DISCOVERY& DEVELOPMENT US 2025

October 2 - 3, 2025 | San Diego, CA

Conference Brochure

Equipping biopharma innovators with critical intelligence and partners needed to discover, develop and deliver first-in-class drugs rapidly



8 Content Tracks Across Two Days



200+
Pre-Arranged
Meetings



35+ Partners



1,000+ Attendees

Composed of 3 Focus Areas





Automation & Robotics



Formulation & Delivery

Meet Your KEYNOTE SPEAKERS



DAY ONE | 09:15

STEVE TURNER

Chief Executive Officer,
Pangio Biosystems &
Former Founding Chief
Executive Officer, PacBio

DAY TWO | 09:00 %

RAY DESHAIES

Former Senior Vice President of Global Research, Amgen



Book Now! Complimentary Industry, Academic or Investor Guest Passes Available

Register 💥

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Co-located with Biologics US 2025 > More Information



Discovery & Development US 2025

Welcome to Discovery & Development US 2025 – where we unify 3 stand-out focus areas under one roof – Drug Discovery, Formulation & Delivery and Automation & Robotics.

Join us for two days of high-level discussion & tailored content, spotlighting novel modalities & emerging computational tools to foster innovation. Our Drug Discovery program highlights the latest breakthroughs in target identification & validation, medicinal & Alguided computational chemistry & hit identification & screening. Join our exclusive track shining a light on degraders & molecular glues – where you'll hear case studies on the current technology & platforms for their discovery & design.

Our featured automation & robotics sessions features all things AI & data-powered, from streamlining lab workflows with automation and FAIR data management & digital transformation to implementing robotics within labs to improve productivity & efficiency.

Finally, join our Formulation & Delivery program to learn about the latest advancements in the formulation of large and small molecule-based drugs, developments in RNA, peptide and oligonucleotide design; and explore the innovations in the delivery of complex and challenging

therapeutics.

This year, Discovery & Development US goes even further - co-located with Biologics US 2025, we offer 1000+ attendees the unique opportunity to engage across both communities and forging the path to the next generation of therapies that will transform lives and reshape the future of healthcare.

Charlotte Catley

Portfolio Director – Discovery & Development Brand. Oxford Global



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WHAT'S NEW

Benefits Of Attending



Stay ahead of the curve: uncover how novel modalities like PROTACs, molecular glues, bispecific antibodies and beyond are reshaping the innovation landscape.



The latest in computational R&D: explore how AI-guided chemistry, advanced modeling and predictive tools are accelerating decision-making from discovery to human dose prediction.



Benchmark your lab against industry leaders: discover how automation & robotics can improve productivity for your workflows & exponentially decrease development time.



The use of Al – is it hype or is it justified? Make your decision by hearing from the experts using Al, data & digital transformation to drive drug discovery!



Discover the formulation breakthroughs redefining what's possible: from lipid and polymeric nanoparticles to automation technology and high-dosage therapies.



Keep your finger on the pulse of next-gen drug delivery:

hear how innovators are advancing RNA therapeutics, long-acting injectables, nanoparticle carriers, implants and inhalation approaches.

DON'T MISS THESE

Interactive Features

16:15 | Day 1
Panel Discussion:

The Status of AI/ML in Drug
Discovery: Where Are We and Where
Are We Going?

08:30 | Day 2

Roundtable Discussion:

IP Considerations in Early Discovery

12:10 | Day 2
Panel Discussion:

& Development

Sustainability Impact of Different Drug Delivery Methods

16:15 | Day 1
Industry Workshop:

The Changing Landscape of RNA
Design & Delivery

08:30 | Day 2
Roundtable Discussion:

From Drug Product Development to Overcoming Regulatory Hurdles

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At Oxford Global, our mission is to curate personalised experiences that foster community and inspire innovation.

We believe in the power of networking, connection, and knowledge to deliver quality products and services that exceed expectations. Partnering with Oxford Global means having a dedicated team committed to helping you achieve your goals and navigating the industry's ever-changing landscape.

Arrange 1-1 Meetings

Benefit from guaranteed one-to-one face time with your key prospects, with detailed pre-meeting information provided to enable effective and productive conversations.

Speaking Opportunities

Showcase your company's recent work to a relevant and highly engaged audience.

✓ Host Panel & Roundtable Discussions

Feature alongside key opinion leaders to discuss current hot topics and highlight your company's expertise.

Organise Workshops

Demonstrate best practice within the industry in front of your peers with case studies from your clients.

Exhibit your Products & Solutions

Promote your offerings and ensure delegates know where to find you with a prominent brand presence in the exhibition hall.

Digital Marketing & Lead Generation

Accessing the Oxford Global database, amplify your thought leadership and branding messaging through our digital content initiatives.





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Start Up Zone













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Big Attendance Meets Intimate Connections



Received the 2023 **American Chemical Society** (ACS) Philadelphia Section Award

'Innovation & Breakthroughs Under the Stars' Garden Party

Join us for our 'Innovations & Breakthroughs Under the Stars' Garden Party at the end of Day 1 at the Eventide Gardens, Sheraton San Diego by the Marina. This is your chance to mix and mingle with the brightest minds in the industry in a vibrant, open-air setting, complete with gourmet food stations, refreshing drinks, and a relaxed garden-party vibe.

Whether you're coming to connect, unwind, or simply enjoy a summer evening, this is the can't - miss celebration of the season, all under the



Day 1 Opening Keynote Address: The Single-Molecule Revolution Makes Landfall In Drug Discovery: Where We Are Now & What's To Come

- STEVE TURNER, Chief Executive Officer, Pangio Biosystems & Former Founding Chief Executive Officer, PacBio
- Don't miss Steve opening the event with his talk at 9:15 on Day 1.

Day 2 Opening Keynote Address: How 'Proximity **Medicines' Are Transforming Pharmacotherapy**

- RAY DESHAIES, Former Senior Vice President of Global Research, Amgen
- Ray's talk will be taking place at 9:00 on Day 2.

Interactive Sessions

Get up close and personal with industry trailblazers in our variety of roundtables and panel discussions. Hear firsthand how leaders are tackling big challenges in implementing novel technologies such as AI/ML, driving innovation, and shaping the future of drug development. Expect meaningful debates, bold ideas, and valuable insights you won't find in a typical presentation

The Start Up Zone Program

Step into the spotlight at Discovery & Development US 2025's Start-Up Zone – a dynamic launchpad for 10+ cutting-edge start-ups in drug discovery and development. Pitch your vision, showcase your tech, and connect with key players driving innovation. From Al-driven discovery to lab automation and next-gen therapeutics, this is your stage to shine.

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Program Agenda at a Glance

Explore Discovery & Development US 2025: North America's leading event for drug discovery, preclinical development and formulation returns in October 2025, uniting experts in both small molecule and biologics drug development to deliver first-in-class drugs at pace.

		Day One		Day Two
Cross- Programme Features		Opening Keynote Address Interactive Networking Breaks: ster Presentations & 1-2-1 Meetings Networking Drinks Reception Garden Party	Inter	pening Keynote Address Breakfast Roundtables ractive Networking Breaks: Presentations & 1-2-1 Meetings
Drug Discovery	Track 1	Novel Target ID for Small & Large Molecules	Track 1	Part 1: DMPK, Pharmacology & Safety Part 2: Hit Identification, Optimization & Screening
Program	Track 2	Medicinal & AI-Guided Computational Chemistry	Track 2	Part 1: Implementing Lab Automation & Digitization Part 2: Degraders & Molecular Glues
Formulation &	Track 3	Small Molecule Formulation	Track 3	Large Molecule Formulation
Delivery Program	Track 4	Part 1: Large Molecule Delivery & RNA Design Part 2: Large Molecule Formulation & RNA Design	Track 4	Part 1: Small Molecule Delivery Part 2: Large Molecule & RNA Delivery
Start Up Zone Afternoon: Start-Up Zone Presentations				

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AGENDA AT A GLANCE

Discovery & Automate US

Join leading scientific experts through our annual Discovery and Automation & Robotics programs, addressing critical strategies, scientific breakthroughs and the latest technological advancements in novel modalities & computational tools, new case studies on PROTACs & molecular glues and the utilization of AI, robotics & automation in early R&D.

