

# Benjamin Cohen-Stead

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## EDUCATION

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<b>University of California, Davis</b> Ph.D. in Physics	Davis, Ca 2016–2022
<ul style="list-style-type: none"><li>• Advisor: Prof. Richard Scalettar</li><li>• Thesis: A Hybrid Quantum Monte Carlo Method for Electron-Phonon Models</li></ul>	
<b>Whitman College</b> B.A. in Physics	Walla Walla, Wa 2010–2014

## PROFESSIONAL EXPERIENCE

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<b>University of Tennessee, Knoxville</b> Research Assistant Professor	Knoxville, Tn 2024–present
<b>University of Tennessee, Knoxville</b> Postdoctoral Research Associate	Knoxville, Tn 2022–2024
<ul style="list-style-type: none"><li>• Principal Investigator: Prof. Steven Johnston</li></ul>	
<b>Los Alamos National Laboratory</b> Graduate Student Researcher	Los Alamos, NM 2020–2022
<ul style="list-style-type: none"><li>• Mentor: Dr. Kipton Barros</li><li>• Project: Langevin Methods for Quantum Electron-Phonon Simulations</li></ul>	
<b>Picarro, Inc.</b> Associate Data Scientist	Santa Clara, Ca 2014–2016
<ul style="list-style-type: none"><li>• Responsibilities: software development, algorithm design, data analysis</li></ul>	
<b>Picarro, Inc.</b> Data Science Intern	Santa Clara, Ca Summer 2014
<b>University of Rochester, Department of Physics</b> Physics REU Student, Mentor: Professor Stephen Teitel	Rochester, NY Summer 2013
<ul style="list-style-type: none"><li>• Research: numerical investigation of two-dimensional granular systems</li></ul>	

## GRANTS & AWARDS

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<b>Scientific Software Research Faculty (SSRF) Award</b> Organization: Simons Foundation	2024–2029
<ul style="list-style-type: none"><li>• Principle Investigator: Dr. Benjamin Cohen-Stead</li><li>• Project: User-friendly and Extensible Quantum Monte Carlo Related Tools</li></ul>	
<b>UC-National Lab In-Residence Graduate Fellowship</b> Organization: Los Alamos National Laboratory (LANL)	2020–2022
<ul style="list-style-type: none"><li>• Mentor: Dr. Kipton Barros</li><li>• Project: Langevin Methods for Quantum Electron-Phonon Simulations</li></ul>	

## OPEN SOURCE CODES

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**SmoQyDQMC.jl:** <https://github.com/SmoQySuite/SmoQyDQMC.jl.git>

Flexible implementation of the determinant quantum Monte Carlo method for simulating Hubbard and electron-phonon interactions.

**SmoQyKPMCore.jl:** <https://github.com/SmoQySuite/SmoQyKPMCore.jl.git>

A lightweight and flexible implementation of the Kernel Polynomial Method (KPM) for approximating function of Hermitian matrices.

**SmoQySynthAC.jl:** <https://github.com/SmoQySuite/SmoQySynthAC.jl.git>

A package for generating synthetic noisy correlation function data intended to mimic that generated by a quantum Monte Carlo simulation. This package is useful for testing efficacy of various analytic continuation methods.

**SmoQyHankelCorrCleaner.jl:** <https://github.com/SmoQySuite/SmoQyHankelCorrCleaner.jl.git>

Julia package exporting methods for denoising imaginary time correlation data using the Hankel projection method introduced in arXiv:2403.12349

**JDQMCFramework.jl:** <https://github.com/SmoQySuite/JDQMCFramework.jl.git>

Julia package exporting a suite of types and routines useful for writing a determinant quantum Monte Carlo code.

**JDQMCMeasurements.jl:** <https://github.com/SmoQySuite/JDQMCMeasurements.jl.git>

Julia package implementing various correlation measurements that are frequently made in determinant quantum Monte Carlo simulations.

**StableLinearAlgebra.jl:** <https://github.com/SmoQySuite/StableLinearAlgebra.jl.git>

Exports numerically stable linear algebra routines used in determinant quantum Monte Carlo codes.

**Checkerboard.jl:** <https://github.com/SmoQySuite/Checkerboard.jl.git>

Implements the checkerboard approximation for representing exponentiated kinetic for tight-binding models.

**LatticeUtilities.jl:** <https://github.com/SmoQySuite/LatticeUtilities.jl.git>

Julia package for representing arbitrary periodic lattice geometries.

**MuTuner.jl:** <https://github.com/cohensbw/MuTuner.jl.git>

Implements a method for tuning the chemical potential in grand canonical Monte Carlo simulations to achieve a target particle density.

## PATENTS

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**Aggregate leak indicator display systems and methods**

US Patent Number: 10962437

Assignee: Picarro, Inc.

Date of Patent: March 30, 2021

- Inventors: A. Nottrott, S. MacMullin, S.M. Tan, B. Cohen-Stead, C. Rella

## PUBLICATIONS

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- [1] B. Cohen-Stead, S. M. Costa, J. Neuhaus, A. T. Ly, Y. Zhang, R. Scalettar, K. Barros, and S. Johnston, “SmoQyDQMC.jl: A flexible implementation of determinant quantum Monte Carlo for Hubbard and electron-phonon interactions”, *SciPost Phys. Codebases*, p. 29, 2024. DOI: 10.21468/SciPostPhysCodeb.29.
- [2] B. Cohen-Stead, K. Barros, R. Scalettar, and S. Johnston, “A hybrid Monte Carlo study of bond-stretching electron–phonon interactions and charge order in BaBiO<sub>3</sub>”, *npj Computational Materials*, vol. 9, no. 1, p. 40, 2023. DOI: 10.1038/s41524-023-00998-6.

