# Agenda of the International Workshop on Data-Driven Computational and Theoretical Materials Design (DCTMD2024)

(Dated on July 1st, 2024)

**Dates: October 9-13, 2024** 

Venue: Grand Central Hotel Shanghai, Shanghai, China

and Shanghai University (SHU), Shanghai, China

#### Overview

6 Plenary talks, 30 mins each, including 5 min discussion
28 Invited talks, 20 mins each, including 3 min discussion
16 Contributed talks, 15 mins each, including 2 min discussion
50+ Posters (~20 poster awards)
1 Panel discussion 1, 40 mins including 4 questions
1 Tutorial, 5 hours

#### Day 0: October 9, 2024 - Registration and Reception, Grand Central Hotel Shanghai

Time	Arrangement
1:00 PM - 8:00 PM	Arrival and registration at the hotel
5:00 PM - 8:00 PM	Buffet dinner at the hotel

### Day 1: October 10, 2024, Shanghai University (SHU)

**Morning Session** (Plenary talk 1, invited talk 4)

Time	Arrangement
9:00 AM - 10:00 AM	Moving from the hotel to SHU (~1 hour by conference bus or subway)
10:00 AM - 10:25 AM	Opening Ceremony & Welcome Remarks
10:25 AM - 10:55 AM	Plenary 1: Title
	Roberto Car (Princeton University, USA)
10:55 AM - 12:15 PM	Session 1-1: Machine-learned interatomic potential
	<b>I1-1</b> : Title
	T. Daniel Crawford (Virginia Tech, USA)
	<b>I1-2</b> : Title
	Linfeng Zhang (DP Technology, China)
	I1-3: "Accurate and efficient biomolecular dynamics enabled by
	machine-learned force fields"
	Alexandre Tkatchenko (Luxembourg University, Luxembourg)
	I1-4: "First-principles artificial intelligence"
	Yong Xu (Tsinghua University, China)

## Afternoon Session (Plenary talk 1, invited talk 4, Panel discussion 1, Poster 50)

Time	Arrangement
12:15 PM - 1:30 PM:	Lunch Break; Video play at noon
1:30 PM - 2:00 PM	Plenary 2: Title
	Yanming Ma (Jilin University, China)
2:00 PM - 3:20 PM	Session 1-2: Automatic, autonomous, self-driving experiments
	I1-5: "Creating Synergies between Experimental and Computational
	Approaches in Advanced Materials Design: Importance and Challenges
	of Clean Data"
	Annette Trunscke (FHI-Berlin, Germany)
	I1-6: Title
	Jun Jiang (USTC, China)
	<b>I1-7</b> : Title
	Keisuke Takahashi (Hokkaidao University, Japan)
	I1-8: Title
	Jungho Shin (KRICT, South Korea)
3:20 PM - 4:00 PM	Panel Discussion: "Unlocking the AI future of Materials Science"
	(40 min Panel discussions, four topics: Databases, Computations, AI
	algorithms, Autonomous experiments, 10 mins each)
	(Collection of oral discussions and written comments to form a future
	perspective on Data-driven materials design, possibly submitted a
	journal, e.g. Journal of Materials Informatics)
	Speakers: Hong Wang (Shanghai Jiaotong University),
4:00 PM - 5:00 PM	Poster session (~50+ posters)
5:00 PM - 6:30 PM	Moving from SHU to the Shanghai Bund (~1 hour by conference bus or
	subway)
6:30 - 9:00 PM	Dinner at the Shanghai Bund

## Day 2: October 11, 2024, Grand Central Hotel Shanghai

Morning Session (Plenary talk 1, Invited talk 5, Contributed talk 4)

Time	Arrangement
8:45 AM - 9:15 AM	Plenary 3: "Data Science to Optimize Metal-Organic Frameworks for
	Carbon Capture Applications"
	Berend Smit (EPFL, Switzerland)
9:15 AM - 10:35 AM	Session 2-1: AI-guided materials design
	<b>I2-1</b> : "How Landau theory can guide data science to find materials with
	targeted response"
	Turab Lookman (AiMaterials Research LLC, USA)
	<b>12-2</b> : "Polymer Informatics: Algorithmic Advances & Materials Design"

	Rampi Ramprasad (Georgia Tech, USA)
	<b>I2-3</b> : Title
	Timon Rabczuk (The Bauhaus-Universität Weimar, Germany)
	<b>I2-4</b> : "Machine learning based multiscale exploration and
	characterization of 2D materials"
	Xiaoying Zhuang (Leibniz University Hannover, Germany)
10:35 AM - 11:00 AM	Coffee Break
11:00 AM - 12:20 PM	Session 2-2:
	<b>12-5</b> : "What do we mean by new? Quantifying structural uniqueness in
	the era of generative crystal structure prediction"
	Taylor Sparks (The University of Utah, USA)
	C2-1: "From computational screening to the synthesis of a promising
	OER catalyst"
	Zhenpeng Yao (Shanghai Jiaotong University, China)
	C2-2: "Language Data-Driven Machine Learning Design of New
	Materials"
	Lei Zhang (Nanjing University of Information Science and Technology,
	Nanjing, China)
	C2-3: Title
	Speaker
	C2-4: "Progress in Machine Learning Studies for High-Entropy Alloys"
	Guangcun Shan (Beihang University, Beijing, China)