VPs, Directors & Senior Managers from leading pharma & biotech companies and research institutions in the following fields and more:

- PROTACs
- Medicinal Chemistry
- Lead Generation
- Screening
- Small Molecule Discovery

- · AI/ML
- Lab Automation
- Digital Transformation
- DMPK
- Molecular Glues

Formal and informal meeting opportunities offer delegates the chance to discuss key solutions with leading service providers. Formal 1:1 meeting opportunities will be available to arrange prior to the event which take place during the dedicated networking breaks covering:

- Library Optimisation
- Screening Technologies
- Target Validation
- Organoid Discovery
- Protein Degradation Tools

- Automation
- Computational Tools
- Animal Models
- Inhibitor Discovery
- AI/ML

Day One

Day 1 Track 1: Novel Target ID for Small & **Large Molecules**

- Technological advancements for the discovery of novel new targets
- New paradigms beyond the rule of five
- Al to unblock drug discovery, drug repurposing and target validation
- Complementary use of genomic & proteomic technologies with Al
- 3D organoids to pursue 'undruggable' targets
- How can we streamline translation of basic scientific findings into human applications
- Disease-centric biologic insights and dealing with increasingly complex target biology
- New modalities that go after the target identification phase: ADCs, cell therapies, bispecific antibodies
- Validating novel biomarkers to improve success in early drug development

Track 2: Medicinal & Al-Guided Computational Chemistry

- New approaches in computer-aided drug design
- Quantum mechanics in drug discovery
- · Fragment & structure-based drug design, including Cryo-EM applications
- Biocatalysis for pharmaceutical synthesis & drug discovery
- · Utilising high-quality chemical probes
- Progressing 'intractable targets' from a medicinal chemistry & DMPK perspective

Day Two

Track 1:

Part 1: DMPK, Pharmacology & Safety

- Predicting & Measuring PK/PD, including computational models
- Novel models for ADME-Tox research & disease modelling - including organ-on-a-chip, organoids, iPSC-derived cells and humanized mouse models
- Development of animal efficacy models
- Safety assessment & safety pharmacology
- Human dose prediction

Part 2: Hit Identification, Optimization & Screening

- Phenotypic screening
- Affinity/binding assays for hit ID, including mass spectrometry, biophysical approaches, DNA-encoded libraries, high content imaging &

Track 2:

Part 1: Implementing Lab Automation & Digitization

- Automating laboratory workflows through discovery & development to improve productivity
- Lab robotics & tools for drug discovery

Part 2: Degraders & Molecular Glues

- using:
- » Proteomics & mass spec
- » Biophysical techniques
- New technologies & platforms for glue discovery
- PROTACs & targeted protein degradation
- Exploring new degradation pathways & modalities

- Molecular target-based screening
- computational screening

Digital transformation & digitization strategies

- · Discovery & optimization of novel degraders
- » Screening technologies

- Rational design & mechanistic understanding of molecular glues

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AGENDA AT A GLANCE

Formulation & Delivery

Join world-leading experts in drug product development and delivery for this in-person meeting on the West Coast's scientific hub, San Diego. You'll learn about the latest advancements in the formulation of large and small molecule-based drugs, developments in RNA and explore the innovations in the delivery of complex and challenging therapeutics.

VPs, Directors & Senior Managers from leading pharma & biotech companies and research institutions in the following fields and more:

- Formulation Science
- Drug Delivery
- Bioanalysis
- Device Development
- RNA Delivery

- Pharmaceutics
- Pharmaceutical Sciences
- Nanomedicine
- Drug Formulation
- Sustainable Formulation

Formal and informal meeting opportunities offer delegates the chance to discuss key solutions with leading service providers. Formal 1:1 meeting opportunities will be available to arrange prior to the event which take place during the dedicated networking breaks covering:

- Contract Development
- Characterization Tools
- Raw Materials
- Analytical Services
- Solubility & Stability Testing

- Dosage Forms
- Delivery Devices
- Formulation Development
- Quality Control
- Controlled Release

Day One

Track 3: Small Molecule Formulation

- Overcoming challenges in small molecule drug formulation and delivery including non-oral routes
- Recent advances to improve bioavailability: permeability, solubility
- Digitalization and automation of processes: AI/ML, in-silico modelling

Track 4: Part 1: Large Molecule Delivery & RNA Design

- Moving beyond injectable delivery of biologics including inhalation
- Overcoming oral delivery challenges
- Innovative technologies developing personalized drug delivery systems
- Novel strategies for RNA design
- Harnessing inherent potency of peptides with long-acting injectables and depositions

Track 4 Part 2: Large Molecule Formulation & RNA Design

- Overcoming challenges in RNA delivery
- Lipid nanoparticles to deliver RNA

Day Two

Track 3: Large Molecule Formulation

- Innovative formulation strategies for ADCs, cell & gene therapies, peptides and more
- Overcoming challenges in large molecule drug formulation
- Formulating therapies for pre-filled syringe delivery
- Optimizing high dosage biologics

Track 4: Part 1: Small Molecule **Delivery**

- Oral dosage form improvement
- Novel routes of administration for small molecules: pulmonary, transdermal, topical, ocular, nasal
- Patient-friendly delivery and dosage forms
- Challenges of delivering sustainability

Track 4: Part 2: Large Molecule & **RNA Delivery**

- Lipid nanoparticles to deliver biologics
- Pre-filled syringes
- · Use cases of utilizing viral and nonviral delivery modalities
- Overcoming challenges in RNA delivery
- Advanced techniques for delivery beyond the liver

Attending

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AWARD WINNING SPEAKERS ON BOARD

Shaping Drug Discovery Now

DR. RAY DESHAIES

Former Senior Vice President of Global Research, Amgen

Dr. Deshaies received the 25th Jacob and Louise Gabbay Award in Biotechnology and Medicine for his pioneering work on PROTACs (proteolysis targeting chimeras), a novel approach in targeted protein degradation.

TOR. RICK EWING

Dr. Beth Hoffman, Dr. Rick Ewing, Dr. Ray Deshaies, Dr. Robert Fremeau

Vice President, Head of Chemistry, Rapafusyn Pharmaceuticals

Dr. Ewing was honored with the 2023 American Chemical Society (ACS) Philadelphia Section Award, recognizing his distinguished career in medicinal chemistry.

TOR. ROBERT FREMEAU

Founder & Chief Executive Officer, BrainStorm Therapeutics

Dr. Fremeau serves as the Principal Investigator for a National Science Foundation (NSF) Small Business Innovation Research (SBIR) Phase I award, focusing on developing a novel Parkinson's disease drug discovery platform.

TOTAL SETH HOFFMAN

President & Chief Executive Officer, Origami Therapeutics, Inc.

Earlier in her career, Dr. Hoffman was awarded the Anna-Monika Foundation Prize for Advancement in Depression Research and an NSF Pre-doctoral Fellowship.

