





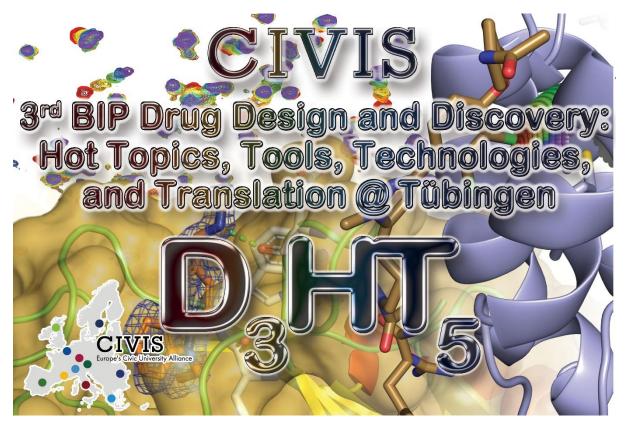






Virtual component of the CIVIS Blended Intensive Program: D₃HT₅

3rd BIP Drug Design and Discovery: Hot Topics, Tools, Technologies, and Translation @ Tübingen 2024



Planned Program (as of 19th Oct. 2023)

		Time slots (CEST)	Lectures, Speakers
Week 1	Tuesday, 30 th April	14:00-17:00	From Charges to π -systems: Classical Molecular Interactions – Familiar, yet still Surprising Non-Classical Molecular Interactions: New Tools, New Rules
			Prof. Dr. Frank M. Böckler Molecular Design & Pharm. Biophysics, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
	Friday, 3 rd May	14:00-17:00	The Drug Discovery Process & Technologies - Strategies, Issues, Costs, Outlook
			Thermodynamics in Drug Discovery (in Hit2Lead / LeadOpt)
			Molecular Complexity & Fragment-based Drug Discovery
			Hit-to-Lead optimization, Co-optimizing Ligand Properties, Avoiding late Attrition
			Prof. Dr. Frank M. Böckler Lab for Molecular Design & Pharm. Biophysics, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen















		Time slots (CEST)	Lectures, Speakers
Week 2	Wednesday, 8 th May	14:00-15:30	In Vitro Cell Culture Models and Evaluation of Drug Effects Dr. Miruna Stan, Faculty of Biology, Bucharest University
		15:30-17:00	Neurodegenerative Diseases and Drug Discovery Prof. Dr. Maria Cano-Abad , Facultad de Medicina, Universita Autonoma Madrid
	Friday, 10 th May	14:00-15:00	Immunomodulation and Drug Impact on Cytokine Production and on the M1/M2 Macrophage Phenotype Dr. Gratiela Gradisteanu, Research Institute of Bucharest University
		15:00-17:00	Evaluation Drug-induced Responses in Cultured Cells at Specific Time Points (Cell Cycle Analysis and Apoptosis). Screening of Drug-induced Reactive Oxygen Species Production in Leukocytes Subpopulations in Whole Blood (Ex-vivo) Prof. Dr. Luminita Marutescu, Faculty of Biology, Bucharest University
Week 3	Wednesday, 15 th May	14:00-17:00	Microcalorimetry (ITC, DSC) and NanoDSF: Application to Anticancer Drugs Screening and Thermodynamic Characterization Prof. Dr. Francois Devred , Faculty of Pharmacy, Aix-Marseille Université
	Friday, 17 th May	14:00-17:00	Introduction to NMR Spectroscopy and Related Screening Methods *Prof. Dr. Emmanuel Mikros, Faculty of Pharmacy, National Kapodistrian University of Athens
Week 4	Wednesday, 22 th May	14:00-17:00	Pharmacokinetics in Drug Discovery and Development: Physiologically-based PK (PBPK) Models Prof. Dr. Florence Gattacceca, Faculty of Pharmacy, Aix-Marseille Université
	Friday, 24 th May	14:00-17:00	Facts and Principles in Drug Discovery and Development: the Molecular Drug Space Selection and Optimization Prof. Dr. Sandrine Alibert, Faculty of Pharmacy, Aix-Marseille Université
Week 5	Wednesday, 29 th May	14:00-15:00	Animal Models (Ethical Considerations, Experimental Design, <i>In vivo</i> Analysis of Drug Effects) Dr. Gratiela Gradisteanu, Research Institute of Bucharest University
		14:00-17:00	Visualizing K ⁺ Ion Signaling Pathways in Cell Metabolism Prof. Dr. Robert Lukowski , Experimental Pharmacology, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
	Friday, 31 st May	14:00-15:30	Nanobody Technology Platform Prof. Dr. Ulrich Rothbauer , Pharm. Biotechnology, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
		15:30-17:00	Mass Spectrometry and Bioanalytics Prof. Dr. Michael Lämmerhofer , Pharm. Analytics, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen







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		Time slots (CEST)	Lectures, Speakers
Week 6	Wednesday, 5 th June	14:00-15:30	Peptides by Design – RiPP Engineering Technology JProf. Dr. Silja Mordhorst, Pharm. Biology, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
		15:30-17:00	Genome Mining Enables a Renaissance of Natural Product Drug Discovery Prof. Dr. Harald Groß, Pharm. Biology, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
	Friday, 7 th June	14:00-15:30	Targeted Covalent Inhibitors in Drug Discovery JProf. Dr. Matthias Gehringer, Medicinal Chemistry, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
		15:30-17:00	Academic Drug Discovery: Fiction, Facts or Fantasy Prof. Dr. Stefan Laufer, Medicinal Chemistry, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
Week 7	Wednesday, 12 th June	14:00-15:30	Introduction to Epigenetics and Related Drugs Prof. Dr. Antonello Mai, Faculty of Pharmacy, La Sapienza University
		15:30-17:00	Introduction to Ligand-based Methods: Pharmacophore Modeling, (3D) QSAR Prof. Dr. Rino Ragno Faculty of Pharmacy, La Sapienza University, Rome
	Friday 14 th June	14:00-15:30	Introduction to Structure-based Methods in Drug Design: Molecular Docking Prof. Dr. Rino Ragno Faculty of Pharmacy, La Sapienza University, Rome
		15:30-17:00	Machine Learning and Artificial Intelligence in Drug Discovery Dr. Markus Zimmermann, Molecular Design & Pharm. Biophysics, Department of Pharmacy and Biochemistry, Eberhard Karls University Tübingen
Week 8	Wednesday, 19 th June	14:00-15:30	Case Studies in Industrial Anti-cancer Drug Discovery: KRAS G12C and HIF2α Dr. Rainer Wilcken, Director, Computational Chemistry / Computer-Aided Drug Design at Flare Therapeutics
		15:30-17:00	Fragment-based Approaches towards PROTAC Development and Free Energy Methods in Industry Dr. Matthias R. Bauer, Senior Expert II Data Science at Novartis Institutes for BioMedical Research (NIBR)





National and Kapodistrian

University of Athens











		Time slots (CEST)	Lectures, Speakers
	Friday 21 st June	14:00-15:30	Embedding of Nanoparticles in Solid (Orodispersible) Matrices Prof. Dr. Denise Steiner, Pharmaceutical Technology, Department of Pharmacy, Westfälische Wilhelms-Universität-Münster (previously: Eberhard Karls University Tübingen)
		15:30-17:00	Clinical Trials Prof. Dr. Evangelos Terpos , School of Medicine, National Kapodistrian University of Athens