

	Friday Jun 2	Saturday Jun 3 General Techniques	Sunday Jun 4 Biophysics and binding	Monday Jun 5 Informatics	Tuesday Jun 6 Small molecule chemistry	Wednesday Jun 7	Thursday Jun 8 Biologics	Friday Jun 9 Case Studies	Saturday Jun 10	
8:30		Introduction (Directors & IT) (15') FEBS MIC (15')								
9	ARRIVALS (Dinner provided)	Historical Perspective (T. Blundell)	Measuring binding affinities (F. von Delft)	Machine learning in drug discovery (C. Deane)	General MedChem (M. Congreve)	Conformational landscaping of the SARS-COV-2 spike protein (G. Verkhivker)	Intro to Biologics (C. Deane)	Structure-based drug discovery targeting HIV reverse transcriptase (E. Arnold)	DEPARTURES	
9:45		Introduction to crystallography (C. Lesburg)	Protein Conformational Landscapes (M. Fischer)	Target selection drug discovery TBD	Compound design (A. Bradley)	Accelerating drug discovery (A. Kotecha)	Rational design (S.J. Fleishman)	Case study 3 (S. Jacob)		
10:30		<i>Coffee</i>	<i>Coffee</i>	<i>Coffee</i>	<i>Coffee</i>		<i>Coffee</i>	<i>Coffee</i>		
11:00		CryoEM in drug discovery (G. Scapin)	Computing Drug Binding Kinetics (R. Wade)	Data integration & pharmacology (A. Bradley)	In Silico SB optimization for accelerated drug discovery (E. Davis)	Talks from abstracts	Biologics Developability (R. Buick)	Talks from abstracts		
11:45		High-throughput and automation (F. von Delft)	Targeted Allosteric Modulators (G. Verkhivker)	Docking, Free energy calculation and Molecular Dynamics (S. Riniker)	3D structures, interactions, and implications (AY Sheikh)		Atoms up protein engineering (J. Williams)			
12:30		<i>Lunch</i>	<i>Lunch with poster preview</i>	<i>Lunch</i>		Excursion	<i>Lunch with poster preview</i>	<i>Lunch</i>		
2:30		Project Moonshot A. VonDelft	WS	WS	WS		WS	WS		Fragment-based design of MAT2 inhibitors (M. Schimpl)
3:15		Machine learning and protein structure prediction: present and future (C. Outeiral Rubiera)								Cu' havi 'na bona vigna, havi pani, vinu e ligna" – Startups (J. Williams)
4:00		<i>Coffee</i>	<i>coffee</i>	<i>Coffee</i>	<i>coffee</i>		<i>coffee</i>	<i>coffee</i>		
4:30		WS Intro	WS	WS	WS		WS	WS		Where will SB be in 10 years? (S. Jacob) + group discussion
5:15	Ice Breaker	Closing Remarks								
6:00	Intro to Erice	Poster Session Odd Numbers					Poster Session Even Numbers			

8:00		Welcome buffet	Dinner at posters				Dinner at posters	Farewell dinner	
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Workshop schedule

Sunday – 4 June		Monday – 5 June		Tuesday – 6 June		Thursday – 8 June	
2:30-4:00	4:30-6:00	2:30-4:00	4:30-6:00	2:30-4:00	4:30-6:00	2:30-4:00	4:30-6:00
Ligand crystallography /Hands on compound building <i>C. Lesburg</i>	Ligand crystallography /Hands on compound building <i>C. Lesburg</i>	General Machine Learning <i>Fergus Boyles</i>	General Machine Learning <i>Fergus Boyles</i>	Biologics design – Rosetta <i>Sarel Fleishman</i>		Hands-on single-particle cryoEM data analysis with cryoEDU <i>M. Herzik</i>	
General Machine Learning <i>Fergus Boyles</i>	General Machine Learning <i>Fergus Boyles</i>	Virtual Screening at Giga Scale <i>Steve Muchmore (OpenEye)</i>	Virtual Screening at Giga Scale <i>Steve Muchmore (OpenEye)</i>	In silico SB optimization for Accelerated Drug Discovery <i>Schroedinger</i>		In silico SB optimization for Accelerated Drug Discovery <i>Schroedinger</i>	
Hands-on single-particle cryoEM data analysis with cryoEDU <i>M. Herzik</i>	Hands-on single-particle cryoEM data analysis with cryoEDU <i>M. Herzik</i>	CCDC	CCDC	SmartEPU and G2 <i>A. Kotecha (TFS)</i>	SmartEPU and G2 <i>A. Kotecha (TFS)</i>	Biologics design – SabDab <i>Charlotte Deane</i>	Biologics design – SabDab <i>Charlotte Deane</i>