> Click Below to Preview our Award Winning Speaker's **Presentations**

> > Full **Programme:** Day Two

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FEATURED SPEAKERS

Shaping Drug Development Now

ERIC MUNSON,

Professor, Purdue University

IPEC Americas Foundation: Ralph Shangraw Memorial Award 2020

ANDREW GEALL,

Co-Founder & Chief Development Officer, Replicate Bioscience

BINDHU RAYAPROLU,

Director, Research & Development, Regeneron

MARIA WINZELL,

Senior Director Drug Development, Camurus

DIMITRIOS ZISOULIS,

Head of RNA Therapeutics,
Janssen Pharmaceutical Companies of Johnson &
Johnson



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SPEAKERS ACROSS ALL PROGRAMS

Discovery & Development US

STEVE TURNER

Chief Executive Officer, Pangio Biosystems & Former Founding Chief Executive Officer, PacBio

RAY DESHAIES

Former Senior Vice President of Global Research, Amgen

SHAH NAWAZ

Former Vice President, Head of Technology & Digital Transformation, Regeneron

ROBERT FREMEAU

Founder & Chief Executive Officer, BrainStorm Therapeutics

BETH HOFFMAN

President & Chief Executive Officer, Origami Therapeutics, Inc.

SAM SINAI

Co-Founder, Vice President, & Head of Machine Learning, Dyno Therapeutics

RICK EWING

Vice President, Head of Chemistry, Rapafusyn Pharmaceuticals

STEFANO ZANOTTI

Senior Vice President, Neuromuscular Research, Dyne Therapeutics

DAVID BUCKLER

Executive Director – Research Program Management, Regeneron

PRASHI JAIN

Director, Drug Discovery, Iterion Therapeutics

KAUSHEEK NANDY

Digital Transformation Director – Research, Boehringer Ingelheim

SHANTHI NAGARAJAN

Director, Eli Lilly

JOSH BAUGHMAN

Scientific Director, Bristol Myers Squibb

LEI JIA

Associate Director, Drug Discovery Data Science, Johnson & Johnson

SURENDRA NAYAK

Scientific Associate Director, Bristol Myers Squibb

ABHINAV SHARMA

Translational Human Models Lead, Dermatology/Immunology Discovery, AbbVie

FABIEN VINCENT

Consultant, Former Pharmacology Lab Head, Pfizer

PHIL COX

Research Fellow, Discovery Research, AbbVie

JEAN GE

Patent Attorney, Wolf Greenfield

SHUANGWEI LI

Director of Toxicology, Calibr-Skaggs Institute for Innovative Medicines

KATHERIN PATSCH

Director, Drug Discovery, Ellison Medical Institute

TOBIAS RAABE

Director, Translational Genome Editing Laboratory, Perelman School of Medicine, University of Pennsylvania

JOSEPH WU

Professor & Director, Stanford Cardiovascular Institute & Co-Founder, Greenstone Biosciences

NAGENDRA CHEMUTURI

Senior Director, ADME, Eli Lilly

YUE WEBSTER

Vice President, Model Driven Drug Discovery Platforms, Eli Lilly

PEGGY THOMPSON

Vice President, Head of Biology, Plexium

KISHORE PASIKANTI

Senior Director, DMPK, Grünenthal Group

DAVID LOYND

Chief Executive Officer, EnduRx

PETER HEIFETZ

Chief Executive Officer, Orpro Therapeutics

ANDREW GEALL

Co-Founder & Chief Development Officer, Replicate Bioscience

CRAIG BEHNKE

Executive Vice President, Production & Development, Lumen Bioscience

SEAN SULLIVAN

Vice President of Technology Innovation & Development, Arcturus Therapeutics

MARIA WINZELL

Senior Director Drug Development, Camurus

BINDHU RAYAPROLU

Director, Research & Development, Regeneron

JAMIE TSUNG

Director, Alnylam Pharmaceuticals

STEPHEN FRAMIL

Corporate Global Head of Accessibility, Merck

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SPEAKERS ACROSS ALL PROGRAMS

Discovery & Development US

DIMITRIOS ZISOULIS

Head of RNA Therapeutics, Janssen Pharmaceutical Companies of Johnson & Johnson

SUNITA YADAV

Associate Director, Novartis

SAYONI RAY

Principal Scientist, Janssen Pharmaceutical Companies of Johnson & Johnson

JITENDRA KUMAR

Lead Scientist, Chemistry & Process Development, Entos Pharmaceuticals

ERIC MUNSON

Professor, Purdue University

ENRICO ZAMBELLI

Sustainable Formulations Unit Manager, Chiesi

ELENA LIANG

Principal Research Scientist, AbbVie

ASHLEY DAVALOS

Assistant Director, Ionis

KALPITA MEHTA

Principal Combination Production & Medical Device Expert, Hikma

IACK ROGERS

Director of Neurochemistry, Harvard University

CATHERINE CAHILL

Assistant Professor, Harvard Medical School

SEVERIN SCHNEEBELI

Associate Professor, Purdue University

ANISHA DSOUZA

Research Fellow, Massachusetts Eye and Ear Infirmary, Harvard Medical School

SHILPA THOSAR

Former Executive Director, Cara Therapeutics

RAMESH KASHI

Former Senior Scientific Director, Bristol Myers Squibb

MARK ECCLESTON

Chief Executive Officer, ValiRx

VIVEK GUPTA

Associate Professor, St. John's University

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Day One | October 2 2025

Welcome & Keynote Address Coral 5

Oxford Global's Welcome Address By Eszter Sutowski Nagy, Senior Production Director

09:00

OPENING KEYNOTE ADDRESS & Q&A IN CORAL 5

The Single-Molecule Revolution Makes Landfall In Drug Discovery: Where We Are Now & What's To Come

STEVE TURNER, Chief Executive Officer, Pangio Biosystems & Former Founding Chief Executive Officer, PacBio



	Q&A session & transition time between conference rooms					
	DRUG DISCOV	ERY PROGRAM	FORMULATION & D	ELIVERY PROGRAM		
	NOVEL TARGET ID FOR SMALL & LARGE MOLECULES	MEDICINAL & AI-GUIDED COMPUTATIONAL CHEMISTRY	SMALL MOLECULE FORMULATION	LARGE MOLECULE DELIVERY & RNA DESIGN		
	CORAL 1	CORAL 2	SEAGLASS	CORAL 3		
	Morning Track Chair: ABHINAV SHARMA, Translational Human Models Lead, Dermatology/Immunology Discovery, AbbVie	Morning Track Chair TBA	Morning Track Chair TBA	Morning Track Chair: ERIC MUNSON, Professor, Purdue University		
	Track Keynote Address: Utilizing Human Genetics Technology For Antibody Discovery	Track Keynote Address: Targeting TBL1: Molecular & Computational Strategies To Modulate Complex Protein Interactions In Wnt Signaling	Track Keynote Address: Strategies For RNA Therapy For AD/PD With Implications For Treating SARS-CoV-2	Track Keynote Address: Navigating the Next Frontier: Industry Insights Into Emerging Trends In Biologics Delivery		
09:35		 This talk highlights mechanistic studies describing the inhibition of TBL1 by tegavivint and the development of second-generation small-molecule inhibitors using structural biology and predictive computational approaches Tegavivint is a small-molecule inhibitor of TBL1 that is currently in clinical trials for multiple oncology indications. Using in silico predictive modeling, we have identified a series of novel small molecules that bind to TBL1 and inhibit Wnt signaling. We will present novel strategies employed to identify these compounds and discuss the structural studies that guided lead optimization against this complex protein target 	We reported that the broad-acting Alzheimer's disease clinical drug Posiphen inhibited translation of the amyloid precursor protein as well as the mRNA encoding alpha-synuclein, the major culprit protein of Parkinson's disease & the Huntingtin protein. However, our new pipeline of translation blockers seeks to prevent neurodegeneration with potent drugs that interfere at higher selectivity with these 5'untranslated regions of the neurodegenerative disease-associated transcripts, i.e. the APP/ amyloid of AD and PD specific asyn. Recently, we discovered a line of therapy to inhibit APP translation linked to providing RNA-directed inhibitory actions to limit COVID by lowering the translation rates of the replicase in SARS-COV-2.	Biologics, which include complex molecules like proteins, monoclonal antibodies, and gene therapies, are revolutionizing modern medicine. As the industry continues to innovate, biologics delivery systems are evolving to meet the growing demand for precision, efficiency, and patient-centric care. This presentation explores the emerging trends shaping the future of biologics delivery, focusing on cutting-edge advancements, challenges, and opportunities.		
	DAVID BUCKLER, Executive Director – Research Program Management, Regeneron	PRASHI JAIN, Director, Drug Discovery, Iterion Therapeutics	CATHERINE CAHILL, Assistant Professor, & JACK ROGERS, Director of Neurochemistry, Harvard University	BINDHU RAYAPROLU, Director, Research & Development, Regeneron		
10:00		MORNING BREAK	& REFRESHMENTS 💝			
- 11:00			Poster D	Displays Page 1		

