

「分子認識と分子計算」分野 口頭発表
Selected Oral Presentations
(Molecular recognition and molecular modeling)

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1. **[P1-06] Flexible Docking using Replica-Exchange Molecular Dynamics Simulation**
Suyong Re
RIKEN Quantitative Biology Center (QBiC)

2. **[P1-10] Validation of Epigenetic Therapeutic Target Proteins for Homogenous Assay Performance**
Masato Yonezawa
Active Motif Inc., Carlsbad CA, USA

3. **[P1-12] Finite-size effect on the charging free energy in the alchemical perturbation and “warp drive” method**
Toru Ekimoto
Graduate School of Medical Life Science, Yokohama City University

4. **[P1-23] Comparing two molecular dynamics simulation trajectories in terms of residue-residue interaction**
Chie Motono
Molecular Profiling Research Center for Drug Discovery, AIST

5. **[P1-36] Interaction Analysis between Beta-Secretase and its Inhibitors by Fragment Orbital Method**
Norihito Kawashita
Faculty of Science and Engineering, Kindai University

6. **[P1-39] Characteristics of Biomolecule Dynamics under the Crowding Environment of Cytoplasm Discovered by Massive All-atom simulation and Big-data analysis**
Isseki Yu
iTHES Research Group, RIKEN

7. **[P1-44] Structure and Dynamics of RNA Aptamer to Human Immunoglobulin G**
Hisae Yoshida
College of Engineering., Nihon University