:55-15:15 **A G Zhamharyan**, Yerevan State Medical University, Armenia

Title: New developments in docking and scoring

15:15-15:35 **Istvan J Enyedy**, Biogen Idec, USA

Title: Comparative molecular modeling study between the pre-fusion and post-fusion conformations of newcastle disease virus: Homology modeling, and virtual screening

15:35-15:55 **Mohammed A Khedr**, King Faisal University, KSA

Coffee Break 15:55-16:10 @ Foyer

Title: Peptide-displaying phage technology in breast cancer diagnosis

16:10-16:30 **Thaise Gonçalves Araújo**, Federal University of Uberlandia, Brazil

Title: Discovery of small molecule blockers of protein-protein interactions using DNA-encoded small-molecule libraries

16:30-16:50 **Nils Hansen**, VipergenApS, Denmark

Title: Neuroprotective properties of compound isolated from *Dianthus superbus* L.

16:50-17:10 **Choong Je Ma**, Kangwon National University, Korea

Title: Investigation of human CA1-induced toxicity in the *Drosophila* model of motor neuron degeneration

17:10-17:30 **Jian Liu**, Xi'an Jiaotong-Liverpool University, China

Title: Applications of Proteochemometrics - From Species Extrapolation to Cell Line Sensitivity Modelling

17:30-17:50 **Isidro Cortes-Ciriano**, Bioinformatique Structurale Institut Pasteur, France

Title: Data and Compute Intensive eScience Approaches in Computational Medicinal Chemistry

17:50-18:10 **Scott James Lusher**, Netherlands eScience Center, Netherlands

Title: Synthesis, Biological Evaluation and 3d Qsar of some Novel Benzimidazole Derivatives as Anti Microbial Compounds

18:10-18:30 **Sonal Dubey**, Krupanidhi College of Pharmacy, India

18:30-19:30 Cocktails sponsored by Journal of Medicinal Chemistry @ Tiburon-Sausalito

Day 2 December 09, 2014

Sierra Ballroom

Keynote Forum

09:30-09:55 Pierre Falson

IBCP, France

Track 5: Cancer Research in Medicinal Chemistry

Track 6: New Approaches in Drug Discovery

Session Chair: Andrew B McElroy, Eligochem Ltd., UK

Session Introduction

09:55-10:15 Title: Research and development for preparation and preclinical evaluation of novel low-molecular-weight phospho sugar antitumor agents targeting IER5/Cdc25B

Mitsuji Yamashita, Shizuoka University, Japan

10:15-10:35 Title: Strategies to optimize and use bacterial cytochromes p450 for drug discovery & development

Nico P E Vermeulen, VU University Amsterdam, Netherlands

10:35-10:55 Title: New agonists of the CB₂ cannabinoid receptor: Discovery of a new class of analgesic compounds

Pier Giovanni Baraldi, Università di Ferrara, Italy

Coffee Break 10:55-11:10 @ Foyer

11:10-11:30 Title: Polar drugs

Andrew B McElroy, Eligochem Ltd., UK

11:30-11:50 Title: Drug discovery against category A-C pathogens through MEP pathway

Prabakaran Narayanasamy, University of Nebraska Medical Center, USA

11:50-12:10 Title: Discovery of novel lead compounds by large scale diverse encoded chemical libraries

Jin Li, HitGen Ltd., China

12:10-12:30 Title: *In silico* approach to predict ADME-Tox properties of small organic molecules: Challenges and opportunities for drug discovery

Maria Miteva, University Paris Diderot, France

Track 7: Drug Development and Delivery System

Track 10: Receptors and Inhibitors

Session Chair: Niren Murthy, University of California, USA

Session Co-Chair: Peter Teriete, Sanford-Burnham Institute for Medical Research, USA

Session Co-Chair: Nico P E Vermeulen, VU University Amsterdam, Netherlands

Session Introduction

12:30-12:50 Title: *In vivo* delivery of transcription factors with chemically modified oligonucleotides

Niren Murthy, University of California, USA

12:50-13:10 Title: Eradication of asbestos tumors *in vivo* with histone deacetylase inhibitors-polymer conjugate nanoparticles for acid-responsive drug delivery

Philippe Bertrand, Institut de Chimie des Milieux et Matériaux de Poitiers, France

Lunch Break 13:10-14:00 @ Tiburon-Sausalito

14:00-14:20 Title: The σ_1 receptor as target for novel drugs

Bernhard Wünsch, Westfalian Wilhelms-University of Münster, Germany

14:20-14:40 Title: Using Computer-Aided Drug Design (CADD) techniques to optimize the natural product-derived phenylmethylidene-hydanto in scaffolds as promising antimetastatic leads

Mudit Mudit, D'Youville College School of Pharmacy, USA

14:40-15:00 Title: Design and development of small peptidomimetics of RXFP1 for the treatment of acute heart failure

Akhter Hossain, University of Melbourne, Australia

15:00-15:20 Title: Exploring pharmacological potential of Brazilian plants: SAM database-A tool for recording and comparison of molecules isolated from plants of the Brazilian semiarid

Bruno Andrade, State University of Southwest Bahia, Brazil

Coffee Break 15:20-15:35 @ Foyer

15:35-15:55 Title: Novel chemistry-based tools to study epigenetic enzymes in inflammation