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	Day One October 2 2025				
	DRUG DISCOVERY PROGRAM		FORMULATION & DELIVERY PROGRAM		
	NOVEL TARGET ID FOR SMALL & LARGE MOLECULES	MEDICINAL & AI-GUIDED COMPUTATIONAL CHEMISTRY	SMALL MOLECULE FORMULATION	LARGE MOLECULE DELIVERY & RNA DESIGN	
	CORAL 1	CORAL 2	SEAGLASS	CORAL 3	
	Solution Provider Presentation	Solution Provider Presentation	Solution Provider Presentation	Solution Provider Presentation	
11:00	For sponsorship opportunities please contact sponsorship@oxfordglobal.com	Xtal≥i	ALMAC	P Pfanstiehl	
		MICHAEL BELLUCCI, Senior Director of R&D, Xtalpi	Senior Representative, Almac	Senior Representative, Pfanstiehl	
		Q&A session & transition tin	ne between conference rooms		
	Al-Powered Human Brain Organoid Platform To Transform CNS Drug Discovery	Chemical Space Generation, Exploration, And Visualization	How Digital Accessibility Is Critical For Digital Health Solutions	Polymeric Carriers Enabling Multi- Route Non-Invasive Delivery Of Large Therapeutic Molecules	
11:25	 Pioneering a new paradigm in CNS drug discovery Integrating patient-derived brain organoids with Al to accelerate therapeutic discovery Building Al foundation models to map disease-relevant gene networks Deploying biomarker-based screening to identify efficacious compounds in human brain disease models at the outset of drug discovery Delivering a scalable platform with broad applicability across complex neurodevelopmental and neurodegenerative disorders 	The number of theoretically possible drug-like molecules is vast and estimated to be >1060. Currently it is not practically feasible to enumerate such vast numbers, however it is possible to use abbreviated methods that encode the entirety of space using reaction SMARTS and associated monomer building blocks. This talk will focus on how we used such a method to generate and explore Nebula, AbbVie's UL chemical space.	 Personalized drug delivery systems Digital accessibility Formulation considerations for digital health soultions 	 Understand the key challenges and physiological barriers associated with non-invasive delivery of large molecule therapeutics Explore the design principles and material considerations behind engineering multifunctional carriers Evaluate preclinical data demonstrating the pharmacokinetic, biodistribution, and therapeutic performance of these engineered carriers 	
	ROBERT FREMEAU, Founder & Chief Executive Officer, BrainStorm Therapeutics	PHIL COX, Research Fellow, Discovery Research, AbbVie	STEPHEN FRAMIL, Corporate Global Head of Accessibility, Merck	VIVEK GUPTA, Associate Professor, St. John's University	
		Q&A session & transition tin	ne between conference rooms		
	Gold Solution Provider Presentation	Solution Provider Presentation	Solution Provider Presentation	Solution Provider Presentation	
1:50	For sponsorship opportunities please contact sponsorship@oxfordglobal.com	© Chemaxon	APRECIA	⊗ NOF	
		Senior Representative, Chemaxon	Senior Representative, Aprecia	Senior Representative, Nof Corporation	
2:15		ਊ LUNCH	BREAK		
- 3:35	1-2-1 Me	etings x4	Poster	Displays 💾	

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	Day One October 2 2025				
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	NOVEL TARGET ID FOR SMALL & LARGE MOLECULES CORAL 1	MEDICINAL & AI-GUIDED COMPUTATIONAL CHEMISTRY CORAL 2	SMALL MOLECULE FORMULATION SEAGLASS	LARGE MOLECULE DELIVERY & RNA DESIGN CORAL 3	
	Afternoon Track Chair: TOBIAS RAABE, Director, University of Pennsylvania	Afternoon Track Chair TBA	Afternoon Track Chair TBA	Afternoon Track Chair: ERIC MUNSON, Professor, Purdue University	
	Gold Solution Provider Presentation	Solution Provider Presentation	Solution Provider Presentation	Small Particles, Big Impact: Maximizing Therapeutic Impact Through Nanoparticle-Enabled Delivery	
13:35	For sponsorship opportunities please contact sponsorship@oxfordglobal.com	optibrium	XtalPi	Learn how key drug delivery challenges such as targeting, bioavailability, and circulation time can be overcome through nanoparticle-enabled formulations.	
				FORTIS LIFE SCIENCES	
		Senior Representative, Optibrium	Senior Representative, Xtalpi	JESS WIDMAN, Senior Product Development Scientist, Fortis Life Sciences	

Q&A session & transition time between conference rooms

	The FORCETM Platform Delivers Functional Improvement In DM1 & DMD	Utilizing AI/ML Models To Optimize Screening Techniques	Formulation Of CLX-001, A Next Generation Oncolytic Peptide For Cancer Therapy	FAST Protein-Enabled Proteolipid Vehicles: A Novel Platform For Advanced Gene Therapy Delivery
14:00			 CLX001 – a novel oncolytic peptide Membrane-lytic peptide based intravenous nanotherapeutics Oncolytic peptides – synergistic agents for small molecule and Immunotherapy 	 FAST protein integration into proteolipid vehicles (PLVs) enables efficient delivery of therapeutic DNA and RNA directly into target cell cytoplasm by overcoming endocytosis barriers FAST protein-enabled PLVs demonstrate broad biodistribution and low immunogenicity in animal models, maintaining sustained gene expression for extended periods without significant inflammatory responses The PLV platform, leveraging FAST proteins, offers a durable and versatile system capable of delivering large genetic payloads, including CRISPR-Cas9, with significant implications for future gene therapies
	STEFANO ZANOTTI, Senior Vice President, Neuromuscular Research, Dyne Therapeutics	SAM SINAI, Co-Founder, Vice President & Head of Machine Learning, Dyno Therapeutics	MARK ECCLESTON, Chief Executive Officer, ValiRx	JITENDRA KUMAR, Lead Scientist, Chemistry & Process Development,, Entos Pharmaceuticals

