

ICCS2025: back in Noordwijkerhout

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Abstract

This week the 13th International Conference on Chemical Structures took place (see also this Scholia overview or this overview of the full ICCS history). This is the conference I first joined 20 years ago as a PhD student presenting a poster (see these past blog posts). Of course, I am actually co-organizer nowadays (actually, co-treasurer). Organizing a meeting with just over 200 participants, and I like to thank Gerard and Willem in

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This week the 13th [International Conference on Chemical Structures](#) took place (see also [this Scholia overview](#) or [this overview of the full ICCS history](#)). This is the conference I first [joined 20 years ago](#) as a PhD student [presenting a poster](#) (see [these past blog posts](#)). Of course, I am actually co-organizer nowadays (actually, co-treasurer). Organizing a meeting with just over 200 participants, and I like to thank [Gerard](#) and [Willem](#) in particular, but also [Pieter](#), [Marcus](#), [Frank](#), [Jenke](#), and Frans Koeman who has helped us during the past three events.

The meeting started, as usual, with the [CSA Trust Mike Lynch Award](#), this year awarded to Prof. [Val Gillet](#) (see also [this press release](#)).

[This time](#), there were the following themes, where the first was by far the most dominant theme:

- Artificial Intelligence, Machine Learning, and QSAR (five sessions)
- New Modalities and Large Chemical Data Sets (one session)
- Advanced Cheminformatics Techniques (two sessions)
- Integrative Structure-Based Drug Design (two sessions)

My contribution this time was a poster for the [VHP4Safety](#) project, but more about that later.

Like last time, I have been annotating speakers with identifier and accounts, if they provided those:

Wednesday, June 4, 2025

8:30 – 10:00 **Advanced Cheminformatics Techniques 1** Rotonde
Chair: Egon Willighagen   , Maastricht University

8:30 – 9:00 **Transformers for molecular property prediction: Domain adaptation efficiently improves performance**
Afnan Sultan , Saarland University

9:00 – 9:30 **Navigating Synthon Space: Property-Driven Molecular Optimization for Pharmacokinetics**
Rafał Adam Bachorz, Simulations Plus

9:30 – 10:00 **Scaffold Hopping with Generative Reinforcement Learning**
Luke Rossen , Eindhoven University of Technology

10:00 – 10:30 BREAK Atrium

10:30 – 12:30 **Advanced Cheminformatics Techniques 2** Rotonde
Chair: Greg Landrum

10:30 – 11:00 **Honey, I shrunk the database: Making multi-billion compound libraries as small as possible** 
John Wilkinson Mayfield , NextMove Software

11:00 **Docking-based geometric graph models for kinase-ligand affinity**

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As you can see, it also includes PDFs, for both [talks](#) and [posters](#). At the time of writing, I collected PDFs of two presentations and five posters. Additions are still most welcome, ideally with DOI, so that they can be cited (doi:[10.5281/zenodo.15494630](https://doi.org/10.5281/zenodo.15494630), doi:[10.13140/RG.2.2.27441.90720](https://doi.org/10.13140/RG.2.2.27441.90720) doi:[10.5281/zenodo.15614295](https://doi.org/10.5281/zenodo.15614295), and doi:[10.13140/RG.2.2.36774.23365](https://doi.org/10.13140/RG.2.2.36774.23365))!

Finally, I like to remind everyone that there is again a [proceedings collection in the Journal of Cheminformatics](#), and presenters of oral and poster presentations are invited to submit their presented work to this collection.