

**Programme (draft) of the 13th Bologna's convention on Crystal Forms - CF@Bo n. 13
University of Bologna 7-9 September 2025**

	SUNDAY 7 September	Aula Prodi Complesso di San Giovanni in Monte University of Bologna	Abstract
11:30-14:00	REGISTRATION		
	Chairing Dario Braga	Opening of the convention	
14:00-14:30	Fabrizia Grepioni Department of Chemistry G. Ciamician, University of Bologna, Italy	Crystal Forms @ Bo	O1
14:30-15:00	Alfred Lee MSD Research Laboratories Rahway, New Jersey 07065, USA	Solid Form Discovery: Past, Present and Future.	O2
15:00-15:30	Philippe Fernandes Associate Director at Johnson and Johnson	Key Insights from Computational Tools in Solid-State Pharmaceutical Development	O3
15:30-16:00	Franziska Emmerling Federal Institute for Materials Research and Testing (BAM) Department Materials Chemistry, Berlin	Real-Time Monitoring and Temperature Control for Optimized Polymorph Engineering	O4
16:00-16:30	Ivo B. Rietveld SMS Laboratory University of Rouen Normandy	Preparation and stabilisation of metastable polymorphs for the improvement of API bioavailability.	O5
16:30-17:00	Coffee break		
	Chairing Teresa Duarte		
17:00-17:30	Gareth Williams Department of Chemistry, University College London, UK	Accessing new polymorphs via solid solutions	O6
17:30-18:00	Helen Wheatcroft APS Crystallisation Scientist, AstraZeneca,UK	Crystallisation and Particle Control of a Multi-component API Crystal Form: From Structure to Process	O7
18:00-18:30	Elena Simone Department of Applied Science and Technology (DISAT) Politecnico di Torino, Torino, Italy	Can crystal engineers make food? A few examples of crystallization strategies for the design of food formulations	O8
18:30-19:00	Gerard Coquerel Laboratoire Sciences et Méthodes Séparatives. University of Rouen, France	Interferences between Polymorphism and Solid Solutions	O9
19:00	Aperif in the courtyard	Complesso di San Giovanni in Monte University of Bologna	
	MONDAY 8	Hotel Aemilia	

	Chairing Gazhala Sadiq		
8:30-9:00	Jon Steed Department of Chemistry, Durham University, Durham, UK	Crystals as Intellectual Property	O10
9:00-9:30	Susan Reutzel-Edens SuRE Pharma Consulting, LLC	Turning polymorph challenges into patent opportunities	O11
9:30-10:00	Vania André IMS Researcher at CQE-IMS/IST	Unlocking New Antibiotic Forms: Crystal Engineering and Supramolecular Strategies for Polymorphs, Cocrystals, and Beyond	O12
10:00-10:30	Sarah (Sally) Price FRS, Department of Chemistry, University College London, UK	Pharmaceutical Digital Design: Can we go from Chemical Structure through Crystal Polymorph to Conceptual Crystallization Process?	O13
10:30-11:00	Coffee break		
	Chairing		
11:00-11:30	Marcus A. Neumann CEO Avant-garde Materials Simulation Deutschland GmbH, Merzhausen, Germany	A conceptual framework for the crystallizability of organic compounds	O14
11:30-12:00	Doris Braun Ass.-Prof. Priv.-Doz. Dr. Institute of Pharmacy, Christian Doppler Laboratory for Advanced Crystal Engineering Strategies in Drug Development, University of Innsbruck	Hybrid Approaches in Solid Form Design: Virtual Screening and Experimental Validation	O15
12:00-12:30	Joost van den Ende Roche Pharma Research and Early Development, Therapeutic Modalities, Basel, Switzerland	Machine Learning within CSP: from one crystal energy landscape to another	O16
12:30-13:00	Miglani Bhardwaj, Rajni Associate research Fellow, Pfizer	An integrated approach combining experimental and computational for solid form design and selection	O17
13:00-14:00	Lunch and poster session		
14:00-16:00	Poster session	POSTER	P
16:00-16:30	Coffee break		
	Chairing		
16:30-17:30	GOLD sponsor presentations	GOLD sponsor presentations	S
	Chairing		