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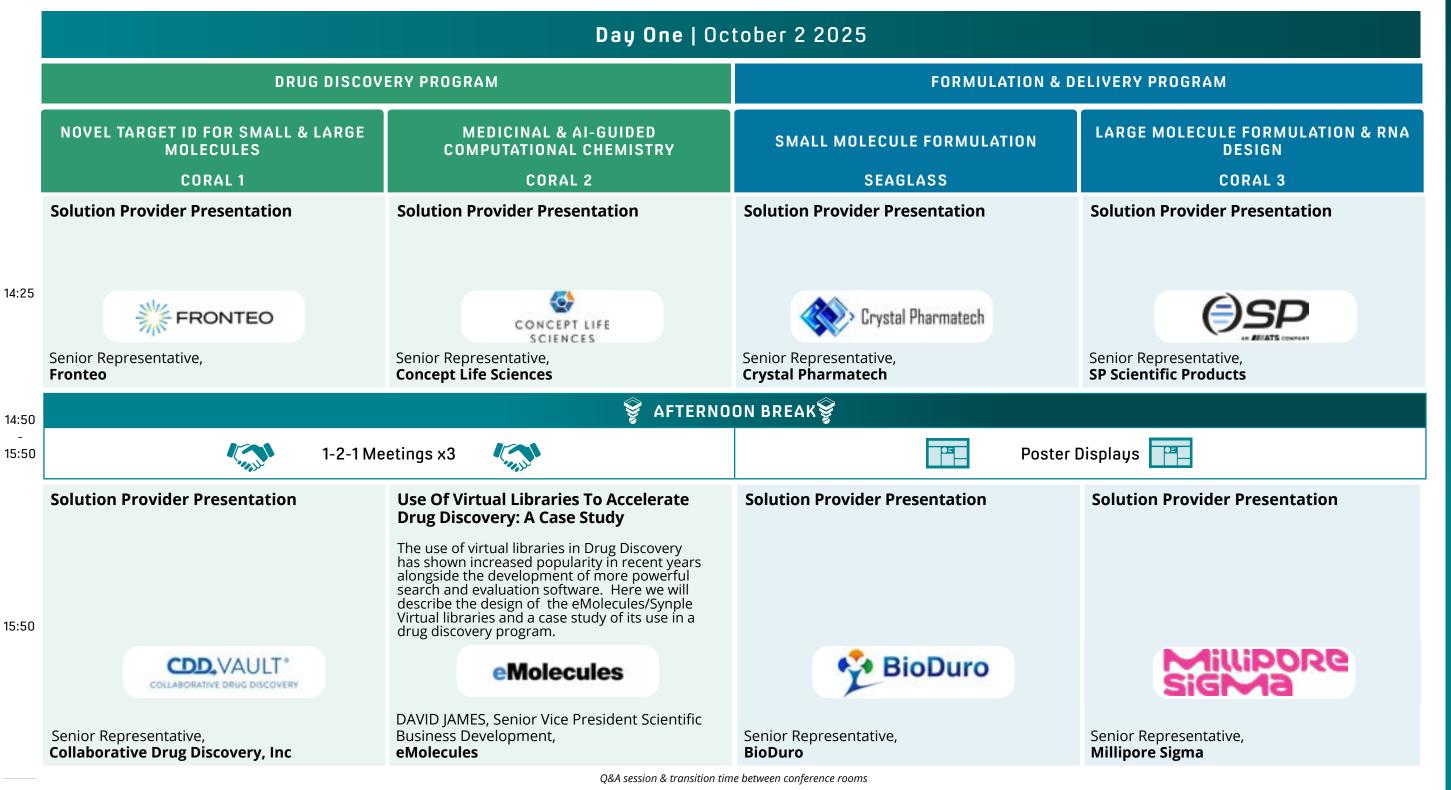
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NOVEL TARGET ID FOR SMALL & LARGE MOLECULES	START UP ZONE	SMALL MOLECULE FORMULATION	LARGE MOLECULE FORMULATION & RNA DESIGN
CORAL 1	CORAL 2	SEAGLASS	CORAL 3
Complex In Vitro Models For Drug Discovery & Development Of Immunological Disorders ABHINAV SHARMA, Translational Human Models Lead, Dermatology/Immunology Discovery,	PANEL DISCUSSION: The Status Of Al/ML In Drug Discovery: Where Are We And Where Are We Going? • Examples of successful application / acceleration of discovery programmes using Al/ML • Best practices • Limitations & challenges Panellists: PHIL COX, Research Fellow, Discovery Research, AbbVie MATTHEW SEGALL, Chief Executive Officer, Optibrium KATHERIN PATSCH, Director, Drug Discovery,	WORKSHOP: Tides Analytical Development: Innovations In Oligonucleotides & Peptide Analysis & Delivery Presentation 1: New Computational & Experimental Insight Into Oral Peptide Delivery With Permeation Enhancers Oral peptide formulations improve patient compliance by eliminating injections, especially in early-stage disease where long-term adherence is crucial. Permeation enhancers have transformed oral delivery by enabling peptides to cross the intestinal barrier safely, allowing effective systemic absorption. Despite these advances, further refinement and understanding of ophancer.	WORKSHOP: The Changing Landscapes Of RNA Design & Delivery Presentation 1: Translation Of An Optimized Self-Replicating RNA (srRNA) Platform With An Expanded Clinical Therapeutic Index • Demonstration that next generation srRNA vectors with optimized non-coding regions achieve protective immunity at doses up to 1,000,000-fold lower than conventional mRNA in preclinical models of influenza and rabies • Phase I clinical data with RBI-4000, an improved srRNA vector encoding the rabies virus glycoprotein and encapsulated in a novel lipid nanoparticle. Demonstrating the high therapeutic
Stem Cells, Genomics & Al For Drug Discovery	Ellison Medical Institute ime between conference rooms Start Up Zone Presentations x2	refinement and understanding of enhancer mechanisms are needed. This presentation offers new insights into their function, supported by constant pH molecular dynamics simulations and experimental data from relevant model systems. SEVERIN SCHNEEBELI, Associate Professor, Purdue University	index of next generation srRNA vectors and the feasibility of both low dose and single dose administration for vaccine applications • Potential of srRNA for applications beyond vaccines ANDREW GEALL, Co-Founder & Chief
• Learn recent announcements of 'New Alternative Methodologies' by FDA, NIH and EPA • Learn how stem cells, organoids, genomics & Al can be integrated for drug discovery		Presentation 2: Universal Assay Determination Method For Antisense Oligonucleotides Via Slope Spectroscopy Developed a universal assay method for antisense oligonucleotide (ASO) drug products using slope spectroscopy with the SoloVPE system, offering a simple, robust alternative for in-process control (IPC) of ASO concentration. The method was optimized across multiple parameters and validated for linearity, accuracy, precision, reproducibility, and robustness—with excellent performance across different ASO chemistries and formulations. The method proved transferable across sites (Ionis and CMOs), is independent of specific ASO attributes, and was recognized as a universal IPC assay platform for broad application in ASO drug product manufacturing.	Presentation 2: Lipid Nanoparticle (LNP) Formulation & mRNA Manufacturing Development • Development of an ionizable lipid library • Drug product formulation development • Optimization of drug substance manufacturing unit operations SEAN SULLIVAN, Vice President of Technology Innovation & Development, Arcturus Therapeutics
JOSEPH WU, Professor & Director, Stanford Cardiovascular Institute & Co-Founder, Greenstone Biosciences		ASHLEY DAVALOS, Assistant Director,	

16:15

16:40

Greenstone Biosciences

Q&A session & transition time between conference rooms

Ionis

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NOVEL TARGET ID FOR SMALL & LARGE MOLECULES CORAL 1

Lipid Nanoparticles Enable Efficient Cell Type Specific Gene Targeting In A Patient Liver - Derived Fibrosis Model

- •We developed a lipid nanoparticle-based platform that enables >80% gene targeting efficiency in non-dividing hepatocytes, cholangiocytes, myofibroblasts, and proinflammatory macrophages within a fully numan 3D organoid model in just two days. This system supports cell type-specific knockdown, overexpression, or editing for multi-target liver fibrosis therapies
- Our modular, patient-derived organoid model allows parallel screening across diverse fibrotic liver diseases, including MASH, cholangiopathies, alcohol-related cirrhosis, and viral fibrosis

TOBIAS RAABE, Director, Translational Genome Editing Laboratory, Perelman School of Medicine,