# Afternoon Session (Plenary talk 1, Invited talk 5, Contributed talk 4)

Time	Arrangement
12:20 PM - 1:30 PM	Lunch Break
1:30 PM - 2:00 PM	Plenary 4: Title
	Chris Wolverton (Northwestern University, USA)
2:00 PM - 3:20 PM	Session 2-3: AI-assisted computational materials design
	<b>12-6</b> : Title
	Zhipan Liu (Fudan University, China)
	I2-7: "Accurate materials modeling by machine learning and beyond
	DFT methods"
	Carla Verdi (The University of Queensland, Australia)
	I2-8: "Advancing Molecular Simulations with Machine-Learned
	Interatomic Potentials"
	Yangshuai Wang (National University of Singapore, Singapore)
	<b>12-9</b> : Title
	Yuanyuan Zhou (DTU, Denmark)
3:20 PM – 3:50 PM	Coffee Break
3:50 PM - 5:10 PM	Session 2-4:

	<b>I2-10</b> : Title
	Lixue Cheng (Microsoft Research AI for Science Lab)
	C2-5: "Constructing the highly accurate general AI interatomic potentials
	for the study of thermoelectric half-Heuslers"
	Jiong Yang (Shanghai University, Shanghai, China)
	<b>C2-6</b> : Title
	Speaker
	C2-7: Title
	Speaker
	C2-8: Title
	Speaker
6:00-8:00 PM	Banquet at the Wangbaohe Hotel (5 min walk)

Day 3: October 12, 2024, Grand Central Hotel Shanghai

Morning Session (Plenary talk 1, Invited talk 5, Contributed talk 4)

Time	Arrangement
8:45 AM - 9:15 AM	Plenary 5: "Describing Materials Properties and Functions via the
	"Materials Genes" Concept"
	Lucas Foppa (FHI-Berlin, Germany)
9:15 AM - 10:35 AM	Session 3-1: AI-assisted materials discovery
	<b>I3-1</b> : "Symbolic Regression in Materials Informatics: Applications and
	Challenges"
	Runhai Ouyang (Shanghai University, China)
	<b>I3-2</b> : "Finding Descriptors of Catalytic Properties from Data for Catalyst
	Design with the Help of Artificial Intelligence"
	Sergey V. Levchenko (Skotech, Russia)
	I3-3: "AI4Materials: From Simulation to Generation"
	Hongxia Hao (Microsoft Research AI for Science)
	<b>I3-4</b> : Title
	Speaker
10:35 AM - 11:00 AM	Coffee Break
11:00 AM - 12:20 PM	Session 3-2:
	<b>I3-5</b> : Title
	Speaker
	<b>C3-1</b> : Title
	Speaker
	<b>C3-2</b> : Title
	Speaker
	<b>C3-3</b> : Title
	Speaker

<b>C3-4</b> : Title
Speaker

## Afternoon Session (Plenary talk 1, Invited talk 5, Contributed talk 4)

Time	Arrangement
12:20 PM - 1:30 PM	Lunch Break
1:30 PM - 2:00 PM	Plenary 6: "AI Foundation models and Active Learning for Materials
	Discovery and Process Design"
	Xiaonan Wang (Tsinghua University, China)
2:00 PM - 3:20 PM	Session 3-3: Databases and large models
	<b>I3-6</b> : Title
	Junfeng Qiao (EPFL, Switzerland)
	I3-7: "Ontology and the role of semantics in the development of
	knowledge-based materials databases"
	Lauren Takahashi (Hokkaido University, Japan)
	I3-8: "Generating public datasets for chemistry"
	Karmen Čondić-Jurkić (Open Molecular Software Foundation, New
	Zealand)
	<b>I3-9</b> : Title
	Speaker
3:20 PM - 3:50 PM	Coffee Break
3:50 PM - 5:10 PM	Session 3-4:
	<b>I3-10</b> : Title
	Speaker
	C3-5: "Unexpected Failure and Success in Data-Driven Materials
	Science"
	Kangming Li (University of Toronto, Canada)
	C3-6: "Simulating the Microscopic World: From Schrödinger Equation
	to Large Atomic Models"
	Han Wang (Institute of Applied Physics and Computational Mathematics,
	Beijing, China)
	C3-7: "A Large Multi-Modality Model for Chemistry and Materials
	Science"
	Xin Chen (Suzhou Laboratory, Suzhou, China)
	C3-8: Title
	Speaker
5:10 PM - 5:40 PM	Closing Ceremony: Awards, Summary & Farewell Remarks
6:00 PM - 8:00 PM	Buffet dinner at the hotel

Day 4: October 13, 2024 Tutorial (optional) at Shanghai University (SHU) and Departure

Time	Arrangement
8:30 AM - 9:30 AM	Moving from the hotel to SHU (~1 hour by conference bus or subway)
9:30 AM - 5:00 PM	Tutorial: "DeepMD: from algorithms to applications"
	Yibo Wang (DP Technology, China) et al.

Website of the DeepMD tutorial (Introduction and Registration):

https://bohrium.dp.tech/courses/1347727500?tab=courses&lang=en-us

Time	Торіс
09:30 - 10:30	Machine Learning Potentials: From DeePMD to DPA-2
10:30 - 11:30	Introduction to DeePMD-kit & DP-GEN
11:30 - 12:00	DeePMD-kit hands-on
13:00 - 14:30	Applications of DeePMD in Materials Research
14:30 - 16:00	Open Large Atomic Model: Advances in Alloy Research
16:00 - 17:00	Q&A Session

Note: The DeepMD tutorial has limited seats and needs additional registration. The tutorial is free for the registered participants at DCTMD (course materials and instruction in English).

Note: This agenda including the presentation types is subject to adjustments after evaluation by the organization committee. For updates, speaker details, and session topics, please visit <a href="International workshop on data-driven computational and theoretical materials design (scievent.com">International workshop on data-driven computational and theoretical materials design (scievent.com</a>) (https://dctmd2024.scievent.com/).