David R. Slochower

Curriculum Vitae

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Education and positions

2015- **Postdoctoral scholar**, University of California, San Diego

Advisor: Michael K. Gilson, M.D., Ph.D.

Topic: Biophysical modeling and computational chemistry

2014 **Instructor**, University of Pennsylvania

Course: "Molecular physiology and cellular engineering"

2007-2014 Ph.D. in Biochemistry and Molecular Biophysics, University of Pennsvylvania

Advisor: Paul A. Janmey, Ph.D.

Thesis: Multiscale simulations of phosphatidylinositol bisphosphate: understanding its

biological role through physical chemistry

2003-2007 A.B. cum laude in Physics with distinction, Kenyon College

Research: high energy nuclear imaging

Research interests

- Physical chemistry of biological membranes
- Nonequilibrium statistical mechanics of molecular motors
- New methods for computing binding free energies
- Improved simulation force fields based on open source science

Computational design of synthetic peptides

2006-2008 Coded aperture imaging

Advisor: Jeffery Saven, Ph.D. (University of Pennsylvania)

Previous research

2007

2017-	Force field development with the open force field consortium
	Open Force Field Group
2016-	Thermodynamics of host-guest molecular recognition
	Advisor: Michael K. Gilson, M.D., Ph.D. (University of California, San Diego)
2015-	Theory of molecular motors
	Advisor: Michael K. Gilson, M.D., Ph.D. (University of California, San Diego)
2014-2015	Simulations and docking of macrocycles
	Advisors: Ravi Radhakrishnan, Ph.D. (University of Pennsylvania) and Mark A. Lemmon,
	Ph.D. (Yale University)
2009-2014	Quantum, all-atom, and coarse-grained molecular dynamics of membranes
	Advisor: Paul A. Janmey, Ph.D. (University of Pennsylvania)
2008	Simulations of viral entry into cells
	Advisors: William DeGrado, Ph.D. (University of California, San Francisco) and Michael L.
	Klein, Ph.D. (Temple University)
2007	Experimental single molecule biophysics
	Advisor: Yale E. Goldman, M.D., Ph.D. (University of Pennsylvania)

Advisors: John Idoine, Ph.D. (Kenyon College), John Frangioni, M.D., Ph.D. (Harvard University), and Richard Lanza, Ph.D. (Massachusetts Institute of Technology)

2005 Analysis of protein hydration shells in simulations

Advisor: Matthias Buck, Ph.D. (Case Western Reserve University)

Peer-reviewed publications

1. Himmelstein DS, **Slochower DR**, Malladi VS, Greene CS, Gitter, A. "Open Collaborative Writing with Manubot" *Accepted with minor revisions at PLOS Computational Biology* 2019 · GitHub (code) · GitHub (manuscript) · Nature TechBlog



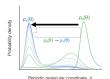
2. Bradley RP*, **Slochower DR***, Janmey PA, Radhakrishan R. "Molecular modeling of divalent cation-specific nano-clusters of phosphoinositides in physiologically-composed bilayers" *Under Review at JACS* 2018 · GitHub



3. Mobley DL, Bannan CC, Bayly CI, Rizzi A, Chodera JD, Lim VT, Lim NM, Beauchamp KA, **Slochower DR**, Shirts MR, Gilson MK, Eastman PK. "Escaping Atom Types in Force Fields Using Direct Chemical Perception" *Journal of Chemical Theory and Computation*, 14 6706-6092, 2018 · GitHub



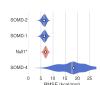
4. **Slochower DR**, Gilson MK. "Motor-like Properties of Nonmotor Enzymes" *Biophysical Journal* 114:9, 2018 · bioRxiv · DOI · GitHub · New and Notable · UCSD In the News



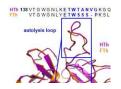
5. Yin J, Henriksen NM, **Slochower DR**, Gilson MK. "The SAMPL5 host-guest challenge: computing binding free energies and enthalpies from explicit solvent simulations by the attach-pull-release (APR) method" *Journal of Computer-Aided Molecular Design* 1:31 133-145, 2017 · DOI



