

AI and Computational Drug Discovery

Thursday 22 June 2023

10.15 - 10.25	Welcome and plan for the day <i>Dr Martin Redhead (Exscientia)</i>
10.25 - 10.50	Demystifying AI: Intuition for how it works, when it works and what to watch out for <i>Dr Angelo Pugliese (BioAscent)</i>
10.50 - 11.15	Biological graph databases – information transfer for AI <i>Dr Alan Bilsland (University of Glasgow & Exscientia)</i>
11.15 - 11.40	Generative models for small molecule drug discovery <i>Aleksandra Kalisz (Exscientia) and Dr Francesca Vianello (Exscientia)</i>
11.40 - 12.00	Break
12.00 - 12.30	Breakout session
12.30 - 12.55	AI to solve difficult kinetic problems in receptor pharmacology <i>Dr Martin Redhead (Exscientia)</i>
12.55 - 13.40	Lunch
13.40 - 14.10	Breakout session
14.10 - 14.35	Model-based target pharmacology assessment (mTPA) – how AI can inform strategies for drug discovery <i>Dr Emile Chen (GlaxoSmithKline)</i>
14.35 - 15.05	Breakout session
15.05 - 15.20	Break
15.20 - 15.45	Quantitative systems pharmacology target modulation and clinical outcomes <i>Dr Valeriu Damian (GlaxoSmithKline)</i>
15.45 - 16.15	Breakout session
16.15 - 16.20	Wrap-up and confirmation of the Learning Outcomes <i>Dr Francesca Vianello (Exscientia)</i>