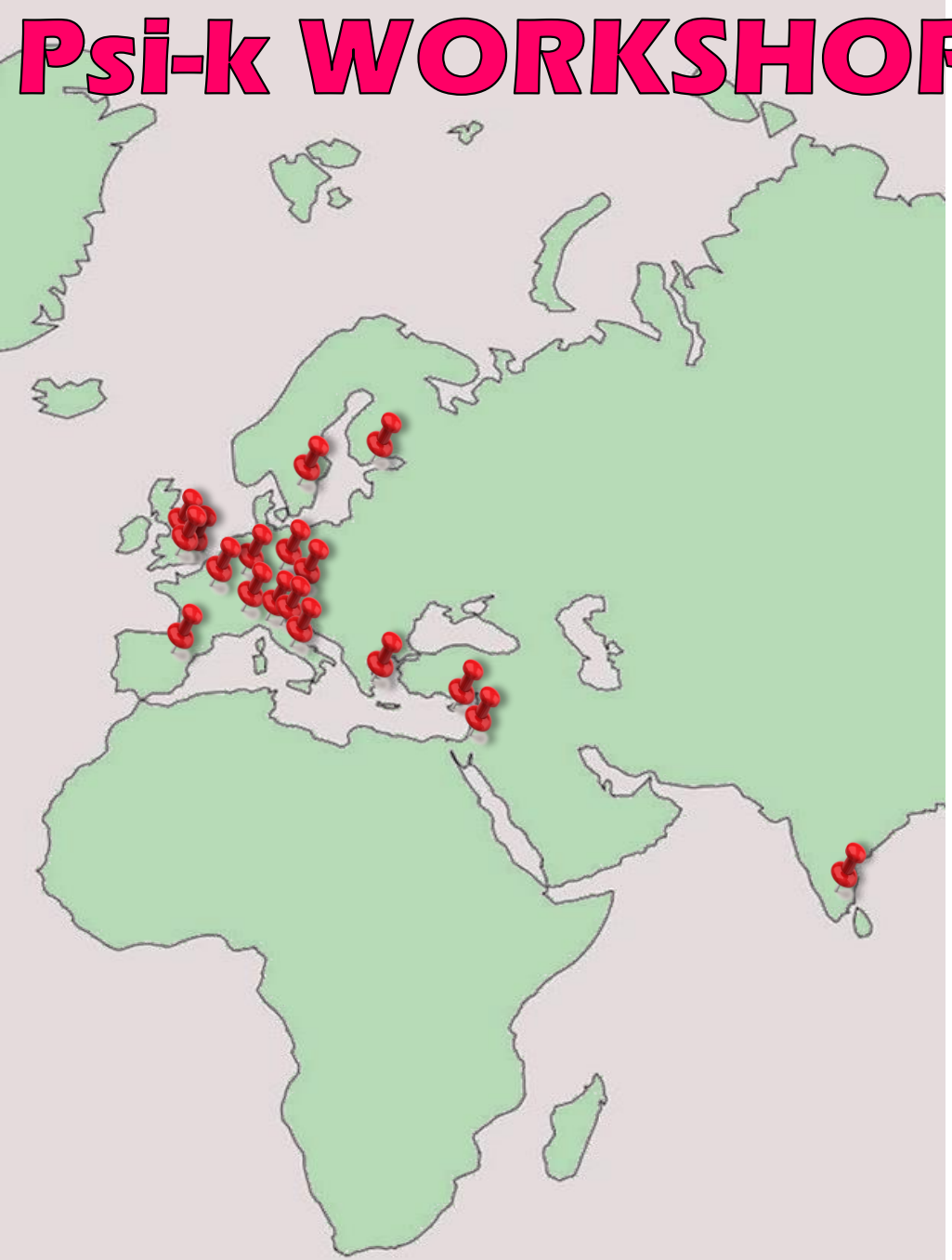
Psi-k WORKSHOPS 2023



For more information and to register for events please visit
www.psi-k.net/workshops

Registered charity number 1126308
Registered company number 06440198

27 February - 3 March, Aalto, Finland - *Actively Learning Material Science*

24-28 April, Daresbury, UK - *Open Science with the Atomic Simulation Environment*

8-12 May, Jülich, Germany - *All electron DFT with Fleur - a Hands-on Tutorial*

22-26 May, Rome, Italy - *Ab initio many-body perturbation theory: from equilibrium to time-resolved spectroscopies and nonlinear optics*

23-26 May, Paris, France- *Young Researcher's School on Theory and Simulation in Electrochemical Conversion Processes*

4-8 June, Tel Aviv, Israel - *Advanced School: Path Integral Quantum Mechanics*

12-16 June, Zaragoza, Spain - *19th ETSF Young Researchers' Meeting 2023*

19-23 June, Berlin, Germany - *CECAM/Psi-k Conference: Bridging length scales with machine learning: from wavefunctions to thermodynamics*

3-5 July, Modena, Italy - *Principles of Light-Induced Charge Transfer for Optogenetics*

24-28 July, Bangalore, India - *New Approaches and Machine learning Methods for Ab initio calculations (NAMMA 2023): Psi-k India Workshop and Conference*

2-10 August, Berlin, Germany - *HoW exciting! Hands-on Workshop on Excitations in Solids 2023 (Short: HoW exciting! 2023)*

7-11 August, Örebro, Sweden - *Progress in Non-equilibrium Green's Functions 8 (PNGF8)*

28 August - 1 September, Pavia, Italy - *Advanced Quantum ESPRESSO school: Hubbard and Koopmans functionals from linear response*

4-7 September, Athens, Greece - *First-principles Green function formalisms: algorithms, method developments and applications to spinorbitronics and magneto-superconductivity*

4-7 September, Freiburg, Germany - *FIT and livMatS biohackers challenge: excited-state dynamics and applications*

18-23 September, London, UK - *MSSC2023 - Ab initio Modelling in Solid State Chemistry – blended/hybrid edition - new users*

20-22 September, Manchester, UK - *Twistronics of 2D materials: from modelling to real systems*

26-29 September, Lausanne, Switzerland - *Atto2Nano: modeling ultrafast dynamics across time-scales in condensed matter*

3-8 October, Paphos, Cyprus - *Towards exascale solutions in Green function methods and advanced DFT*

29 October - 3 November, Tegernsee, Germany - *Ab initio Description of Iron and Steel (ADIS2023): Digitalization and Workflows*

30 October - 10 November, Les Houches, France - *Blending the DFT-based multiple scattering Greens' function approach to spectroscopies with machine learning*

6-10 November, Aalto, Finland - *School on Machine Learning Interatomic Potentials ("ML-IP 2023")*

19 Feb-1 Mar 2024, Lausanne, Switzerland - *Electronic Structure Software Development: Advancing the Modular Paradigm*

11-15 March 2024, Lausanne, Switzerland - *Theoretical Spectroscopy Lectures*

12-14 March 2024, Aalto, Finland - *LOBSTER School on Chemical Bonding Analysis*