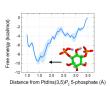
6. Yin J, Henriksen NM, **Slochower DR**, Shirts MR, Chiu MW, Mobley DL, Gilson MK. "Overview of the SAMPL5 host-guest challenge: Are we doing better?" *Journal of Computer-Aided Molecular Design* 1:31 1-19, 2017 · DOI · GitHub



7. Smith JR, Galie PA, **Slochower DR**, Weisshaar CL, Janmey PA, Winkelstein BA. "Salmon-derived thrombin inhibits development of chronic pain through an endothelial barrier protective mechanism dependent on APC" *Biomaterials* 80 96-105, 2016 · DOI



8. **Slochower DR**, Wang Y-H, Radhakrishnan R, Janmey PA. "Physical chemistry and membrane properties of two phosphatidylinositol bisphosphate isomers" *Physical Chemistry Chemical Physics* 17:19 12608-12615, 2015 · DOI



9. Wang Y-H, **Slochower DR**, Janmey PA. "Counterion-mediated cluster formation by polyphosphoinositides" *Chemistry and Physics of Lipids* 182 38-51, 2014 · DOI



10. **Slochower DR**, Wang Y-H, Tourdot RW, Radhakrishnan R, Janmey PA. "Counterion-mediated pattern formation in membranes containing anionic lipids" *Advances in Colloid and Interface Science* 208 177-188, 2014 · DOI



11. Janmey PA, **Slochower DR**, Wang Y-H, Wen Q, Ceber A. "Polyelectrolyte properties of filamentous biopolymers and their consequences in biological fluids" *Soft Matter* $10:10\ 1439-1449$, $2014\cdot DOI$



12. **Slochower DR**, Huwe PJ, Radhakrishnan R, Janmey PA. "Quantum and All-Atom Molecular Dynamics Simulations of Protonation and Divalent Ion Binding to Phosphatidylinositol 4,5-Bisphosphate (PIP₂)" *The Journal of Physical Chemistry B* 117:28 8322-8329, 2013 · DOI



13. Fujii, H, Idoine JD, Gioux S, Accorsi R, **Slochower DR**, Lanza R, Frangioni JV. "Optimization of Coded Aperture Radioscintigraphy for Sentinel Lymph Node Mapping" *Molecular Imaging and Biology* 14:2 173-182, 2012 · DOI



14. Donald JE, Zhang Y, Fiorin G, Carnevale V, **Slochower DR**, Gai F, Klein ML, De-Grado WF. "Transmembrane orientation and possible role of the fusogenic peptide from parainfluenza virus 5 (PIV5) in promoting fusion" *Proceedings of the National Academy of Sciences* 108:10 3958-3963, 2011 · DOI



15. Moravcevic K, Mendrola JM, Schmitz KR, Wang Y-H, **Slochower DR**, Janmey PA, Lemmon MA. "Kinase Associated-1 Domains Drive MARK/PAR1 Kinases to Membrane Targets by Binding Acidic Phospholipids" *Cell* 143:6 966-977, 2010 · DOI



Book chapters

1. **Slochower DR**, Wang Y-H, Radhakrishan R, Janmey PA. "Lipid membrane shape evolution and the actin cytoskeleton" in *Handbook of Lipid Membranes, Molecular and Materials Aspects*, Eds. Safinya C, Rädler, J. (2018)



Ph.D. thesis

Slochower DR. "Multiscale simulations of phosphatidylinositol bisphosphate: understanding its biological role through physical chemistry" *University of Pennsylvania*, 2014



Software packages and contributions

pAPRika	GitHub	Free energy calculations with AMBER and OpenMM
speakeasy	GitHub	Automates the conversion of SMIRNOFF parameters to AMBER force field files
smirnovert	GitHub	Convert host-guest systems into SMIRNOFF force fields
manubot	GitHub	Automated scholarly publishing
BioPhysCode	GitHub	Tools for building and analyzing membrane simulations

^{*} These authors contributed equally.

