

# Christopher Maffeo

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<http://bionano.physics.illinois.edu/People/chris.html>

## Education

### University of Illinois

Ph.D., Physics ..... 8/2014  
M.S., Physics ..... 5/2009

### University of California: Santa Barbara

B.S., Physics with highest academic honors ..... 6/2006

## Experience

**Research Assistant** ..... **12/2006–8/2014**  
University of Illinois at Urbana–Champaign ..... Urbana, IL

### Theoretical and computational modeling of biological systems

- Quantitatively studied DNA/protein interactions using all-atom and coarse-grain molecular dynamics simulations and analytical modeling
- Created my own adaptation of the umbrella sampling and weighted histogram analysis methods to obtain estimates of standard binding free energies
- Fitted custom chemical rate equations to simulation data to extract kinetic constants
- Led discussions of theoretical and computational challenges with fellow students and postdocs
- Supervised and supported the research activities of incoming students
- Contributed to research proposals that successfully secured an average of ~10 M cpu hours (worth roughly 500,000 USD) annually 2007–2014

### Workshop Instruction

- Designed and taught coursework for the Center for the Physics of Living Cells Summer School for biophysics graduate students and postdocs taking introductory and advanced courses (2010, 2011, 2013, 2014)

**Teaching Assistant** ..... **8/2006–12/2007**  
University of Illinois at Urbana–Champaign ..... Urbana, IL

- Taught physics laboratory and discussion sections for both technical and non-technical students
- Received “Teaching Excellence” award

**Grader** ..... **1/2005–5/2006**  
University of California: Santa Barbara ..... Goleta, CA

- Tutored small groups of students weekly

**Summer researcher** ..... **6/2003–8/2003**  
University of California: Santa Barbara ..... Goleta, CA

- Assisted experimental group (repaired BNCs, prepared bone samples, produced AFM images)

## Technical Skills

### Common Software

Molecular Dynamics software (NAMD, VMD)  
Adobe CS Suite (Photoshop, Illustrator)  
MS Office Suite (PowerPoint, Word, Excel)  
Development tools (emacs, gdb, git, SVN)

### Computer Languages

TCL, BASH, Matlab ..... *expert*  
C/C++, Perl, GNU Make ..... *proficient*  
Python, Mathematica ..... *proficient*

## Computer Environments

Unix/Linux, Windows, Apple  
High-Performance Computing Systems

## Languages

English ..... *native*  
German ..... *intermediate*

## Analytics

Construct coarse-grained models  
Solve ordinary and partial differential equations  
Work with probability, Bayesian inference  
Perform error analysis, parameter estimation  
Select appropriate approximations

## Publications & Presentations

### Research articles

**Maffeo C**, Ngo T, Ha T, and Aksimentiev A. (2014) A Coarse-Grained Model of Unstructured Single-Stranded DNA Derived from Atomistic Simulation and Single-Molecule Experiment. *Journal of Chemical Theory and Computation*. doi:10.1021/ct500193u

**Maffeo C**, Luan B, and Aksimentiev A. (2012) End-to-end attraction of duplex DNA. *Nucleic Acids Research*, 40(9), 3812-21. doi:10.1093/nar/gkr1220

**Maffeo C**, Schöpflin R, Brutzer H, Stehr R, Aksimentiev A, Wedemann G, and Seidel R. (2010) DNA–DNA interactions in tight supercoils are described by a small effective charge density. *Physical Review Letters*, 105(15), 158101-5. doi:10.1103/PhysRevLett.105.158101

**Maffeo C**, and Aksimentiev A. (2009) Structure, dynamics, and ion conductance properties of the phospholamban pentamer. *Biophysical Journal*, 96, 4853-4865. doi:10.1016/j.bpj.2009.03.053

### Review articles

**Maffeo C**, Jejoong Yoo and Jeffrey Comer and David Wells and Binqun Luan and Aleksei Aksimentiev. (2014) Close encounters with DNA. *Journal of Physics: Condensed Matter*

**Maffeo C**, Bhattacharya S, Yoo J, Wells D, and Aksimentiev A. (2012) Modeling and simulation of ion channels. *Chemical Reviews*, 112(12), 6250-84. doi:10.1021/cr3002609

### Other publications

**Maffeo C**, Carr R, Aksimentiev A. (2010) Introduction to MD simulation of DNA–protein systems (tutorial). <http://bionano.physics.illinois.edu/Tutorials>

**Maffeo C**, and Aksimentiev A. (2010) Single molecule force measurements: Insights from molecular simulations. *Physics of Life Reviews*, 7(3), 353-354.

### Presentations

“Coarse-grained modeling of single-stranded DNA and single-stranded DNA binding protein SSB,” (talk) *CPLC Post-doc and Graduate Student Symposium*, Urbana, Illinois, Feb. 2011.

“Bound to amaze: Molecular dynamics simulations of SSB,” (talk) *CPLC Post-doc and Graduate Student Symposium*, Urbana, Illinois, Feb. 2012.

“Molecular mechanics of DNA unwinding from single-stranded DNA binding protein” (poster) *CPLC Summer School*, Urbana, Illinois, Jun. 2011.

“Bound to amaze: Molecular dynamics simulations of SSB,” (talk) *CPLC Post-doc and Graduate Student Symposium*, Urbana, Illinois, Feb. 2011.

“Side-by-side and end-to-end attraction of double-stranded DNA” (poster) *Biophysical Society Annual Meeting*, Boston, Massachusetts, Feb. 2009.

“Towards a model of the replisome: beta clamp and friends,” (talk) *CPLC Post-doc and Grad Student Symposium*, Urbana, Illinois, Feb. 2009.

“Wetting, ionic conductance and structural stability of the phospholamban pentamer” (poster) *Biophysical Society Annual Meeting*, Long Beach, California, Feb. 2008.

## References

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### **Dr. Ralf Seidel**

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