

# Erik Nordquist

✉ [enordquist@rx.umaryland.edu](mailto:enordquist@rx.umaryland.edu)

🌐 [eriknordquist.com](http://eriknordquist.com)

20 Penn St, Baltimore, MD 21201

## Education

2023 **Ph.D.** Chemistry, University of Massachusetts Amherst. Advisor: Jianhan Chen, Ph.D.  
2018 **B.S.** Chemistry and Physics, The College of Idaho

## Appointments

2023– **Postdoctoral Fellowship**, University of Maryland, Baltimore, Department of Pharmaceutical Sciences. Advisor: Alexander D. MacKerell, Jr., Ph.D.

## Fellowships and Awards

2024 **Best of Biophysical Journal 2023**, publication titled “Inner pore hydration free energy...”  
Talk award, Research Symposium, UM Greenebaum Comprehensive Cancer Center  
2023– **NIH T32 Postdoctoral Fellowship**, University of Maryland, Baltimore  
2022 **Paul H. Terry Endowment Award**, Chemistry Dept., University of Massachusetts Amherst  
**Teaching Fellowship**, College of Natural Sciences, University of Massachusetts Amherst  
2020–22 **NIH T32 Predoctoral Fellowship**, University of Massachusetts Amherst  
2020 **William E. McEwen Poster Award**, Chemistry Dept., University of Massachusetts Amherst

## Peer-Reviewed Publications

- Patel R, **Nordquist E**, Polli J. Prediction of surfactant-mediated dissolution of poorly soluble drugs from drug particles. **Eur. J. Pharm. Sci.** 2025, 107052. DOI: [10.1016/j.ejps.2025.107052](https://doi.org/10.1016/j.ejps.2025.107052)
- **Nordquist E**<sup>#</sup>, Zhao M<sup>#</sup>, Kumar A, MacKerell A. Physics- and Machine-Learning Based Method to Identify Druggable Binding Sites Using SILCS-Hotspots. **J. Chem. Inf. Model.** 2024, 64, 19, 7743-7757. DOI: [10.1021/acs.jcim.4c01189](https://doi.org/10.1021/acs.jcim.4c01189) <sup>#</sup>Contributed equally.
- **Nordquist E**, Jia Z, Chen J. Small Molecule NS11021 Promotes BK Channel Activation by Increasing Inner Pore Hydration. **J. Chem. Inf. Model.** 2024, 64, 19, 7616-7625. DOI: [10.1021/acs.jcim.4c01012](https://doi.org/10.1021/acs.jcim.4c01012)
- **Nordquist E**<sup>#</sup>, Zhang G<sup>#</sup>, Barethiya S, Ji N, White K, Han L, Jia Z, Shi J, Cui J, and Chen J. Incorporating physics to overcome data scarcity in predictive modeling of protein function: a case study of BK channels. **PLOS Comput. Biol.** 2023 19(9): e1011460. DOI: [10.1371/journal.pcbi.1011460](https://doi.org/10.1371/journal.pcbi.1011460) <sup>#</sup>Contributed equally.
- Zhang L, Barethiya S, **Nordquist E**, Chen J. Machine Learning Generation of Dynamic Protein Conformational Ensembles. **Molecules** 2023, 28(10), 4047. DOI: [10.3390/molecules28104047](https://doi.org/10.3390/molecules28104047)
- **Nordquist E**, Zhiguang J, Chen J. Inner pore hydration free energy controls activations of the big potassium channel and its mutants. **Biophys. J.** 2023, 122, 1158-1167. DOI: [10.1016/j.bpj.2023.02.005](https://doi.org/10.1016/j.bpj.2023.02.005) Selected for Best of 2023 Edition.
- **Nordquist E**<sup>#</sup>, Schultz S<sup>#</sup>, and Chen J. Using Metadynamics To Explore the Free Energy of Dewetting in Biologically Relevant Nanopores. **J. Phys. Chem. B** 2022, 126, 34, 6428-6437 DOI: [10.1021/acs.jpcb.2c04157](https://doi.org/10.1021/acs.jpcb.2c04157) <sup>#</sup>Contributed equally.
- **Nordquist E**, Clerico E, Chen J, Gierasch L. Computational Modeling of Hsp70-Client Interactions: Past, Present, and Future. **J. Phys. Chem. B** 2022, 126, 36, 6780–6791 DOI: [10.1021/acs.jpcb.2c03806](https://doi.org/10.1021/acs.jpcb.2c03806)

- **Nordquist E**, English C, Clerico E, Sherman W, Gierasch L, Chen J. Physics-based modeling provides predictive understanding of selectively promiscuous substrate binding by Hsp70 chaperones. **PLOS Comput. Biol.** 2021, 17 (11): e1009567. DOI: [10.1371/journal.pcbi.1009567](https://doi.org/10.1371/journal.pcbi.1009567)
- Gong X, Chiricotto M, Liu X, **Nordquist E**, Feig M, Brooks CL, Chen J. Accelerating the generalized born with molecular volume and solvent accessible surface area implicit solvent model using graphics processing units. **J. Comput. Chem.** 2020, 41, 830–838. DOI: [10.1002/jcc.26133](https://doi.org/10.1002/jcc.26133)

## Teaching and Mentoring

---

- 2022 **Instructor of record**, First-Year Seminar, self-designed, title: “Reconciling Atomic Chaos and Human Order”. Funded by UMass College of Natural Sciences Teaching Fellowship. ([link](#))
- 2020–24 **Guest lectures**  
 Computer-aided Drug Design in UMB Graduate Cancer Biology course (2x);  
 Discussion seminar moderator on AlphaFold2 at Amherst College Biophysics course;  
 Lecture on molecular mechanics, additive force fields in UMass Graduate Stat. Mech. Course
- 2018–19 **Lab TA**, General Chemistry I Lab

