

「インシリコ創薬」分野 口頭発表
Selected Oral Presentations
(*In silico* drug discovery)

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1. [P2-05] Improving 3D-QSAR in AutoGPA using Gaussian-distributed physicochemical descriptors
Yoshirou Kimura
Life Science Department, MOLSYS Inc.
2. [P2-07] Construction of FMO IFIE-database
Chiduru Watanabe
RIKEN Center for Life Science Technologies
3. [P2-09] Using multiple molecular fingerprints for improvement of drug activity predictions
Yusuke Matsuyama
Tokyo Institute of Technology
4. [P2-12] Quick and precise homology modeling method of GPCRs
Mika Nabeno
Mitsubishi Tanabe Pharma Corporation
5. [P2-13] KampoDB: An integrated platform for mode-of-action analysis and repositioning of natural medicines
Ryusuke Sawada
Medical Institute of Bioregulation, Kyushu University
6. [P2-17] The Contribution of Entropy in Drug-Protein Binding
Takeshi Tanaka
Interprotein Corporation
7. [P2-19] Computer simulation-based prediction of drug-induced arrhythmia by evaluating repolarization reserve
Shingo Murakami
Dept. of Physiology Sch. of Medicine, Toho Univ.
8. [P2-21] A Genetic Approach to Deep Learning in Prediction of Molecular Properties
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