School on **Electron-Phonon Physics** from First Principles

19 - 23 March 2018 Trieste, Italy

The School addresses senior PhD students and experienced researchers with prior working knowledge of DFT.

Theoretical and hands-on training will focus on ab-initio calculations of many properties relating to the electron-phonon interaction, for applications in condensed matter physics, materials physics, and nanoscience.

Starting from an introduction to the background on electron-phonon physics and related materials properties from the point of view of abinitio calculations, we will show the participants how to perform cuttingedge electron-phonon calculations using a suite of electronic structure codes, including EPW, Wannier90, Quantum ESPRESSO, and ABINIT.

Topics:

- Density functional perturbation theory
- Electron-phonon coupling
- Phonon-assisted optical absorption
- Maximally-localized Wannier
- Temperature dependence of the electronic bandstructure
- Electronic transport using the Boltzmann transport equation
- Phonon-driven superconductivity

Further information: http://indico.ictp.it/event/8301/ smr3191@ictp.it

ICTP

Directors:

S. Poncé, University of Oxford E.R. Margine, Binghamton University - SUNY F. Giustino, University of Oxford

Local Organizers:

R. Gebauer, ICTP N. Seriani, ICTP

Lecturers:

- S. de Gironcoli, SISSA
- P. Giannozzi, Universiy of Udine
- F. Giustino, University of Oxford
- X. Gonze, Univ. Catholique of Louvain
- E. Kioupakis, University of Michigan
- E. R. Margine, Binghamton University-SUNY
- G. Pizzi, EPFL
- S. Poncé, University of Oxford

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How to apply:

Online application: http://indico.ictp.it/event/8301/

Female scientists are encouraged to apply.



A limited number of grants are available to support the attendance of selected participants, with priority given to participants from developing countries. There is no registration fee.

Deadlines: 1 December 2017

for those needing visa and/or financial support

15 January 2018

otherwise









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