17:30-18:00	Enrico Modena <i>Associate Director Science and Technology</i> Novartis	Solid state, Polymorphism and the pharmaceutical	O18
18:00-18:30	Delia A. Haynes Stellenbosch University	Co-crystals, salts and sublimation	O19
18:30-19:00	Thomas Rades and Inês Martins Department of Pharmacy Faculty of Health and Medical Sciences University of Copenhagen	Amorphous forms of drugs: from preparation to polyAmorphism	O20
19:00-19:30	Amy Woods-Ryan Durham University (PhD student) and GSK (Investigator)	HEPES of conformational, multi-zwitterionic polymorphs	O21
	<i>Free evening in Bologna</i>		
	TUESDAY 9 September	Hotel Aemilia	
	Chairing		
8:30-9:00	Anna Slater University of Liverpool, Department of Chemistry and Materials Innovation Factory	Continuous flow chemistry as a tool for crystallisation of porous organic materials	O22
9:00-9:30	Edyta Pindelska Department of Pharmaceutical Chemistry and Biomaterials, Medical University of Warsaw, Poland	Mechanisms of Cocrystal Formation and Coformer Exchange in Ethenamide Systems – From <i>In Situ</i> Studies to Pharmaceutical Applications	O23
9:30-10:00	Matteo Daldosso Chief Scientific and Innovation Officer at PolyCrystalline	API: Don't Forget the I. From the Right Molecule to the Right Particle	O24
10:00-10:30	Federica Lazzari Solid state and crystallization scientist at Chemessentia Srl (part of Chemo group)	Challenges in crystallization scale-up of an API nanocrystalline form	O25
10:30-11:00	<i>Coffee break</i>		
	Chairing		
11:00-11:30	Joop ter Horst Tiofarma, NL	Optimizing complex multicomponent solid form discovery & crystallization process design	O26
11:30-12:00	Irene Bassanetti Senior Scientist in Analytics and Early Formulation Department, Preclinical R&D Chiesi Farmaceutici	The Critical Role of 3D Molecular and Biomolecular Structures in Innovative Drug Discovery	O27
12:00-12:30	Martin Viertelhaus Principal Scientist BASF SE - Analytical and Material Science	Solubility – Easy Parameter with Hurdles in Determination and Interpretation	O28
12:30-13:00	Matteo Lusi	Crystalline Solutions for Pharmaceutical Problems	O29

	Dpt of Chemical Science & Bernal Institute, University of Limerick, Ireland		
13:00-14:00	Lunch		
	Chairing		
14:00-14:30	Luc Aerts Head Solid State Development Sciences, UCB Pharma	Co-crystallisation as a versatile tool in pharmaceutical development	O30
14:30-15:00	Pietro Sacchi CCDC Research and Application Scientist – The Cambridge Crystallographic Data Centre	Computational approaches for the prediction of particle properties of organic molecular materials.	O31
15:00-15:30	Joe Lubach Distinguished Scientist at Genentech, Inc., department of Synthetic Molecule Pharmaceuticals.	Insights into Pharmaceutical Drug Substance and Product Using Multinuclear Solid-State NMR Spectroscopy	O32
15:30-16:00	Helen Blade Astrazeneca	To be confirmed	O33
16:00-16:30	Coffee break and Poster Prizes		
16:30-17:30	NICE event : Nature Inspired Crystal Engineering		
	Alessia Bacchi University of Parma	A NICE PoEM: cocrystals from liquid ingredients	O34
	Giuseppe Resnati Polytechnic of Milan	Pharmaceutical cocrystals via halogen bond	O35
	Michele Remo Chierotti University of Torino	Advanced Solid-State NMR tools for Crystal Engineering: From Structure Elucidation to Phase Purity Assessment	O36
	Pavel Zolotarev University of Milan	Study of pillared MOFs with Zn-paddlewheel state switching	O37

Registrations will open on March 1. Abstracts of lectures, posters and sponsor presentations will be available for download from MCE and PCL websites