

Seminar “Molecular modeling in drug discovery”

5th June 2015, Conference Hall, Latvian Institute of Organic Synthesis

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| 9.45-10.00 | <i>Registration and coffee</i> |
| 10.00-10.05 | Welcome remarks |
| 10.05-10.50 | Amedeo Caflich (University of Zürich, Zürich, Switzerland) Molecular dynamics in drug design |
| 10.50-11.35 | Christoph Rademacher (Max Planck Institute of Colloids and Interfaces, Berlin, Germany) Fragment-based ligand design of C-type lectin ligands |
| 11.35-12.00 | <i>Discussions and coffee</i> |