

THE PROGRAM

of the 10th edition of the International Workshop of Materials Physics

14 May, 2025

INVITED ORAL PRESENTATIONS

08:15	Registration – Oteteleşanu Hall
08:45	Official opening – Lucian PINTILIE, Scientific Director of NIMP
Session 1: Chairman – Lucian Pintilie	
9:00	Jorge INIGUEZ Luxembourg Institute of Science and Technology Hafnia ferroelectrics: a singular challenge for theory and simulation
9:40	Sergiu CLIMA IMEC, Leuven, BELGIUM First-principles screening for OTS/SOM chalcogenide materials
10:20	Santanu SAHA University of Limoges and Institut de Recherche sur les Céramiques (IRCER), Limoges, FRANCE Adapting Physico-Chemical Affinities in First Principles Design of New Materials
11:00	Coffee break
11:15	Adolfo O. FUMEGA Academy of Finland, Aalto University, FINLAND Computational modeling of emergent phases in moiré materials
11:55	Andrei BERNEVIG Princeton University, USA Moiré Systems as Quantum Simulators of Strongly Correlated Hamiltonians
12:35	Lunch
Session 2: Chairman – Marius Husanu	
14:00	Bilal TANATAR Department of Physics, Bilkent University, Ankara, TURKEY Studies of localization within Aubry-Andre Model
14:40	Krzysztof WOHLFELD Institute of Theoretical Physics of the Faculty of Physics of the University of Warsaw, POLAND Altermagnetism from a strong-coupling perspective
15:20	Jan MINAR New Technologies Research Center (NTC), University of West Bohemia, Pilsen, CZECHIA Quantum Materials and Magnetic Phenomena Studied by Spin-Resolved ARPES: Theoretical perspectives
16:00	Coffee break
16:15	Andrei MANOLESCU Reykjavik University, ICELAND Magnetic-flux periodicity in core/shell semiconductor/superconductor nanowires
16:55	Cătălin PASCU MOCA University of Oradea, ROMANIA Kondo Compensation and Quantum Criticality in Superconducting and Fractionalized Systems
17:35	Poster Sessions
19:30	Dinner

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Session 3: Chairman – Alin Velea	
9:00	Stephan ROCHE ICREA Institució Catalana de Recerca i Estudis Avancats, Barcelona, SPAIN Linear scaling quantum simulation methodologies
9:40	Marius STAN University of Chicago, USA Artificial Intelligence for Material Design
10:20	Alin M. ELENA Scientific Computing Department, Science and Technology Facilities Council, Daresbury Laboratory, Daresbury, UK Machine Learned Interatomic Potentials for modelling porous materials
11:00	Coffee break
11:15	Felix MOCANU Department of Materials, University of Oxford, Oxford, UK Computational design of high energy density cathode materials for Li-ion batteries
11:55	Larisa VON RIEWEL Heraeus Noblelight, Global Innovation, Bayern, GERMANY Microscopic investigation of the slurry drying process and binder migration in Li-ion battery anodes using computational chemistry
12:35	Lunch
Session 4: Chairman – Cristian Teodorescu	
14:00	Claudiu GENES Max Planck Institute, Erlangen, GERMANY Quantum optics with atoms and molecules: analytical approaches
14:40	Vidar GUDMUNDSSON Science Institute, University of Iceland, ICELAND Arrays of quantum dots or rings in a FIR-photon cavity
15:20	Coffee break
15:35	Krzysztof WIECZERZAK Department of Materials Science, Faculty of Mechanical Engineering and Aeronautics, Rzeszow University of Technology, Rzeszow, POLAND ONLINE PRESENTATION Exploring CuAgZr metallic glasses for biomedical use: A study using combinatorial synthesis, high-throughput experiments, and machine learning
16:15	G. A. NEMNES Physics Faculty, Bucharest University Hybrid perovskite materials: from perovskite solar cells to memristive elements
17:30	Poster Sessions
19:30	Dinner