

「創薬データサイエンス」分野Selected Oral Presentations
(Drug Discovery Data Science)モデレーター: 池田 和由 Kazuyoshi Ikeda
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口頭発表演題 (1演題 10分:発表 8分、質疑 2分)

1. [P7-08] Prediction of Compound-Protein Interactions and Visualization Based on Graph Convolutional Networks
池口 茉莉恵 Marie Ikeguchi
Kyoto University
2. [P7-03] Predicting drug indications and therapeutic target modules based on disease similarity by interpretable machine learning models
澤田 隆介 Ryusuke Sawada
Kyushu Institute of Technology
3. [P7-09] Development of data curation and integration protocol for the chemical library in early drug discovery
清水 祐吾 Yugo Shimizu
Keio University
4. [P7-10] Clustering therapeutic drugs based on similarities of indications and side effects reported in public database
新川 龍太郎 Ryutaro Shinkawa
Mie University School of Medicine
5. [P7-11] Drug Discovery Raid Battle 2018: an open challenge to discover PD-1/PD-L1 small-molecule inhibitors
山本 一樹 Kazuki Yamamoto
Isotope Science Center, University of Tokyo
6. [P7-05] The Multiple Representation of Protein Sequence Motifs Using Sequence Binary Decision Diagrams
加藤 博明 Hiroaki Kato
National Institute of Technology, Hiroshima College
7. [P7-12] Automatic Reading of Tables and Figures in Scientific Papers
進藤 裕之 Hiroyuki Shindo
Nara Institute of Science and Technology