- [3] B. Cohen-Stead, O. Bradley, C. Miles, G. Batrouni, R. Scalettar, and K. Barros, “Fast and scalable quantum Monte Carlo simulations of electron-phonon models”, *Phys. Rev. E*, vol. 105, p. 065302, 6 Jun. 2022. DOI: 10.1103/PhysRevE.105.065302.
- [4] B. Cohen-Stead, K. Barros, Z. Meng, C. Chen, R. T. Scalettar, and G. G. Batrouni, “Langevin simulations of the half-filled cubic holstein model”, *Phys. Rev. B*, vol. 102, p. 161108, 16 Oct. 2020. DOI: 10.1103/PhysRevB.102.161108.
- [5] B. Cohen-Stead, N. C. Costa, E. Khatami, and R. T. Scalettar, “Effect of strain on charge density wave order in the holstein model”, *Phys. Rev. B*, vol. 100, p. 045125, 4 Jul. 2019. DOI: 10.1103/PhysRevB.100.045125.
- [6] M. Naamneh, E. C. O’Quinn, E. Paris, D. McNally, Y. Tseng, W. R. Pudłko, D. J. Gawryluk, J. Shamblin, B. Cohen-Stead, M. Shi, M. Radovic, M. K. Lang, T. Schmitt, S. Johnston, and N. C. Plumb, “Persistence of small polarons into the superconducting doping range of  $\text{Ba}_{1-x}\text{K}_x\text{BiO}_3$ ”, *Phys. Rev. Res.*, vol. 7, p. 043082, 4 Oct. 2025. DOI: 10.1103/s3p1-cy1s.
- [7] Y. Zhang, P. M. Dee, B. Cohen-Stead, T. A. Maier, S. Johnston, and R. Scalettar, “Optimizing the critical temperature and superfluid density of a metal-superconductor bilayer”, *Phys. Rev. B*, vol. 112, p. 064510, 6 Aug. 2025. DOI: 10.1103/lcgr-bqcv.
- [8] A. Tanjaroon Ly, B. Cohen-Stead, and S. Johnston, “Antiferromagnetic and bond-order-wave phases in the half-filled two-dimensional optical su-schrieffer-heeger-hubbard model”, *Phys. Rev. B*, vol. 111, p. 245138, 24 Jun. 2025. DOI: 10.1103/2bnf-tmtc.
- [9] P. Mai, B. Cohen-Stead, T. A. Maier, and S. Johnston, “Fluctuating charge-density-wave correlations in the three-band hubbard model”, *Proceedings of the National Academy of Sciences*, vol. 121, no. 50, e2408717121, 2024.
- [10] J. Neuhaus, N. S. Nichols, D. Banerjee, B. Cohen-Stead, T. Maier, A. Del Maestro, and S. Johnston, “Smoqydeac.jl: A differential evolution package for the analytic continuation of imaginary time correlation functions”, *SciPost Physics Codebases*, p. 039, 2024.
- [11] S. Malkaruge Costa, B. Cohen-Stead, and S. Johnston, “Kekulé valence bond order in the honeycomb lattice optical su-schrieffer-heeger model and its relevance to graphene”, *Phys. Rev. B*, vol. 110, p. 115130, 11 Sep. 2024. DOI: 10.1103/PhysRevB.110.115130.
- [12] S. Malkaruge Costa, B. Cohen-Stead, A. T. Ly, J. Neuhaus, and S. Johnston, “Comparative determinant quantum monte carlo study of the acoustic and optical variants of the su-schrieffer-heeger model”, *Phys. Rev. B*, vol. 108, p. 165138, 16 Oct. 2023. DOI: 10.1103/PhysRevB.108.165138.
- [13] S. Karakuzu, B. Cohen-Stead, C. D. Batista, S. Johnston, and K. Barros, “Flexible class of exact hubbard-stratonovich transformations”, *Phys. Rev. E*, vol. 107, p. 055301, 5 May 2023. DOI: 10.1103/PhysRevE.107.055301.
- [14] O. Bradley, B. Cohen-Stead, S. Johnston, K. Barros, and R. T. Scalettar, “Charge order in the kagome lattice Holstein model: A hybrid Monte Carlo study”, *npj Quantum Materials*, vol. 8, no. 1, p. 21, 2023. DOI: 10.1038/s41535-023-00553-y.
- [15] P. M. Dee, B. Cohen-Stead, S. Johnston, and P. J. Hirschfeld, “Charge correlations suppress unconventional pairing in the holstein model”, *Phys. Rev. B*, vol. 107, p. 104503, 10 Mar. 2023. DOI: 10.1103/PhysRevB.107.104503.
- [16] C. Miles, B. Cohen-Stead, O. Bradley, S. Johnston, R. Scalettar, and K. Barros, “Dynamical tuning of the chemical potential to achieve a target particle number in grand canonical Monte Carlo simulations”, *Phys. Rev. E*, vol. 105, p. 045311, 4 Apr. 2022. DOI: 10.1103/PhysRevE.105.045311.
- [17] G. Paleari, F. Hébert, B. Cohen-Stead, K. Barros, R. Scalettar, and G. G. Batrouni, “Quantum monte carlo study of an anharmonic holstein model”, *Phys. Rev. B*, vol. 103, p. 195117, 19 May 2021. DOI: 10.1103/PhysRevB.103.195117.