

**「医薬品研究と ADMET」分野**  
**Selected Oral Presentations**  
(Information and computing approach for drug design and ADMET study)

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**口頭発表演題** (1 演題 20 分: 発表 15 分、質疑 5 分)

1. **[P5-12] Cluster Gauss-Newton method for efficiently estimating multiple sets of parameters: Application to Physiologically-Based Pharmacokinetic models**  
**速水 謙 Ken Hayami**  
国立情報学研究所, 総合研究大学院大学  
National Institute of Informatics, SOKENDAI
  
2. **[P5-13] Prediction of Health Effects of Food Peptides and Elucidation of The Mode-of-action Using Multi-task Graph Convolutional Neural Networks**  
**福永 一貴 Itsuki Fukunaga**  
九州工業大学  
Kyushu Institute of Technology
  
3. **[P5-15] Free Energy Landscapes of Cyclic Hexapeptide Diastereomers by Multicanonical Molecular Dynamics Simulations**  
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