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Ph.D in Physics, works with computational physics/chemistry and developing scientific software. My research interests are in the development of new methods for molecular simulation and the use of molecular simulation and quantum mechanics calculations to solve problems in physical chemistry, biochemistry and materials science.

Education and positions

- 2018- **Postdoctoral researcher, Institute of Physics, University of São Paulo**
Development and applications of Configurational Bias Monte Carlo Method in dyes of interest to organic solar cells
- 2015-2018 CAPES Fellowship - Supervisor: Dr. Kaline R. Coutinho
Ph.D. in Physics, Institute of Physics, University of São Paulo
Implementation and Development of Efficient Algorithm for Intramolecular Deformation with the Monte Carlo Method
- 2013-2015 CNPq Ph.D. Fellowship - Advisor: Dr. Kaline R. Coutinho
Master's in Computational Physics, São Carlos Institute of Physics, University of São Paulo
Implementation of the Parallel Tempering Monte Carlo Method to the Study of Thermodynamic Properties of Nanoclusters
- 2009-2012 FAPESP Master Fellowship - Advisor: Dr. Juarez L. F. Da Silva
B.S. in Computational Physics, São Carlos Institute of Physics, University of São Paulo
Undergraduate research entitled: Fluid Dynamics in Porous Media with Lattice Boltzmann
CNPq Scientific Initiation Fellowship - Advisor: Dr. Leonardo P. Maia

Complementary Education

- 2018 **Short Course:** I GPU Computing Workshop (Duration: 11h). University of São Paulo.
2016 **Short Course:** 4th Workshop HPC - USP/Rice (Duration: 8h). University of São Paulo.
2015 **Coursera Online Course:** Algorithms: Design and Analysis, Part 1. (Duration: 36h). Stanford University.
2006-2008 **Technical Course:** Informatics. Colégio Divino Salvador.

Awards

- 2017 **Best poster** award at the XIX Simpósio Brasileiro de Química Teórica.
2015 **Honorable mention** at the best poster award of the II Workshop on Biomolecular Theory-Experiment Interplay.

Teaching Experience

- 2019 **Física I:** Teaching the introductory physics class for the Mechanical Engineering course at the Escola Politécnica of the University of São Paulo.
2017 **Física Moderna I:** Teaching assistant for the introductory modern physics class of the Physics course of University of São Paulo, under the supervision of Prof. Dr. Sylvio Canuto.

Research overview

My research is mainly focused on the development of algorithms and computer code to efficiently sample properties of systems with intricate potential energy surfaces. Specifically, I have experience implementing and using Monte Carlo methods and extended ensemble methods such as replica exchange. During my Master I have implemented a parallel tempering Monte Carlo algorithm with several options in GOTNano, and have used this same software, together with FHI-aims, to investigate the properties of nanoclusters. In my PhD I developed a Configurational Bias Monte Carlo method for the efficient simulation of flexible molecules in solution, showing that the method may perform better than standard molecular dynamics for the sampling of systems with high energy barriers between conformers.

Even though most of my research has been in developing and implementing computational methods, I also have interest in combining different techniques to solve problems in physical chemistry and materials science. In my career I have used different methods such as molecular dynamics for the simulation of molecular systems, global optimization algorithms to search for the putative global minimum configurations, clustering algorithms to analyze data from simulation and density functional theory to investigate the electronic structure and absorption spectra of molecules and transition metal nanoclusters. All these methods form a useful toolset that can be used to investigate matter in the solid, liquid and gas phases.

Computational Skills

I am proficient with the programming languages (from more to less proficient): Fortran (legacy and modern), C, Python, C++ and Bash. I have experience with OpenMP parallelization, code profiling and optimization, and also have knowledge in message passing parallelization with MPI and GPU and shared memory parallelization with OpenACC. I know how to use the Git and SVN version control systems.

Concerning scientific software, I am familiar with DICE, GOTNano and GROMACS. For electronic structure calculations I have used Gaussian and FHI-AIMS.

Software development

Apart from several scripts and smaller software, my major contributions are developing:

Cezar HM, Canuto S, Coutinho K
DICE

A Monte Carlo code for the molecular simulation of liquids and gases using classical force fields. My main contribution is the coding and development of a Configurational Bias Monte Carlo method used to sample the internal degrees of freedom of the molecules. The code is written in Fortran and has OpenMP parallelization.

Rondina GG, Cezar HM, Da Silva JLF
GOTNano

A code for global optimization and sampling of thermodynamic properties of nanoclusters. My main contribution is the implementation of the Parallel Tempering Monte Carlo algorithm used to sample the thermodynamic properties. The code is written in C and has OpenMP parallelization.

Cezar HM

Clustering Trajectory

<https://github.com/hmcezar/clustering-traj>

A parallel Python script that performs clustering over an atomistic molecular simulation trajectory, searching for the minimum RMSD between each structure and using one of the many clustering algorithms available in the machine learning library sklearn.

Cezar HM

DICEtools

<https://github.com/hmcezar/dicetools>

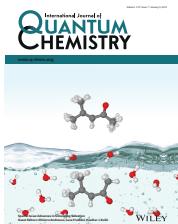
A package containing several scripts used to prepare DICE inputs and analyze simulation data.

