

MARCO BERNARDI

CONTACT INFORMATION

E-mail: bmarco@caltech.edu
Phone: 626-395-2515
<http://bernardi.caltech.edu>

ADDRESS

California Institute of Technology
1200 E California Blvd. MC 107-81
Pasadena CA, 91125

RESEARCH INTERESTS

- Materials physics, electronic structure theory, computational physics
- Interactions, transport, and nonequilibrium dynamics in conventional and quantum materials
- Materials for electronics, energy, and quantum technologies

EDUCATION

- **Massachusetts Institute of Technology**, Cambridge, MA, USA
Ph.D. in Materials Science, June 2013. Advisor: Prof. Jeffrey C. Grossman
- **University of Rome Tor Vergata**, Rome, Italy
M.Sc. (Laurea Specialistica) in Materials Science, January 2008. *Summa Cum Laude*
- **University of Rome La Sapienza**, Rome, Italy
B.Sc. (Laurea Triennale) in Chemistry, November 2004. *Summa Cum Laude*

RESEARCH AND PROFESSIONAL EXPERIENCE

- **Professor**
Department of Applied Physics and Materials Science, and Department of Physics
California Institute of Technology. Pasadena, CA, USA
Dates: June 2021 –
Leads a research group focused on understanding electron interactions and dynamics in materials
- **Assistant Professor**
Department of Applied Physics and Materials Science
California Institute of Technology. Pasadena, CA, USA
Dates: September 2015 – May 2021
- **Post-Doctoral Fellow**
Physics Department, University of California, Berkeley and
Materials Science Division, Lawrence Berkeley National Laboratory, Berkeley, CA, USA
Dates: September 2013 – August 2015
Supervisors: Prof. Steven G. Louie, Prof. Jeffrey B. Neaton
Developed new first-principles methods for electron dynamics in materials
- **Research Assistant**
Massachusetts Institute of Technology, Cambridge, MA, USA
Department of Materials Science and Engineering
Dates: September 2008 – August 2013
Supervisor: Prof. Jeffrey C. Grossman
Studied novel materials and physical processes for solar energy conversion

HONORS AND AWARDS

- 2020 ISSNAF “Franco Strazzabosco” Young Investigator Award
- 2019 Emerging Young Investigator Award at the 4th Functional Oxide Thin Films Conference
- 2018 NSF CAREER Award
- 2017 Air Force Young Investigator Award
- 2015 Psi-K Volker Heine Young Investigator Award
- 2011 Intel Ph.D. Fellowship
- 2009 MIT DMSE 1st Year Graduate Student Exceptional Performance Award
- 2007 Australian Endeavor Research Fellowship

TEACHING

- Structure and Bonding in Materials (MS 131). Taught every year from 2016 to 2023
- Introduction to Computational Methods for Science and Engineering (MS 141). Taught from 2020 to 2023
- Computational Solid State Physics and Materials Science (APh 256). Taught in 2018

CURRENT GROUP MEMBERS

Graduate students

Applied Physics: Dhruv Desai, Yao Luo, David Abramovitch, Shaelyn Iyer. *Materials Science:* Sergei Kliavinek.
Physics: Jia (Kelly) Yao, Ina Sorensen. *Chemistry:* Khoa Le, Laurie Tan

Postdocs

Shiyu Peng

PROFESSIONAL ACTIVITIES AND AFFILIATIONS

- **Reviewer:** Nature, Science, Nature Materials, Nature Physics, Nature Nanotechnology, Nature Communications, Science Advances, Physical Review Letters, Physical Review X, Physical Review B, Physical Review Materials, Nano Letters, Materials Today Physics, Journal of Physical Chemistry Letters (among others).
- **Memberships:** American Physical Society, American Chemical Society, Materials Research Society.
- **Organizer:** Organized invited symposia on ultrafast dynamics at the 2020, 2021, and 2024 APS March Meetings.
- **Consulting:** Reviewed high-performance computing proposals for ALCC in the US and PRACE in Europe.
- **Editor:** Section editor for the 2019 Handbook of Materials Modeling (Springer).