Software skills

- Proficient programming in Python and Jupyter Notebooks with knowledge of Bash, C++, FOR-TRAN, and Perl.
- Active user of GitHub and continuous integration platforms.
- Experience writing unit, integration, and regression tests.
- Expert with molecular dynamics simulations using AMBER and OpenMM; familiar with Gromacs, CHARMM, and LAMMPS.

Invited talks, posters, abstracts, and workshops (selected)

2019	Talk	Binding free energy calculations using the attach-pull-release method 2019 AMBER developers' meeting
2019	Talk	Benchmarking emerging force fields with binding thermodynamics calculations
		2019 University of California Chemical Symposium
2017	Talk	Directional motion in chiral molecules out of equilibrium
		253rd American Chemical Society Meeting
2017	Poster	Directional and driven motion in enzymes out of equilibrium
		61st Annual Biophysical Society Meeting
2014	Talk	Multiscale modeling of polyphosphoinositides
		University of California, San Diego
2014	Talk	Physical chemisty of phosphatidylinositol isomers
		University of California, Irvine
2013	Talk	Membranes: Polyphosphoinositides
		Friday Research Discussions, University of Pennsylvania
2013	Poster	Quantum and All-atom Molecular Dynamics Simulations of Proton Binding to
		Phosphatidylinositol 4,5-bisphosphate (PIP ₂)
		57th Biophysical Society Meeting
2012	Talk	Molecular Dynamics Simulations of Ion Binding and Protonation of Phos-
		phatidylinositol Bisphosphate (PIP ₂)
		244th American Chemical Society Meeting
2012	Talk	Simulations of membrane electrostatics with PtdInsP ₂
		George W. Raiziss 30th Annual Retreat
2012	Poster	Molecular Dynamics Simulations of Phosphatidylinositol Bisphosphate (PIP ₂)
		American Physical Society, March Meeting
2011	Talk	Molecular Dynamics Simulations of Membranes
		47th New England Complex Fluids Workshop
2011	Poster	Molecular Dynamics Simulations of Monolayers and Membranes with Phos-
		phatidylinositol Bisphosphate
• • • •	- 11	55th Biophysical Society Meeting
2010	Talk	Simulating highly charged monolayers
		Mechanistic Studies in Membrane Biophysics: Experiments and Theory, Tel-
2010	ъ.	luride Science and Research Workshop
2010	Poster	Simulations of Monolayers with Phosphatidylinositol Bisphosphate
		Gotham-Metro Condensed Matter Meeting

Awards and grants

2012-2013	NIH T32 Structural Biology Training Grant
2011	Juan Grana Graduate Teaching Assistantship
2010-2012	NIH T32 Interdisciplinary Cardiovascular Training Grant
2007	Distinction in Physics (best research), Kenyon College
2007	Sigma Xi, The Scientific Honor Society
2004-2007	Dean's List, Kenyon College
2005	Best Summer Project (biophysics), Case Western Reserve University
2001	Science Olympiad, National Champion Team

Teaching and mentoring experience

Fall 2014 Molecular Physiology & Cellular Engineering – University of Pennsylvania I was responsible for creating the syllabus, giving lectures, designing project assignments, and grading for one-half of this course.

Spring 2011 Macromolecular Biophysics II – University of Pennsylvania

I was in charge of arranging lectures, holding office hours and regular review sessions, grading homework and exams for first and second year graduate students.

2009- Mentored high school, undergraduate, and graduate students in research.

Spring 2005 Programming I – Kenyon College

I designed regression tests for weekly project assignments, graded, and then posted my own solutions to the class.

Service

- Reviewer for Soft Matter
- Reviewer for European Biophysics Journal
- Reviewer for Scientific Reports
- Reviewer for Nature Structural & Molecular Biology
- Member, Biophysical Society
- Member, American Chemical Society

References

Michael K. Gilson, M.D., Ph.D.
Professor and Chair in Computer-Aided Drug Design
Co-Director UC San Diego Center for Drug Discovery Innovation
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