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### Giuseppe Barbalinardo

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**location**: Berkeley, CA

### Skills

**Technologies**

* Python (Numpy, Tensorflow, Scikit Learn, Pandas, Matplotlib, PySpark, MPI4py)
* DB (PostgreSQL / MySQL)
* Infrastructure (Docker / Kubernetes / Google Cloud)
* Mobile (Objective C, Swift, Android)
* Others (HPC / MPI / CUDA / Linux / JAVA SE)
* Software Development (Design Patterns / Algorithms)
* Advanced Math Tools (Statistics / Probability / Linear Algebra / Stochastic Methods / Econophysics)

**Machine Learning and AI**

* Optimization methods
* Dimensionality reduction
* Regularization
* Neural Networks
* Clusterization Algorithms
* NLP (Sentiment Analysis, Entity extraction, Transformer/BERT)
* Time series analysis (ARMA, ARIMA, LSTM, Wavenet)

**Experience**

**Ph. D. Researcher – University of California, Davis Sept ‘16 - Dec ‘20**

* Lead developer and creator of [kALDo](https://nanotheorygroup.github.io/kaldo), a modern Tensorflow-based open-source software package for heat transport simulations, which can run on CPUs and GPUs. Development from conception to release, including unit-tests, documentation, continuous integration, and deployment using Docker and CircleCI. Implementation of Google Colab examples. Team of 5 developers. Paper published in the Journal of Applied Physics (2020). Advisor: Dr. D. Donadio.
* Development of a novel mathematical model for heat transport, which uses advanced statistical tools to reduce the computational cost of large-scale simulations. Simulations ran on HPC. Published on Nature Communication (2019).
* Application of Neural Networks to model interatomic potentials. Journal of Applied Physics (2019).
* Teaching Assistant for the graduate class of Mathematical Methods for Scientists. Numerical algorithms using Python, including: Optimization Methods (Descent and BFGS), Partial Differential Equations, and Regularization, Dimensionality Reduction, Penalized Regressions, LASSO, Fast Fourier Transform, and Linear Algebra.
* Recipient of the Software Development Investment Fellowship from the National Science Foundation - Molecular Sciences Software Institute.
* Recipient of the 2020 Peter A. Rock Graduate Fellowship in Chemical Physics by UC Davis.

**Software Developer and Engineering Manager – Grio, San Francisco May ‘14 – Aug ‘16**

***Engineering Manager Jan ‘16 - Aug ‘16***

* Managed a team of 6-8 software developers across several simultaneous projects, while continuing hands-on coding.
* Designed and implemented the apprentice program and mentored junior developers.
* Organized the company’s first hackathon.

***Software Developer May ‘14 – Aug ‘16***

* Contributed to projects with diverse technology stacks, coding in Java, Objective C, SQL, and Python.
* Developed the new Target iPad app in an Agile-driven team of 12 people.
* Collaborated with the marketing and business team at Twitter. Developed an AngularJS dashboard to convert proprietary meta-language to Ruby and later application to over 10 marketing campaigns.
* Presented 4 Tech Talks at the company all-hands meeting.

### Projects

Co-founder of [Ergo](https://www.searchergo.com/) (June 2019), an AI-powered dashboard that pulls the latest news stories across media sources and highlights relevant content to combat the spread of misinformation.

* Implementation of the main machine learning algorithms, including Sentence Transformer (SBERT), Dimensionality reduction using PCA, Entity Extraction, and Clusterization, using Python, Numpy, Tensorflow and Pytorch.
* Development of the main stack, using Postgres and Cube, VueJS, Grafana, Docker, Kubernetes, and Google App Cloud.

### Education

**Ph.D. Computational Chemical Physics - University of California, Davis Sept ‘16 - Dec ‘20**

Relevant coursework: Artificial Intelligence / Natural Language Processing / Computational Methods / Statistical Mechanics

**M.Sc. Theoretical Physics - University of California, San Diego Apr ‘12 – Dec ’13**

Relevant coursework: Linear Algebra / Stochastic Methods / Differential Equations / Markov Chain Monte Carlo