

1st Nordic Conference on Computational Chemistry 2022

February 1-2, Gothenburg, Sweden

Preliminary program

Tuesday, February 1

09:30 **Registration**
10:00 Welcome and Introduction

MD

10:10 *Erik Lindahl*, Stockholm University
11:10 Selected talk
11:30 Flash poster presentations
12:30 *Lunch*

Structure based and Ligand based design

13:30 *Anna Linusson*, Umeå University
14:10 *Jens Carlsson*, Uppsala University
14:50 Selected talk
15:10 Posters and coffee

16:00 *Ruth Brenk*, Bergen University
16:40 Selected talk
17:00-17:20 Selected talk

19:00 Dinner

Wednesday, February 2

Advances in QSAR and Machine Learning

9:00 *Per-Ola Norrby, AstraZeneca*

10:00 *Nadine Schneider, Novartis*

10:40 Coffee and posters

Modelling of complex systems

11:10 *Lucie Delemotte, SciLifeLab*

11:50 *Giulio Tesei, University of Copenhagen*

12:30 Lunch

Protein and peptide design

13:30 *Simone Fulle, NovoNordisk*

14:10 *Ingemar André, Lund University*

14:50 Selected talk

15:10 Concluding remarks and poster prize

15:20 Coffee

15:40-16.40 Tour of the AstraZeneca site with the amazing journey

Sponsor:

