

Adrian Domínguez-Castro, PhD

Data Scientist | Machine Learning | Data Analysis

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SUMMARY

Data Scientist with a PhD in Physics and 10 years of experience. Advanced knowledge in machine learning and statistical analysis. Skilled in Python and SQL. Proven track record of optimizing and implementing computational simulations and data-driven solutions for science and technology.

SKILLS

• Python • Machine Learning • Deep Learning • Data Science • SQL • Statistical Analysis

WORK EXPERIENCE

Postdoctoral Research Fellow

Vanderbilt University

Feb 2023-Jan 2024

Nashville, TN, USA

- Designed, developed and implemented a machine learning system that predicts physical-chemistry observables with more than 85% accuracy, resulting in 1 publication.
- Managed simulations and data science projects related to Computational Chemistry and Machine Learning.

Doctoral Research Assistant

University of Bremen

Oct 2018-Dec 2022

Bremen, Germany

- Completed 3 multi-year research projects, resulting in 4 publications and improved data science solutions resulting in a 50% reduction in data acquisition and processing time.
- Optimized data modeling in developing advanced simulations and predictive models for new technologies in nanoscience, and communicated insights in international scientific meetings.
- Built teams and forged collaborations across five countries (Germany, Spain, China, USA, France)

Graduate Research Fellow

University of Groningen

Sept 2017-Jul 2018

Groningen, The Netherlands

- Managed research on Theoretical Chemistry, and mentored junior team members.
- Optimized code for simulations and data analysis.

Visiting Research Scholar

Thompson Rivers University

Jan 2016-May 2016

Kamloops, Canada

- Developed computational modeling, data science and programming for Quantum Chemistry.
- Built teams and forged collaborations across two countries (Canada and Mexico), resulting in 1 publication.

Associate Junior Researcher

Higher Institute for Technology and Applied Sciences (InSTEC)

Sept 2015-Jul 2017

Havana, Cuba

- Led a team to meet project milestones and optimized computational modeling performance to automate preprocessing, reducing time taken by 50%.
- Developed and contributed to advanced predictive models for Computational Chemistry, resulting in 2 publications.

Assistant Professor

Technological University of Havana (CUJAE)

Sept 2014-Jul 2015

Havana, Cuba

- Delivered Lectures and Tutorials of Calculus, Differential Equations, and Numerical Mathematics (1st and 2nd years Electrical Engineering)

EDUCATION

Ph.D. (Dr. rer. nat.)

University of Bremen

Dissertation: Theoretical investigation on light- driven ultrafast dynamics in nanoscale

(Grade: magna cum laude)

Oct 2018 - Dec 2022

Bremen, Germany

B.Sc. Radiochemistry

Higher Institute for Technology and Applied Sciences (InSTEC)

Dissertation: Essential amino acids interacting with Fumonisin B1: A Theoretical Approach

(Grade: summa cum laude)

Sept 2008 - Jun 2014

Havana, Cuba

PORTFOLIO FOR DATA SCIENCE

Explore my data science projects in detail at <https://github.com/DrAdrianDC/Portfolio-for-Data-Science>.

CERTIFICATIONS

- Machine Learning by Stanford University & DeepLearning.AI on Coursera.
- AI for Everyone by Stanford University & DeepLearning.AI on Coursera.

PUBLICATIONS

Adrian Dominguez-Castro. *DFT and machine learning for predicting hydrogen adsorption energies on rocksalt complex oxides.* **Theor Chem Acc** **143**, **50** (2024) doi:[10.1007/s00214-024-03124-x](https://doi.org/10.1007/s00214-024-03124-x)

Michael Wagstaffe, **Adrian Dominguez-Castro**, Lukas Wenthaus, Steffen Palutke, Dmytro Kutnyakhov, Michael Heber, Federico Pressacco, Siarhei Dziarzhytski, Helena Gleißner, Verena Kristin Gupta, Harald Redlin, Adriel Dominguez, Thomas Frauenheim, Angel Rubio, Andreas Stierle, and Heshmat Noei. *Photoinduced Dynamics at the Water/TiO₂ (101) Interface.* **Phys. Rev. Lett.** **130**, **108001** (2023) doi:[10.1103/PhysRevLett.130.108001](https://doi.org/10.1103/PhysRevLett.130.108001)

Adrian Domínguez-Castro and Thomas Frauenheim. *Impact of vibronic coupling effects on light-driven charge transfer in pyrene-functionalized middle and large-sized metalloid gold nanoclusters from Ehrenfest dynamics.* **Phys. Chem. Chem. Phys.**, **2021**, **23**, **17129–17133** doi:[10.1039/d1cp02890a](https://doi.org/10.1039/d1cp02890a)

Adrian Domínguez-Castro, Carlos R. Lien-Medrano, Khaoula Maghrebi, Sabri Messaoudi, Thomas Frauenheim and Arnaud Fihey. *Photoinduced charge-transfer in chromophore- labeled gold nanoclusters: quantum evidence of the critical role of ligands and vibronic couplings.* **Nanoscale**, **2021**, **13**, **6786–6797** doi:[10.1039/d1nr00213a](https://doi.org/10.1039/d1nr00213a)

Michael Wagstaffe, Lukas Wenthaus, **Adrian Dominguez-Castro**, Simon Chung, Guilherme Dalla Lana Semione, Steffen Palutke, Giuseppe Mercurio, Siarhei Dziarzhytski, Harald Redlin, Nicolai Klemke, Yudong Yang, Thomas Frauenheim, Adriel Dominguez, Franz Kärtner, Angel Rubio, Wilfried Wurth, Andreas Stierle, and Heshmat Noei. *Ultrafast Real-Time Dynamics of CO Oxidation over an Oxide Photocatalyst.* **ACS Catal.** **2020**, **10**, **13650–13658** doi:[10.1021/acscatal.0c04098](https://doi.org/10.1021/acscatal.0c04098)

Adrian Domínguez-Castro, Daniel Hernández, and Fernando Guzmán. *Insights into the interactions of biomolecules with small gold clusters: a theoretical study from a DFTB perspective.* **Theor. Chem. Acc.** (2017) **136**:84 doi:[10.1007/s00214-017-2118-7](https://doi.org/10.1007/s00214-017-2118-7)

Adrian Dominguez-Castro, Fernando Guzman, Yasser Novo-Fernandez. *Adsorption on a nanoporous organic polymer for clean energy applications: A multiscale modelling study using Density Functional Tight Binding Approach.* **Computational and Theoretical Chemistry** **1102** (2017) **30–37** doi:[10.1016/j.comptc.2016.12.037](https://doi.org/10.1016/j.comptc.2016.12.037)

R. Castañeda-Arriaga, **A. Domínguez-Castro**, J. Lee, J. R. Alvarez-Idaboy, N. Mora-Diez. *Chemical Repair of Protein Carbon-Centred Radicals: Long-Distance Dynamic Factors.* **Can. J. Chem.**, **2016**, **94** (12): **1119–1126** doi:[10.1139/cjc-2016-0230](https://doi.org/10.1139/cjc-2016-0230).