RECENT INVITED TALKS

- **2024 Psi-K Workshop: Electronic Structure Simulations for Large-Scale Facilities.** EPFL, Switzerland.
New Directions in Electron-Phonon Calculations.
- **2024 Physics Colloquium.** University of Pisa, Italy.
Precise Calculations of Electron and Phonon Dynamics.
- **2024 NEST Nano Colloquium.** NEST, Pisa, Italy.
Precise First-Principles Calculations of Electron and Spin Dynamics.
- **2024 MRS Spring Meeting.** Seattle, WA.
Precise Calculations of Strong Electronic Interactions and Transport in Oxides.
- **2024 GraFOx Seminar at Paul Drude Institute.** Berlin, Germany.
Building the Computational Toolbox for Quantum Materials:
Precise First-Principles Calculations of Electron and Spin Dynamics.
- **2023 Materials Science and Engineering Colloquium.** Columbia University, New York.
Building the Computational Toolbox for Quantum Materials:
Precise First-Principles Calculations of Electron and Spin Dynamics.
- **2023 35th Annual Workshop on Recent Developments in Electronic Structure Methods.** Merced, CA.
Advances in Electron-Phonon Interactions and Spin Dynamics from First Principles.
- **2023 Quantum Foundry Seminar.** UC Santa Barbara, CA.
Precise First-Principles Calculations of Electron and Spin Dynamics: Building the Toolbox for Quantum Materials.
- **2023 Sanibel Symposium: Spin Workshop.** University of Florida, FL.
Theory and First-Principles Calculations of Spin-Phonon Interactions and Spin Relaxation.
- **2023 2nd Quantum Matters in Materials Science Workshop.** NIST, USA (virtual).
Advances in First-Principles Calculations of Electron and Spin Dynamics in Quantum Materials.
- **2023 APS March Meeting.** Las Vegas, NV.
Theory and First-Principles Calculations of Spin-Phonon Interactions and Spin Relaxation.
- **2023 SIAM Conference on Computational Science and Engineering.** Amsterdam, Netherlands.
Nonequilibrium Dynamics of Interacting Electrons, Phonons and Excitons from First Principles.
- **2023 5th Functional Oxide Thin Films Conference.** Cancun, Mexico.
First-Principles Calculations of Strong Electronic Interactions in Complex Oxides.
- **2022 Vienna Quantum Seminar Lecture.** Vienna, Austria.
Precise and Parsimonious Computational Quantum Physics: From Electrons in Materials to Quantum Circuits.
- **2022 The 23rd Asian Workshop on First-Principles Electronic Structure Calculations (Plenary Talk).** Virtual.
Frontiers of First-Principles Electron-Phonon Interactions: Weak-to-Strong, Correlated, Spinful, and Data-Driven.
- **2022 ETH Zurich Workshop on First-Principles Modeling of Defects in Solids.** Zurich, Switzerland.
Predicting Electronic Interactions and Transport Governed by Polarons and Defects.
- **2022 ICTP Workshop on Thermal Transport.** Virtual.
Advances in Computing Electron Interactions and Dynamics from First Principles.
- **2022 IPAM Workshop on Model Reduction in Quantum Mechanics.** Los Angeles CA, USA.
Precise Quantum Mechanical Calculations of Electron Interactions and Dynamics in Condensed Matter.
- **2022 MRS Spring Meeting.** Honolulu HI, USA.
Nonequilibrium Dynamics of Interacting Electrons, Phonons and Excitons from First Principles.
- **2022 ACS Spring Meeting.** San Diego CA, USA.
Precise First-Principles Tools for Electron Dynamics in Quantum Materials.