Frank J Dekker, Pharmaceutical Gene Modulation, The Netherlands

15:55-16:15 Title: Current SAR on HIV: The flow from phenotypic assays via medicinal chemistry to *in silico* design

Peter Teriete, Sanford-Burnham Institute for Medical Research, USA

16:15-16:35 Title: Structure guided design and synthesis of SAR107375A, a selective and potent dual thrombin and factor Xa inhibitor

Jerome Meneyrol, Sanofi-Aventis R&D, France

16:35-16:55 Title: Molecular modeling approach to investigate the binding mode of 4-nerolidylcatechol into two subtypes of matrix metalloproteinases

Kely Medeiros Turra, University of São Paulo, Brazil

14:30-15:30 Poster Presentations @ Tiburon-Sausalito

17:30-18:30 Cocktails sponsored by Journal of Drug Designing @ Tiburon-Sausalito

Day 3 December 10, 2014

Sierra Ballroom

Track 11: Membrane Proteins as Pharmaceutical Targets

Session Chair: Victor J Hruby, University of Arizona, USA

Session Co-Chair: Pierre Falson, IBCP, France

Session Introduction

- 09:00-09:20 **Title: Tracking anti-tumor drugs: Ruthenium(II)-cyclopentadienyl complexes as promising agents**
M Helena Garcia, Universidade de Lisboa, Portugal
- 09:20-09:40 **Title: TAAR1 ligands as prospective neuroleptics: From D-neuron study**
Keiko Ikemoto, Iwaki Kyoritsu General Hospital, Japan
- 09:40-10:00 **Title: Small-molecule modulators of thiamine transport in pathogenic bacteria**
Anna K H Hirsch, University of Groningen, The Netherlands
- 10:00-10:20 **Title: The behavior of detergents around membrane proteins is more complex than supposed, as revealed by a new method of quantitation**
Pierre Falson, IBCP, France
- 10:20-10:40 **Title: Design of multivalent ligand for the detection and treatment of disease**
Victor J Hruby, University of Arizona, USA
- 10:40-11:00 **Title: Polymer 'ruthenium-cyclopentadienyl' conjugates: A new approach to fight cancer**
Andreia Valente, Universidade de Lisboa, Portugal

Coffee Break 11:00-11:15 @ Foyer

Track 12: Recent Research and Developments

Session Chair: Jetze J Tepe, Michigan State University, USA

Session Co-Chair: Bin Xu, Virginia Tech, USA

Session Introduction

- 11:15-11:35 **Title: Design of mechanistically distinct proteasome inhibitors for the treatment of multiple myeloma**
Jetze J Tepe, Michigan State University, USA
- 11:35-11:55 **Title: Molecular characterization and design of a key new hormone, irisin**
Bin Xu, Virginia Tech, USA
- 11:55-12:15 **Title: Structure-based discovery of new modulators targeting nuclear X receptor alpha for cancer therapy**
Ying Su, Sanford-Burnham Medical Research Institute, USA
- 12:15-12:35 **Title: Privileged heterocycles by palladium-catalyzed aerobic oxidative isocyanide insertion**
Romano V A Orru, VU University Amsterdam, Netherlands

Lunch Break 12:35-13:20 @ Tiburon-Sausalito

- 13:20-13:40 **Title: Visual binding: A radically new concept to support the medicinal chemist's quest for innovative NMEs**
Carsten Detering, BioSolve IT Inc., USA
- 13:40-14:00 **Title: A G-quadruplex/i-motif switch in the HRAS promoter as target for anthrathiophenediones that show a strong anti-proliferative activity in urinary bladder cancer cells**
Luigi E Xodo, University of Udine, Italy
- 14:00-14:20 **Title: Multi-target approach to anti-inflammatory drugs - in silico and medicinal chemistry**
Eugen Proschak, Goethe-University of Frankfurt, Germany
- 14:20-14:40 **Title: Identification and optimization of tertiary sulfonamides as RORc inverse agonists**
Benjamin P Fauber, Genentech, Inc., USA
- 14:40-15:00 **Title: Massive changes to the biophysical properties of DNA upon binding antiviral polyamides**
Gaofei He, University of Missouri-St. Louis, USA

Coffee Break 15:00-15:15 @ Foyer

- 15:15-15:35 **Title: The communion of medicinal chemistry and nanotechnology in anticancer therapeutics**
Debatosh Majumdar, Glycosyn LLC., USA
- 15:35-15:55 **Title: Evaluation of *Satureja hortensis* leaves essential oil pharmacological activities**
A P Manjiky, Yerevan State Medical University, Armenia
- 15:55-16:15 **Title: Phytochemicals and antioxidant capacities from *Dacryodes rostrata* fruits**
Prasad K N, Monash University Malaysia, Malaysia
- 16:15-16:35 **Title: Selection and characterization of RNA aptamers targeting the genomic 5'-UTR in the dengue virus**
Adriana Freitas Neves, Universidade Federal de Goiás, Brazil
- 16:35-16:55 **Title: SurR9C84A exhibits cardioprotective effects against melphalan induced cardiotoxicity in primary human cardiomyocytes**
Ajay Ashok, Deakin University School of Medicine, Australia

Award Ceremony

Bookmark your dates

4th International Conference on

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November 02-04, 2015 Atlanta, USA



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