University of Pennsylvania

START UP ZONE CORAL 2

> **Start Up Zone Presentations** (10 min) Presentation 1: Si-Care Bio



(10 min) Presentation 2: Athens LifeTech Park



Q&A session & transition time between conference rooms

DRUG DISCOVERY PROGRAM

INTERACTIVE

PANEL DISCUSSION: Advanced Tools & Tech For Drug Discovery

- Making use of all the datasets
- New alternative methods

Panellists:

DAVID BUCKLER, Executive Director – Research Program Management, Regeneron

SAM SINAI, Co-Founder, Vie President, & Head of Machine Learning, Dyno Therapeutics

JOSEPH WU, Professor & Director, Stanford Cardiovascular Institute & Co-Founder, **Greenstone Biosciences**

ABHINAV SHARMA, Translational Human Models Lead, Dermatology/Immunology Discovery, **AbbVie**

(10 min) Presentation 1: Insilico Prediction Of In-Vivo Efficacy At The Drug Design Stage



(10 min) Presentation 2: Ontologic



INTERACTIVE **WORKSHOP (CONT.): Tides Analytical**

Development: Innovations In Oligonucleotides & Peptide Analysis & **Delivery**

Panel Discussion: Exploring Alternative Route Of Tides Delivery

SMALL MOLECULE FORMULATION

SEAGLASS

- Oral delivery of peptides
- Nasal delivery
- Delivering oligos beyond the liver
- Pre-filled syringes
- Sustainability and stability considerations

INTERACTIVE

FORMULATION & DELIVERY PROGRAM

WORKSHOP (CONT.): The Changing **Landscapes Of RNA Design & Delivery**

LARGE MOLECULE FORMULATION & RNA

DESIGN

CORAL 3

Panel Discussion: Novel Strategies For RNA Design & Delivery

- RNA Interference (RNAi)
- microRNA
- Antisense Oligonucleotides
- circRNA

Moderator: ERIC MUNSON, Professor, Purdue University

Panellists:

KALPITA MEHTA, Principal Combination Production & Medical Device Expert, Hikma

ASHLEY DAVALOS, Assistant Director, Ionis

SEVERIN SCHNEEBELI, Associate Professor, **Purdue University**

SAYONI RAY, Principal Scientist, Janssen Pharmaceutical Companies of Johnson & Johnson

Moderator: SUNITA YADAV, Associate

Director, Novartis

Panellists:

DIMITRIOS ZISOULIS, Head of RNA Therapeutics, Janssen Pharmaceutical Companies of Johnson & Johnson

ANDREW GEALL, Co-Founder & Chief Development Officer, Replicate Bioscience

SEAN SULLIVAN, Vice President of Technology Innovation & Development, Arcturus

Therapeutics

JAMIE TSUNG, Director, Alnylam **Pharmaceuticals**

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	INTERACTIVE	(10 min) Presentation 1: Epigenetic	INTERACTIVE	
	PANEL DISCUSSION: Advanced Tools & Tech For Drug Discovery	Therapies For Next-Generation Pain Therapeutics & Beyond	PANEL DISCUSSION: Digitization & Automation Of Formulation Processes	
	Making use of all the datasetsNew alternative methods	FERNANDO ALEMAN, Co-Founder & Chief Scientific Officer, Navega Therapeutics	Developments with AI within formulation Utilising the latest technology AI/ML for personalized formulation	
55	Panellists: DAVID BUCKLER, Executive Director – Research Program Management, Regeneron	Navega Thérapeutice		Attendees are welcome to join co-located sessions
	SAM SINAI, Co-Founder, Head of Machine Learning, Dyno Therapeutics		Moderator: SHILPA THOSAR, Former Executive	
	JOSEPH WU, Professor & Director, Stanford Cardiovascular Institute & Co-Founder, Greenstone Biosciences		Director, Cara Therapeutics Panellist: STEPHEN FRAMIL, Corporate Global Head of	
	ABHINAV SHARMA, Translational Human Models Lead, Dermatology/Immunology Discovery, AbbVie		Accessibility, Merck VIVEK GUPTA, Associate Professor, St. John's University	

Grab a drink & make your way to the Visionary Voices Stage for the end of day summaries

Section of the control of the contro

18:20

Innovations & Breakthroughs Under the Stars Garden Party at Eventide Gardens, Sheraton San Diego Welcome Address by Dr Rick Ewing, Vice President, Head of Chemistry, Rapafusyn Pharmaceuticals



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Day Two | October 3 2025



In The Lab

Morning Roundtables with Continental Breakfast In Coral Foyer 📸



ROUNDTABLE DISCUSSION 1:

INTERACTIVE

IP Considerations In Early Discovery & Development

- Establishing effective IP infrastructure
- Crafting a coherent IP strategy that aligns with early discovery and development activities
- Building a "diligence ready" IP portfolio

Moderator: JEAN GE, Patent Attorney, Wolf Greenfield

ROUNDTABLE DISCUSSION 2:

INTERACTIVE

Best Practice Data Management Strategies

- Strategies for data management in labs of the future
- Considerations & challenges

Moderator: YUE WEBSTER, Vice President, Model Driven Drug Discovery Platforms, Eli Lilly

ROUNDTABLE DISCUSSION 3:

Overcoming Regulatory Hurdles

INTERACTIVE From Drug Product Development To

- Regulatory lifecycle
- Aligning product development strategies with regulatory expectations

Moderator: SHILPA THOSAR Former Executive Director, **Cara Therapeutics**

ROUNDTABLE DISCUSSION 4:

INTERACTIVE

LNP Formulation Development Of RNA & DNA

- Tailoring LNPs for different administration routes
- Optimizing formulations

Moderator: SAYONI RAY, Principal Scientist, **Janssen Pharmaceutical Companies of** Johnson & Johnson

Q&A session & transition time between conference rooms



08:30

KEYNOTE ADDRESS IN CORAL 5: How 'Proximity Medicines' Are Transforming Pharmacotherapy

Multispecific medicines - both large and small molecule - are on the ascendancy. Half of the 72 FDA-approved multispecific drugs have gained approval since 2020, and from 2022-24, multispecifics accounted for 1/6th of FDA approvals. Among the different categories of multispecific drugs, proximity medicines are of particular interest. These medicines typically work by either localizing the therapeutic molecule or recruiting a regulatory mechanism in the body to modulate the target. I will discuss the advantages conferred by multispecificity with particular emphasis on proximity medicines, and will describe a category of proximity medicines – LOCKTACs – that work by stabilizing a pre-existing complex.

RAY DESHAIES, Former Senior Vice President of Global Research,

Amgen

Q&A session & transition time between conference rooms

	DRUG DISCOVERY PROGRAM	AUTOMATION & ROBOTICS PROGRAM	FORMULATION & DELIVERY PROGRAM	
	DMPK, PHARMACOLOGY & SAFETY	IMPLEMENTING LAB AUTOMATION & DIGITIZATION	LARGE MOLECULE FORMULATION	SMALL MOLECULE DELIVERY
	CORAL 1	CORAL 2	SEAGLASS	CORAL 3
	Track Chair: SHUANGWEI LI, Director of Toxicology, Calibr- Skaggs Institute for Innovative Medicines	Morning Track Chair TBA	Morning Track Chair TBA	Morning Track Chair TBA
	Solution Provider Presentation	Platinum/Gold Solution Provider Presentation	INTERACTIVE SPONSORED WORKSHOP:	Solution Provider Presentation
09:35	optibrium:	For sponsorship opportunities please contact sponsorship@oxfordglobal.com	Tackling CMC & Manufacturing Challenges Of Complex LNP-Based Biopharmaceuticals For Oncological, Gene Editing & Vaccine Applications	SEVER PHARMA SOLUTIONS
	Senior Representative, Optibrium		SVSTOSIUS	Senior Representative, Sever Pharma Solutions
	Q&A session & transition tim	e between conference rooms		Q&A session & transition time between conference rooms