PUBLICATIONS

Metrics – *h*-index: 40 Citations: 8,000 (see [Google Scholar](#))

Total of 70 peer reviewed publications. For the full publication list, see the [group website](#)

Selected Publications

(see the [magenta links](#) for press articles)

1. Y. Luo, D. Desai, B.K. Chang, J. Park, **M. Bernardi**
Data-Driven Compression of Electron-Phonon Interactions.
Physical Review X **2024** 14 021023. DOI: [10.1103/PhysRevX.14.021023](https://doi.org/10.1103/PhysRevX.14.021023)
See the [press article from Caltech News](#)
2. D. Abramovitch, J.-J. Zhou, J. Mravlje, A. Georges, **M. Bernardi**
Combining Electron-Phonon and Dynamical Mean Field Theory Calculations of Correlated Materials:
Transport in the Correlated Metal Sr₂RuO₄.
Physical Review Materials **2023** 7, 093801 (Editor's Suggestion). DOI: [10.1103/PhysRevMaterials.7.093801](https://doi.org/10.1103/PhysRevMaterials.7.093801)
3. J. Park, J.-J. Zhou, Y. Luo, **M. Bernardi**
Predicting Phonon-Induced Spin Decoherence from First Principles:
Colossal Spin Renormalization in Condensed Matter.
Physical Review Letters **2022** 129, 197201. DOI: [10.1103/PhysRevLett.129.197201](https://doi.org/10.1103/PhysRevLett.129.197201)
See the [press article from Caltech News](#)
4. J.-J. Zhou, J. Park, I. Timrov, A. Floris, M. Cococcioni, N. Marzari and **M. Bernardi**
Ab Initio Electron-Phonon Interactions in Correlated Electron Systems.
Physical Review Letters **2021**, 127, 126404. DOI: [10.1103/PhysRevLett.127.126404](https://doi.org/10.1103/PhysRevLett.127.126404)
See the [press article from EPFL](#)
5. J.-J. Zhou, J. Park, I.-T. Lu, I. Maliyov, X. Tong, and **M. Bernardi**
PERTURBO: A Software Package for Ab Initio Electron-Phonon Interactions, Charge Transport and Ultrafast Dynamics.
Computer Physics Communications **2021** 264, 107970. DOI: [10.1016/j.cpc.2021.107970](https://doi.org/10.1016/j.cpc.2021.107970)
See the [press article in the Caltech News](#)
The code can be downloaded at <https://perturbo-code.github.io/>
6. V. Jhalani, J.-J. Zhou, J. Park, C.E. Dreyer, and **M. Bernardi**
Piezoelectric Electron-Phonon Interaction from Ab Initio Dynamical Quadrupoles:
Impact on Charge Transport in Wurtzite GaN.
Physical Review Letters **2020**, 125, 136602. DOI: [10.1103/PhysRevLett.125.136602](https://doi.org/10.1103/PhysRevLett.125.136602)
7. H.-Y. Chen, D. Sangalli, and **M. Bernardi**
Exciton-Phonon Interaction and Relaxation Times from First Principles.
Physical Review Letters **2020**, 125, 107401. DOI: [10.1103/PhysRevLett.125.107401](https://doi.org/10.1103/PhysRevLett.125.107401)
8. J. Park, J.-J. Zhou, and **M. Bernardi**
Spin-Phonon Relaxation Times in Centrosymmetric Materials from First Principles.
Physical Review B **2020**, 101, 045202. DOI: [10.1103/PhysRevB.101.045202](https://doi.org/10.1103/PhysRevB.101.045202)
9. J.-J. Zhou, and **M. Bernardi**
Predicting Charge Transport in the Presence of Polarons:
The Beyond-Quasiparticle Regime in SrTiO₃.
Physical Review Research **2019**, 1, 033138. DOI: [10.1103/PhysRevResearch.1.033138](https://doi.org/10.1103/PhysRevResearch.1.033138)
See the [press article in the Caltech News](#)

