

# RORY DONOVAN

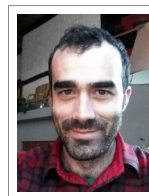
412 Rebecca Ave Ave.

Pittsburgh, PA 15221

☎ (206) 681 5515

✉ donovanr@alumni.reed.edu

🌐 [www.andrew.cmu.edu/user/rdonovan/](http://www.andrew.cmu.edu/user/rdonovan/)



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## EDUCATION

- 2011– **Ph.D., Carnegie Mellon–University of Pittsburgh Program in Computational Biology.**  
Systems Biology & Machine Learning
- 2004–2005 **M.S., University of Washington.**  
Physics
- 2000–2004 **B.A., Reed College.**  
Physics

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## CURRENT RESEARCH

### Network Inference with Graphical Models of Heterogeneous Genomic and Clinical Data.

Our models learn interaction networks between genes, clinical factors, and disease diagnoses, accommodate data that is both continuous and discrete, and are aggressively filtered for false-positive edges via collider detection algorithms.

### Graphical Models for Free Energy Estimation.

Computationally estimating how strongly two biomolecules bind together is often either overly time consuming (e.g. molecular dynamics) or overly empirical (e.g. docking). Alternatively, using graphical models of proteins to compute the Bethe free energy of binding can be both fast and accurate.

### Weighted Ensemble Systems Biology.

By applying weighted ensemble sampling to stochastic models of biological processes, we achieved orders of magnitude speed-ups in simulating events of interest in these complex systems. Our first paper dealt with non-spatial models, and our second paper addresses models of spatially resolved cellular processes.

### Quantitative Evolution and the ATP Synthase.

Using simple state-based models of proton transport and free energy transduction, we probe the optimality of the curiously engineered rotary mechanism of the ATP synthase, using models that are entirely agnostic to structure and are optimized over all unknown parameters that are thermodynamically permissible.

### Stochastic Models of Cellular Heterogeneity.

Simple stochastic models can recapitulate the population-level heterogeneity of protein abundance found in, for example, colonies of *E. coli*. This work employs a non-spatial model of gene expression, stochastically simulated with a modified Gillespie algorithm that takes into account cell division.

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## COMPUTATIONAL SKILLS

**Languages, Python, Julia, R, Matlab, Mathematica, Fortran, C, C++, Bash, HTML & CSS,  $\text{\LaTeX}$ .**

Scientific programming with a focus on simulation algorithms, data analysis, computational efficiency, and machine learning. Strong abilities and interests in:

- Expertise in Python, Matlab, R, and Julia, with additional proficiency in lower-level compiled languages.
- Coding best practices, including version control, maintainability, flexibility, and scalability.
- Machine learning and the automated extraction of knowledge from data.
- Run-time optimization (algorithms and data structures).
- Web design and the visual display of quantitative information.

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## EXPERIENCE

- 2012–2015 **Research Assistant, Carnegie Mellon–University of Pittsburgh Program in Computational Biology.**  
Working with Prof. Daniel Zuckerman, applying the weighted ensemble formalism to systems biology models.
- 2012–2015 **Teaching Assistant, Carnegie Mellon–University of Pittsburgh Program in Computational Biology.**  
Graduate-level computational structural biology.
- 2004–2006 **Research Assistant, University of Washington.**  
Worked with Prof. John Rehr, developing *ab initio* methods for computing Compton scattering profiles.
- 2005–2006 **Teaching Assistant, University of Washington.**  
Introductory and advanced undergraduate physics curriculum.

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## PUBLICATIONS

Rory M. Donovan, Jose-Juan Tapia, Devin P. Sullivan, James R. Faeder, Robert F. Murphy, Markus Dittrich, and Daniel M. Zuckerman. Unbiased rare event sampling in spatial stochastic systems biology models using a weighted ensemble of trajectories. *PLoS Computational Biology*, 2016.

Andrew J. Sedgewick, Ivy Shi, Rory M. Donovan, and Panatiotis V. Benos. Learning mixed graphical models with separate sparsity parameters and stability-based model selection. *BMC Genomics (In Press)*, 2014.

Rory M. Donovan, Andrew J. Sedgewick, James R. Faeder, and Daniel M. Zuckerman. Efficient stochastic simulation of chemical kinetics networks using a weighted ensemble of trajectories. *The Journal of Chemical Physics*, 139(11):115105, 2013.

Rory M. Donovan. *Quantization of constrained systems*. Bachelor's thesis, Reed College, 2004.

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## POSTERS

- 2015 **q-bio**, *Rare Event Sampling in Spatial Stochastic Systems Biology Models Using A Weighted Ensemble Of Trajectories*.
- 2015 **Biophysical Society**, *Rare Event Sampling in Stochastic Systems Biology Models Using A Weighted Ensemble Of Trajectories*.
- 2014 **RECOMB**, *Accelerating systems biology computation: Rapid estimation of equilibrium and kinetic quantities via weighted ensemble sampling*.
- 2013 **q-bio**, *Rare Event Sampling of Chemical Kinetics Networks Using A Weighted Ensemble Of Trajectories*.

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## AWARDS

- 2015 **Best Poster**, ARC 2015, \$500 travel award.
- 2015 **Outstanding Student Research Accomplishment**, CMU-Pitt. Program in Computational Biology.
- 2013 **Best Poster**, CMU-Pitt. Program in Computational Biology, Program Retreat.
- 2012–2014 **NRSA Institutional Training Grant predoctoral fellowship**, CMU-Pitt. Program in Computational Biology.
- 2004–2005 **Graduate Fellow**, University of Washington.

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## MISCELLANEOUS

- 2015 **Web Design and Backend Server**, <http://weightedensemblizer.csb.pitt.edu>. Automates the setup of applying our Weighted Ensemble sampling method to existing MCell simulations.
- 2013–2015 **Student Senator**, CMU-Pitt. Program in Computational Biology. Formed and elected to the student government. Served with faculty on the admissions committee for the Ph.D. program.
- 2013 **Scholarship**, *q-bio Summer School*. Highly selective two-week graduate level summer school / workshop. Worked on decoding stochastic gene regulatory networks from data.