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	DMPK, PHARMACOLOGY & SAFETY	IMPLEMENTING LAB AUTOMATION & DIGITIZATION	LARGE MOLECULE FORMULATION	SMALL MOLECULE DELIVERY	
	CORAL 1	CORAL 2	SEAGLASS	CORAL 3	
10:00	How Close Are We? The Evolving Science of Human PK Prediction This talk will explore the current landscape of human pharmacokinetic (PK) prediction, focusing on our ability to estimate clearance, volume of distribution, and oral bioavailability. We'll also examine a case study addressing non-linearity in PK and its impact on first-in-human dose selection—highlighting practical strategies, limitations, and evolving methodologies in human PK predictions.	Automated, Integrated Wet Lab & Digital Platform This talk outlines Boehringer Ingelheim's strategic initiatives in digital and lab automation in the large molecule research space. It highlights investments in automation systems, digital integration, and lab orchestration to enhance biotherapeutics research, improve efficiency, and drive innovation at the Biotherapeutics department. KAUSHEEK NANDY, Digital Transformation	SPONSORED WORKSHOP: Tackling CMC & Manufacturing Challenges Of Complex LNP-Based Biopharmaceuticals For Oncological, Gene Editing & Vaccine Applications	Formulation Considerations For Polymeric Implants	
	KISHORE PASIKANTI, Senior Director, DMPK, Grünenthal	Director – Research, Boehringer Ingelheim		ELENA LIANG, Principal Research Scientist, AbbVie	
	Q&A session & transition tim	ne between conference rooms		Q&A session & transition time between conference rooms	
10:25	Solution Provider Presentation is now part of Millipore Sigma	Solution Provider Presentation Brooks		Solution Provider Presentation Serán	
	Senior Representative, Hub Organoids	Senior Representative, Brooks Automation		Senior Representative, Seran	
10:50			& REFRESHMENTS 🟺		
- 11:50	1-2-1 Meetings x3 Poster Displays		Displays P		
11:45	Bronze Solution Provider Presentation For sponsorship opportunities please contact sponsorship@oxfordglobal.com	Bronze Solution Provider Presentation For sponsorship opportunities please contact sponsorship@oxfordglobal.com	Senior Representative, Unchained Labs	Senior Representative, Lubrizol	
11:45	For sponsorship opportunities please contact sponsorship@oxfordglobal.com	sponsorship@oxfordglobal.com	Senior Representative,		

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	CORAL 1	CORAL 2	SEAGLASS	CORAL 3	
	Toxicology In Small Molecule Drug Discovery: From Target Safety Assessment To IND	Bridging The Gap Between Automation, Digitalization, & Innovation	Transforming Short-Acting Peptides Into Long-Acting Using Liquid Lipid Depot Formulations	Sustainable Inhalation Devices - The Challenges Of Meeting Net Zero Global Emissions	
12:10	Overview of the critical role toxicology plays in small molecule drug discovery, from early target safety assessment through IND-enabling studies Key strategies for identifying, characterizing, and mitigating safety risks during candidate selection and lead optimization Practical guidance on designing and timing toxicology studies to align with regulatory expectations and support successful IND submissions.		 Presentation of the liquid lipid depot formulation technology – FluidCrystal® Small volume, ready-to-use formulations for improved compliance and better efficacy Examples of a variety of peptide formulations with extended half-lives with clinical data 	 The transition to low global warming potential propellants in metered dose inhalers is essential for reducing the environmental impact of pharmaceutical aerosols This presentation provides a technical overview of the challenges and strategies to address the change and significantly reducing the carbon foot print of MDIs 	
	SHUANGWEI LI, Director of Toxicology, Calibr-Skaggs Institute for Innovative Medicines	SHAH NAWAZ, Former Vice President, Head of Technology & Digital Transformation, Regeneron	MARIA WINZELL, Senior Director Drug Development, Camurus	ENRICO ZAMBELLI, Sustainable Formulation Project Unit Manager, Chiesi	
		Q&A session & transition ti	me between conference rooms		
	Bronze Solution Provider Presentation	Bronze Solution Provider Presentation	Bronze Solution Provider Presentation	Bronze Solution Provider Presentation	
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- 14:00	1-2-1 Mee	etings x3	Poster D	isplays P	

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	CORAL 1	CORAL 2	SEAGLASS	CORAL 3
	Afternoon Track Chair: KISHORE PASIKANTI, Senior Director, DMPK, Grünenthal	Afternoon Track Chair: DAVID BUCKLER, Executive Director, Program Management, Regeneron	Afternoon Track Chair TBA	Afternoon Track Chair TBA
	Novel Models For ADME-Tox Research	RapaGlue™: A Novel Non-Degrading Molecular Glue Modality For Unlocking The Hard To Drug Proteome	Solving Key Function & Formulation Challenges To Enable Development Of A Novel First In Class Thioredoxin-Based Therapeutic	PANEL DISCUSSION: Sustainability Impact Of Different Drug Delivery Methods
14:00		Rapafuysn's platform of non-degrading molecular glues is uniquely positioned to target intracellular proteins and the cytosolic side of transmembrane proteins unlocking the hard to drug proteome. The company's platform of RapaGlue's takes advantage of the cytosolic residing FKBP12 to form ternary complexes with disease target proteins. The presentation will describe successful hit campaign highlighting our progress in inhibiting the GTPase, GSTP1 a target for ferroptosis.	 Thioredoxin is a previously unexploited secreted stress-protectant protein family with anti-inflammatory, anti-oxidative & ferroptosis-inhibiting activity Thioredoxin protects both normal & cancer cells & has demonstrated poor pharmacology with rapid clearance & an association with chemoresistance during cancer treatment Thioredoxin ORP100S has been designed & formulated to greatly improve pharmacokinetics & increase potency, with no evidence of protective function in multiple cancer cell lines 	Inhalation Oral delivery Injection
	(Reserved) JIANQIN LU, Director, Pharmaceutics & Pharmacokinetics Track, University of Arizona	RICK EWING, Vice President, Head of Chemistry, Rapafusyn Pharmaceuticals	PETER HEIFETZ, Chief Executive Officer, Orpro Therapeutics	
		Q&A session & transition time between conference rooms		
	PANEL DISCUSSION: Successful Human Dose Prediction Across Modalities – Considerations & Challenges	Discovery And Characterization Of A Clinical-Stage Molecular Glue Degrader Of Wee1 Kinase, Bms-986463, For The Treatment Of Advanced Solid Tumors	Spirulina As A Next-Generation Biomanufacturing Platform: Low Cost Biologics At Scale	
	 Optimization of small molecule compounds Traditional vs novel modalities Ensuring quality PK predictions 	WEE1 is an essential cell cycle regulator kinase that prevents mitotic division of cells with DNA damage. Some tumors disproportionally depend	Lumen Bioscience has engineered spirulina into a platform for producing oral and intranasal biologic drugs. This talk will highlight the platform's	Moderator: SENIOR REPRESENTATIVE, Quantum Si
14:25	2.13dring quality 11x predictions	upon WEE1 for survival and WEE1-selective inhibitors have shown clinical promise. We describe the discovery and pharmacological characterization of BMS-986463, a first-in-class	molecular tools, manufacturing capabilities, and therapeutic applications, with examples spanning from C. difficile infection to metabolic disease.	Panellists: ENRICO ZAMBELLI, Sustainable Formulation Project Unit Manager, Chiesi
	Panellists: KISHORE PASIKANTI, Senior Director, DMPK,	molecular glue degrader of WEE1 currently in Phase I enrolling patients in multiple oncology indications.		JACK ROGERS, Director of Neurochemistry, Harvard University
	Grünenthal SENIOR REPRESENTATIVE, Chemaxon		CDAIC DELINIVE Francisco Vica David	CATHERINE CAHILL, Assistant Professor, Harvard Medical School
	NAGENDRA CHEMUTURI, Senior Director, ADME, Eli Lilly	JOSH BAUGHMAN, Scientific Director, Bristol Meyers Squibb	CRAIG BEHNKE, Executive Vice President, Production & Development, Lumen Bioscience	VIVEK GUPTA, Associate Professor, St. John's University