Publications

h-index: 2, citations: 15 (16 Apr 2019)

[Publons](#), [Google Scholar](#)

Published

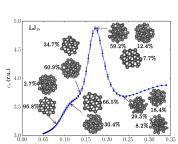


Cezar HM, Canuto S, Coutinho K

Solvent effect on the syn/anti conformational stability: A comparison between conformational bias Monte Carlo and molecular dynamics methods

International Journal of Quantum Chemistry 119:e25688, 2019 · DOI

We compare the Configurational Bias Monte Carlo (CBMC) and molecular dynamics (MD) sampling of the mesityloxide molecule in gas phase and solution, showing that while CBMC performs an ergodic sampling, MD has troubles overcoming the 10 kcal/mol energy barrier between the syn and anti conformations.

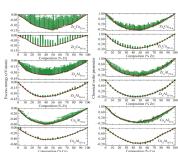


Cezar HM, Rondina GG, Da Silva JLF

Parallel Tempering Monte Carlo Combined with Clustering Euclidean Metrics Analysis to Study the Thermodynamic Stability of Lennard-Jones Nanoclusters

Journal of Chemical Physics 146:064114, 2017 · DOI

We use a clustering algorithm to analyze trajectories from Parallel Tempering Monte Carlo simulations of Lennard-Jones nanoclusters, identifying the phase changes and most frequent structures with great accuracy using a very simple approach.



De Souza DG, Cezar HM, Rondina GG, De Oliveira MF, Da Silva JLF

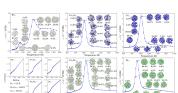
A Basin-hopping Monte Carlo Investigation of the Structural and Energetic Properties of 55- and 561-atom Bimetallic Nanoclusters: the Examples of the ZrCu, ZrAl, and CuAl Systems

Journal of Physics: Condensed Matter 28:175302, 2016 · DOI

We report a basin-hopping Monte Carlo investigation of the structural and energetic properties of bimetallic ZrCu, ZrAl, and CuAl nanoclusters with 55 and 561 atoms, showing how the atoms of the different species are distributed in the structure, and a trend of more spherical structures at 50-50 compositions.

In preparation

I have 4 manuscripts in different steps of preparation, of which I highlight 2 that are under the review of the coauthors and should be submitted in the next weeks.

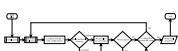


Cezar HM, Rondina GG, Da Silva JLF

Thermodynamic properties of 55-atom Pt-based nanoalloys: phase changes and structural effect on electronic properties

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We investigate the effects of temperature in Pt-based nanoalloys, performing Parallel Tempering Monte Carlo simulations using many-body potentials and DFT calculations.



Cezar HM, Canuto S, Coutinho K

DICE v3.0: A Monte Carlo code for molecular simulation including Configuration Bias Monte Carlo method

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We describe DICE, a software that uses Monte Carlo methods to perform molecular simulation, focusing on solute-solvent systems. We introduce our implementation of the Configurational Bias Monte Carlo method that includes new methodological developments.

Invited and contributed talks

São Carlos, SP
Feb 2019

Monte Carlo Methods

II New Energies Innovation Center - Computational Material Science Division Workshop (Invited talk)

Brasília, DF
Sep 2018

Comparison Between Configurational Bias Monte Carlo and Molecular Dynamics Sampling for Sampling Conformational Stability with High Energy Barriers

VII SeedMol

Livorno, Italy
Aug 2017

Configurational Bias Monte Carlo of Molecules in Solvent and Comparison with Molecular Dynamics Coding Solvation Workshop

Recife, PE
May 2017

Comparsion of the sampling of Molecular Dynamics and Configurational Bias Monte Carlo for 1,2- dichloroethane and octane.
III Advanced School on Biomolecular Simulation

Rio de Janeiro, RJ
Mar 2017

How Indiana Jones is actually a bad explorer when it comes to molecules and how I can do better
Finals of My Thesis in 180 seconds (Swissnex Brazil)

João Pessoa, PB
Sep 2014

Parallel Tempering Monte Carlo Investigation of Phase-Changes in Nanoclusters
XIII Encontro da SBPMat

Poster presentation

São Carlos, SP
Feb 2019

Application of Configurational Bias Monte Carlo method for the simulation of dyes and porous nanoparticles
II New Energies Innovation Center - Computational Material Science Division Workshop

São Carlos, SP
Oct 2018

Implementation of Configurational Bias Monte Carlo method for the simulation of flexible molecules in solution
I New Energies Innovation Center - Computational Material Science Division Workshop

Águas de Lindóia, SP
Nov 2017

Implementation of Configurational Bias Monte Carlo Method to Sample Flexible Solute in Solvent Media
XIX SBQT

Munich, Germany
Aug 2017

Configurational Bias Monte Carlo method to sample molecular flexibility: The case of octane and 1,2-dichloroethane
WATOC 2017

Maresias, SP
Aug 2015

Parallel Tempering Monte Carlo applied to the study of transition-metal nanoclusters
II Workshop on Biomolecular Theory-Experiment Interplay

Trieste, Italy
Aug 2013

Thermodynamic properties of nanoclusters: An investigation with Parallel Tempering Monte Carlo
Density Functional Theory and Beyond: Computational Materials Science for Real Materials

São Paulo, SP
Oct 2012

Estudo de dinâmica de fluidos em meios porosos com métodos de Boltzmann na rede
XX SIICUSP

Conferences organized

Recife, PE
May 2017

III Advanced School on Biomolecular Simulation
Member of the local committee.