## Mentoring Experience with fellow trainees:

**Undergraduates (UMass):** Samantha Schultz (2020-2021); Callie Jillson (2019-2020)

**Graduate students (UMB):** Anthony O'Donnell (2024–)

## Presentations

---

- 2025 **Talk**, Research Day, University of Maryland School of Pharmacy, “Computational modeling of ternary complexes as ensembles using SILCS-PROTAC,” Baltimore, MD.  
**Talk**, Chemistry Department Seminar, George Washington University, “Computational design of PROTACs”, Washington D.C.
- 2024 **Talk**, Institute for Bioscience and Biotechnology Research, University of Maryland Baltimore, Early-Career Research Symposium, “Computational design of PROTACs”, Rockville, MD.  
**Talk**, University of Maryland Greenbaum Comprehensive Cancer Center Research Symposium, “Physics- and machine-learning-based method for identifying druggable binding sites with SILCS-Hotspots.” (Talk award) Baltimore, MD.  
**Talk**, The College of Idaho Natural Science Symposium, “Computer simulations of proteins help understand their function.” Caldwell, ID.  
**Poster**, Biophysical Society Annual Meeting, “Computational mapping of allosteric modulators of the BK channel.” Philadelphia, PA.
- 2023 **Poster**, Biophysical Society Annual Meeting, “A predictive model of voltage gating of BK channels via physical modeling and machine learning.” San Diego, CA.
- 2022 **Talk**, University of Massachusetts Amherst ResearchFest (PH Terry award): “Predicting protein function with physics, experiments and machine learning.” Amherst, MA.  
**Poster**, Biophysical Society Annual Meeting, “Free energy of hydrophobic dewetting in gating of BK channels.” San Francisco, CA.
- 2020 **Talk**, Northeastern Structural Symposium, “Physical origins of selective promiscuity to Hsp70s revealed through physics-based modeling.” Virtual.  
**Poster**, University of Massachusetts Amherst ResearchFest (WE McEwen Award): “Physical origins of selective promiscuity to Hsp70s revealed through physics-based modeling.” Amherst, MA.
- 2019 **Talk**, Biophysics at University of Massachusetts Amherst, “Understanding the origins of DnaK’s selective promiscuity with physics-based modeling”

**Poster**, Molecular Biophysics in the Northeast, “Understanding the origins of DnaK’s selective promiscuity with physics-based modeling.” Boston, MA.

## Service

---

- 2025 **Facilitator**, Skills for Faculty Positions panel, University of Maryland Office of Postdoctoral Scholars  
**Moderator**, Career Panel Discussion, University of Maryland, Baltimore Cancer Retreat
- 2023–24 **Facilitator**, Responsible Conduct of Research NIH training, University of Maryland, Baltimore
- 2024– **Organization committee, Annual Cancer Research Retreat**, University of Maryland, Baltimore (2x)
- 2020–20 **Organization committee, Alumni Networking Symposium**, Chemistry-Biology Interface program, University of Massachusetts Amherst (2x)
- 2020 **Search committee**, Grad. Program Manager, Chemistry Dept. University of Massachusetts Amherst
- 2019–21 **ResearchFest committee** for Chemistry Dept., University of Massachusetts Amherst (3x)

## Journals Refereed for:

Comm. Chem., Biophys. J., J. Chem. Theory Comput., J. Chem. Inf. Model.

## Outreach

---

- 2023-25 **Volunteer instructor**, Baltimore Underground Science Space (BUGSS) ([link](#)), STEM outreach for Baltimore community members in community-accessible lab space
- Volunteer instructor**, RAMP Program for STEM activity for high-schoolers in Baltimore ([link](#)) (2x)
- Poster judge**, Mount Royal Middle School Science Fair, Baltimore (2x)
- Volunteer instructor**, CURE Program University of Maryland, Baltimore, STEM outreach for middle- / high-schoolers in Baltimore ([link](#)) (3x)
- Volunteer**, Halloween STEM outreach event at Community Engagement Center, University of Maryland, Baltimore
- 2020–23 **Reviewer for Journal of Emerging Investigators**, 25 articles by middle- / high-school students ([link](#))
- 2022 **Lab for girls summer science camp**, Eureka! at University of Massachusetts Amherst ([link](#))

## Professional Development

---

- 2025 **Professional Society Memberships:**  
Biophysical Society (2022-), American Chemical Society (2025-)
- 2024 **Writing an Effective Teaching Philosophy Statement**, online CIRTl workshop
- Teaching Biophysics at a PUI workshop**, Biophysical Society Annual Meeting
- 2024 **Safety preparedness trainings**, CPR/AED, Stop the Bleed, Civilian Active Shooter Events, and Fire Extinguishers, University of Maryland, Baltimore ([certificates](#))
- 2022 **CIRTl associate certification**, University of Massachusetts Amherst, achieved through training on evidence-based and inclusive teaching practices ([link](#))
- 2021 **Evidence-based Undergraduate STEM Teaching**, online course ([link](#))
- Inclusive STEM Teaching**, online course ([link](#))

## Book Chapters

---

- **Nordquist E**, Horrigan F, MacKerell A. Computational ligand binding site prediction. **Springer Nature**. (accepted) 2025.