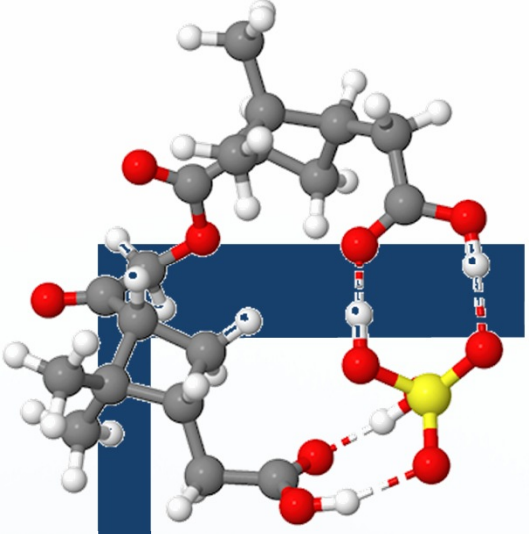
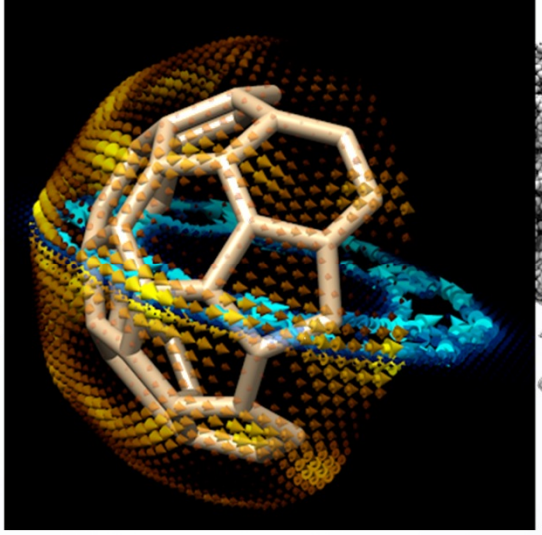
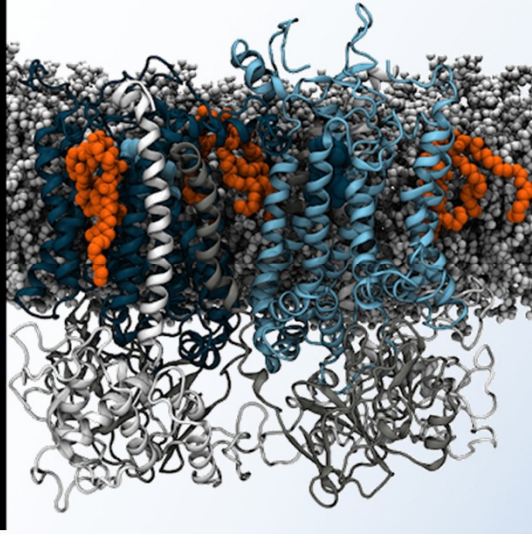

COMPUTATIONAL CHEMISTRY DAYS 2018

University of Helsinki, Finland, 28-29 May 2018

https://www.helsinki.fi/en/conferences/computational-chemistry-days-2018

The annual Computational Chemistry Days bring together all scientists enthusiastic about computational chemistry – including physics, bio, and pharma. Participation of experimental colleagues is strongly encouraged to foster chances for collaboration. The focus themes of this year's meeting include machine-learning techniques, pharmaceutical chemistry linked to industrial interests, catalysis, energy, and computational method development. Welcome!

Plenary Speakers

Tristan Bereau, Max Planck Institute for Polymer Research
Roland Lindh, Uppsala University
Antti Poso, Universities of Tübingen and Eastern Finland
Edina Rosta, King's College London

Invited Speakers

Jaakko Akola, Alex Bunker, Gerrit Groenhof, Karoliina Honkala, Mark Johnson, Kari Laasonen, Kai Nordlund, Esa Räsänen, Dage Sundholm, Juha Vaara, Hanna Vehkamäki

Important Dates

2 May 2018: Abstract submission for contributed talks and posters
7 May 2018: Deadline for registration

Organisers and Contact

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