10. I.-T. Lu, J.-J. Zhou, and **M. Bernardi**
Efficient Ab Initio Calculations of Electron-Defect Scattering and Defect-Limited Carrier Mobility.
Physical Review Materials **2019**, 3, 033804 (Editor's Suggestion). DOI: [10.1103/PhysRevMaterials.3.033804](https://doi.org/10.1103/PhysRevMaterials.3.033804)
11. J.-J. Zhou, O. Hellman and **M. Bernardi**
Electron-Phonon Scattering in the Presence of Soft Modes
and Electron Mobility in SrTiO₃ Perovskite from First Principles.
Physical Review Letters **2018**, 121, 226603. DOI: [10.1103/PhysRevLett.121.226603](https://doi.org/10.1103/PhysRevLett.121.226603)
See the [press article in the Caltech News](#)
12. K. Frohna, T. Deshpande, J. Harter, B. Barker, J. B. Neaton, S. G. Louie, O. Bakr, D. Hsieh and **M. Bernardi**
Inversion Symmetry and Bulk Rashba Effect in Methylammonium Lead Iodide Perovskite Single Crystals.
Nature Communications **2018**, 9, 1829. DOI: [10.1038/s41467-018-04212-w](https://doi.org/10.1038/s41467-018-04212-w)
See the [press article in the Caltech News](#)
13. V. Jhalani, J.-J. Zhou and **M. Bernardi**
Ultrafast Hot Carrier Dynamics in GaN and its Impact on the Efficiency Droop.
Nano Letters **2017**, 17, 5012. DOI: [10.1021/acs.nanolett.7b02212](https://doi.org/10.1021/acs.nanolett.7b02212)
See the [press articles in the Caltech News](#) and [phys.org](#)
14. J.-J. Zhou and **M. Bernardi**
Ab Initio Electron Mobility and Polar Phonon Scattering in GaAs.
Physical Review B (Rapid Communication) **2016**, 94, 201201. DOI: [10.1103/PhysRevB.94.201201](https://doi.org/10.1103/PhysRevB.94.201201)
15. **M. Bernardi**
First-Principles Dynamics of Electrons and Phonons.
European Journal of Physics B **2016**, 89, 239. DOI: [10.1140/epjb/e2016-70399-4](https://doi.org/10.1140/epjb/e2016-70399-4)
16. **M. Bernardi**, J. Mustafa, J. B. Neaton, and S. G. Louie
Theory and Computation of Hot Carriers Generated by Surface Plasmon Polaritons in Noble Metals.
Nature Communications **2015**, 6:7044. DOI: [10.1038/ncomms8044](https://doi.org/10.1038/ncomms8044)
17. M. Palummo*, **M. Bernardi***, and J. C. Grossman
Exciton Radiative Lifetimes in Two-Dimensional Transition Metal Dichalcogenides.
Nano Letters **2015**, 15, 2794. *Equal contributors. DOI: [10.1021/nl503799t](https://doi.org/10.1021/nl503799t)
18. **M. Bernardi**, D. Vigil-Fowler, J. Lischner, J. B. Neaton, and S. G. Louie
Ab Initio Study of Hot Carriers in the First Picosecond after Sunlight Absorption in Silicon.
Physical Review Letters **2014**, 112, 257402. DOI: [10.1103/PhysRevLett.112.257402](https://doi.org/10.1103/PhysRevLett.112.257402)
See the [feature articles in the Berkeley Lab News](#)
19. **M. Bernardi**, M. Palummo, and J. C. Grossman
Extraordinary Sunlight Absorption and 1 nm-Thick Photovoltaics using 2D Monolayer Materials.
Nano Letters **2013**, 13, 3664. DOI: [10.1021/nl401544y](https://doi.org/10.1021/nl401544y)
See the [press articles by the MIT News](#) and [NERSC](#). Cited over 2000 times.