

Program

2nd International Symposium on Computational Materials and Biological Sciences

September 1 - September 2, 2015, Nagoya University, Japan

Tuesday, Sep. 1	
13:30-13:40	Opening: Kenji Yasuoka (<i>Keio University</i>)
Session I (Chair: Kholmirzo Kholmurodov)	
13:40-14:10	Roman Efremov (<i>Institute of Bioorganic Chemistry, Moscow</i>)
Invited #1	Cell Membranes as Stochastic Dynamic Systems: a Computational View.
14:10-14:40	Takao Yoda (<i>Nagahama Institute of Bio-Science and Technology</i>)
Invited #2	Simulation Study on Conformations of Antimicrobial Peptide Cryptdin 4 and its Mutants
14:40-15:40	Poster Previews (Chair: Donguk Suh)
15:40-18:40	Poster Presentations
19:00-21:00	Banquet
Wednesday, Sep. 2	
Session II (Chair: Takao Yoda)	
9:00-9:30	Anton Polyansky (<i>Institute of Bioorganic Chemistry, Moscow</i>)
Invited #3	Single-Point Mutations in PDGFRA Transmembrane Domain Induce Repopulation of an Ensemble of its Homodimer States and Modulate Basal Activity of the Receptor
9:30-10:00	Satoshi Ohmura (<i>Hiroshima Institute of Technology</i>)
Invited #4	Charge Dynamics in Light-Harvesting Organic Supermolecule
10:00-10:30	Vladimir Palyulin (<i>Moscow State University, Moscow</i>)
Invited #5	Molecular Dynamics Simulation and QSAR Studies of Ligand-Receptor Interactions in Ionotropic Glutamate Receptors
10:30-11:00	Coffee Break
Session III (Chair: Satoshi Ohmura)	
11:00-11:30	Yoshimichi Andoh (<i>Nagoya University</i>)
Invited #6	Large-Scale All-Atom Molecular Dynamics (MD) Simulations of Biomolecular Systems Using a General Purpose MD Simulation Program, MODYLAS
11:30-12:00	Noriyoshi Arai (<i>Kinki University</i>)
Invited #7	Self-Assembly of Diblock and Triblock Janus Nanoparticles in Nanotubes
12:00-12:20	Donguk Suh (<i>Keio University</i>)
Oral #1	Molecular Dynamics Simulation of Condensation on Elongated Nanocuboids and Nanotubes
12:20-12:30	Closing: Tomoyuki Yamamoto (<i>Waseda University</i>)