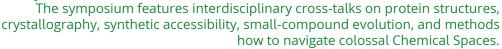


**DrugSpace Symposium** 

## Compounds, Targets, Spaces 21-22 April 2021

In the first virtual BioSolveIT symposium, experts discuss current developments and unprecedented innovations to tackle future challenges in drug discovery.



Time (CEST)	Wednesday 21 <sup>st</sup>	Thursday 22 <sup>nd</sup>
4:00 pm	Welcome address Christian Lemmen BioSolvelT	Welcome address Christian Lemmen BioSolvelT
4:05 – 4:45 pm	Barbara Zdrazil  EMBL-EBI/University of Vienna  "Trending Targets: Monitoring Targets in Drug Discovery through the Years"	Marcus Gastreich BioSolvelT "Fuzzy and Structure-Driven Navigation in Billion-Sized Haystacks"
4:45 – 5:30 pm	Yaroslav Bilokin Otava Chemicals "Human Protein Kinase CK2 Inhibitors: Discovering New Scaffolds"	Serghei Glinca CrystalsFirst "Cheaper, faster, smarter lead generation: Accelerating discovery of active compounds with protein-fragment complexes"
5:30 – 5:40 pm	Yurii Moroz Chemspace "Chemspace Platform: the Largest Catalog of Small Molecules to Search and Buy from!"	Hugues Lemoine Edelris "Navigating in Natural Product inspired Keymical Space with EDEN"
5:40 – 5:50 pm	Yan Xu Affiliation and title tba	<b>Debanu Das</b> Accelero Biostructures "Screening Molecular Compound Libraries Directly by Protein X-ray Crystallography"
5:50 – 6:35 pm	Matthias Rarey Center for Bioinformatics/University of Hamburg "Some New Algorithmic Approaches for Cheminformatics and Modeling"	Zhijie Liu Johnson & Johnson "Construct & Explore Virtual Libraries to Support a Janssen Kinase Project"
6:35 – 7:20 pm	Uta Lessel Boehringer Ingelheim "How to Compare Ultra-Large Chemical Spaces?"	Friedrich Rippmann <i>Merck KGaA</i> "Ultra-Large Tailor-Made Chemical Spaces"
7:20 pm	reception and networking  Details will be announced soon	

## **Our Partners**