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CORAL 1	CORAL 2	SEAGLASS	CORAL 3		
Small Molecule Discovery: Unfolding Precision Medicines Using Phenotypic Screening In Primary Human Cells	Novel Ultra High Density Screening For Molecular Glues	PANEL DISCUSSION: Optimizing Formulation Of High Dosage Biologics	Toward Delivery Of Improved Efficacy, Greater Safety, & Enhanced Patient Experience During & After Treatment		
 Rationale for the use of phenotypic screening Considerations for establishing a high quality phenotypic screen Hit identification and de-convolution in human disease models 		 Viscosity, aggregation, and stability issues Strategies for formulation optimization Scale-up and regulatory alignment 	 Erkki Ruoslahti's "cellular zipcodes" and a novel delivery mechanism of action Improved efficacy: how? Greater safety: how? Enhanced patient experience: how? 		
		Moderator: RAMESH KASHI, Former Senior Scientific Director, Bristol Myers Squibb	Challenges to preclinical development of productFuture vision		
		Panellists:			
		SUNITA YADAV, Associate Director, Novartis			
BETH HOFFMAN, President & Chief Executive	PEGGY THOMPSON, Vice President, Head of	JAMIE TSUNG, Director, Alnylam Pharmaceuticals			
Officer, Origami Therapeutics, Inc.	Biology, Plexium	CRAIG BEHNKE, Executive Vice President, Production & Development, Lumen Bioscience	DAVID LOYND, Chief Executive Officer, EnduRx		

Q&A session & transition time between conference rooms

VS@Lilly: Enablement Of A Dedicated
Platform For Identifying Novel Small-
Molecule Binders

The VS@Lilly project represents a significant advancement in computational chemistry, focusing on the development of a dedicated platform for identifying novel small-molecule binders. This platform supports routine ultralarge-scale virtual screening, leveraging advanced technology to explore commercially available chemical spaces and identify primary or backup chemical starting points. It features an extensive database with over 40 million off-the-shelf compounds and several billion compounds available through on-demand synthesis, all in various formats for efficient screening.

SHANTHI NAGARAJAN, Director, **Eli Lilly**

Predictive Modeling Of Degrader's Oral Bioavailability & In Vivo PK Properties

- Oral bioavailability of TPDs, especially larger sized bifunctional molecules (i.e. PROTAC), is one of the most challenging properties to be optimized for oral drug discovery
- •A comprehensive in vivo data set were collected from public domain and 2D/3D descriptors were developed to establish predictive models for oral bioavailability
- We address the different behavior of predictive models by using data sets from different type of animal models
- Going beyond ML models, mechanistic based method was established to predict in vivo PK parameters

LEI JIA, Associate Director, Drug Discovery Data Science, **Johnson & Johnson**

Formulating Proteins For Maximum Stability

- •Solid-state characterization of mRNA LNPs to identify formulation and lyophilization conditions to ensure long-term stability
- Predicting protein stability, including mAbs, in the solid state – formulation approaches to ensure long-term stability
- New technologies for in situ and non-destructive characterization of aseptic solid formulations

ERIC MUNSON, Professor, **Purdue University**

Q&A session & transition time between conference rooms

BROCHURE CONTENTS

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Featured
Speakers:
Discovery

Featured
Speakers:
Formulation

Formulations For Pre-Filled Syring

Delivery

JAMIE TSUNG, Director,

Alnylam Pharmaceuticals

All Speakers

Full
Programme:
Day One

Full
Programme:
Day Two

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Day Two October 3 2025				
DRUG DISCOVERY PROGRAM		FORMULATION & DELIVERY PROGRAM		
SCREENING & HIT IDENTIFICATION CORAL 1	DEGRADERS & MOLECULAR GLUES CORAL 2	LARGE MOLECULE FORMULATION SEAGLASS	LARGE MOLECULE & RNA DELIVERY CORAL 3	
Donor to Donor Variability in Pharmacology: Mole Hill or Mountain?	Preclinical Evaluation Of Bms-986365: A First-In-Class Dual Ar Ligand-Directed Degrader And Antagonist For Mcrpc Therapy		Delivery Of Oligonucleotides To The Brain Via The Trans-Nasal Route In Neurodegenerative Rat Models	
 Major transition towards using primary cell assays in support of drug discovery programs for autoimmune diseases to improve translation An analysis of >25 primary cell assays used for HTS and SAR will be presented, including an answer to the question of donor to donor variability Specific lessons and strategies learned to facilitate the use of these physiologically relevant assays will be shared 	 BMS-986365 is a heterobifunctional AR degrader and antagonist for wildtype and mutant AR Inhibits AR target gene transcription and cell proliferation more effectively than enzalutamide Maintains low AR protein despite elevated transcript levels Shows robust efficacy in cell line and patient-derived xenograft models Clinically lowers PSA in mCRPC post-ARPI, effective in resistant disease 	Attendees are welcome to join co-located sessions		
FABIEN VINCENT, Consultant, Former Pharmacology Lab Head, Pfizer	SURENDRA NAYAK, Scientific Associate Director, Bristol Myers Squibb		ANISHA DSOUZA, Research Fellow, Massachusetts Eye and Ear Infirmary, Harvard Medical School	



15:40

16:05

F END OF CONFERENCE

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Featured Speakers: Discovery

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All Speakers

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Soak in Stunning Waterfront Views at Sheraton San Diego Resort

Sheraton San Diego Resort dazzles guests with spectacular views of the marina or downtown skyline. Experience San Diego with a central location in the heart of the harbor, yet just minutes from San Diego Airport with a free shuttle.

Enjoy being near downtown, Gaslamp Quarter, San Diego Zoo, and Petco Park, for a fun time in the city. Our newly renovated marina tower shines with coastal hues and modern amenities and our light-filled accommodations have a private balcony with seating. Enjoy picturesque views of the bay or marina in the classic bay tower guest rooms that also boast private balconies with seating. Stop by BreweryX for a refreshing brew or discover authentic Baja dishes at Rumorosa. Enjoy our sprawling outdoor pool complex, with ample opportunities to soak in the San Diego sun or lounge in poolside cabanas, cocktail in hand. Featuring over 130,000 sq ft of venue space accentuated by waterfront views, our San Diego Harbor Hotel is an inspired choice for weddings and meetings.

DIRECTIONS BY AIR

From San Diego International Airport (SAN), the hotel is just 1.5 miles from the airport.

Shuttle: Use the free Sheraton shuttle, available at Terminal 1 & Terminal 2. Look for the "Hotel Courtesy Shuttle" sign. If Shuttle is unavailable: Take the 992 bus from the airport to the Harbor Island Drive & Sheraton Way stop (5-minute ride). Walk 5 minutes to the hotel.

DIRECTIONS BY TRAM/BUS

The total journey from downtown via tram and bus should take around 20-30 minutes.

- Take the San Diego Trolley:
- Use the Green Line or another line that connects to the Santa Fe Depot station.
- Transfer to Bus or Walk:
- » From Santa Fe Depot, take Bus 992 (Airport Route) toward the San Diego International Airport. Exit at the hotel stop near Harbor Island.
- Alternatively, walk approximately 30 minutes (1.5 miles) along Harbor Drive to the hotel.



Sheraton San Diego Resort

1380 Harbor Island Drive, San Diego, CA



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