

# JIATAO LIANG

**Postdoctoral Researcher**

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## EDUCATION

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<b>PhD</b>	Harvard University: Chemistry and Chemical Biology	Feb 2025
<b>BS</b>	University of California, Berkeley: Chemical Biology Highest Honors	May 2019
<b>BA</b>	University of California, Berkeley: Statistics Highest Honors	May 2019

## HONORS AND AWARDS

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<b>UC Berkeley Chemistry Department Citation</b>	2019
<i>Awarded in highest recognition of distinguished undergraduate work to a single graduating senior</i>	

## RESEARCH EXPERIENCE

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<b>University of Pennsylvania</b> , Philadelphia, PA	2025-current
Postdoctoral Researcher, Advisor: Yoseph Barash	
<ul style="list-style-type: none"><li>Implementation of generative machine learning models for the acceleration of directed evolution of non-canonical nucleic acid ribozymes</li><li>Application of computational bioinformatics on long non-coding RNAs to explore biological mechanism towards therapeutic applications</li><li>Consultation for a large language model using contrastive learning to extract key chemical features of small molecules and probes from scientific literature</li></ul>	

<b>Harvard University</b> , Cambridge, MA	2019-2025
Doctoral Candidate, Advisor: Matthew Shair	
Supported by Blavatnik Biomedical Accelerator and Deerfield Lab1636	
<ul style="list-style-type: none"><li>Independently designed and synthesized small molecule inhibitors of ADAR1, an RNA-editing enzyme implicated in cancer pathogenesis</li><li>Conceived and developed novel assays to quantify inhibition profiles of synthesized compounds</li><li>Interpreted assay results in context of <i>in silico</i> docking simulations to establish structure-activity relationships and inform future work</li><li>Collaborated with virtual biotech 3DC in directing contract research organizations in design, synthesis, and assay of key molecules to interrogate inhibition of ADAR1</li><li>Dissertation: Development of Inhibitors of ADAR1</li></ul>	

<b>Lawrence Berkeley National Laboratory</b> , Berkeley, CA	2017-2019
Undergraduate Research Assistant, Advisor: Yi Liu	
<ul style="list-style-type: none"><li>Explored the stability and reactivity of <i>para</i>-azaquinodimethane moiety in extended conjugated systems for use in organic electronic applications</li></ul>	

**Plexxikon, Inc.**, Berkeley, CA  
Medicinal Chemistry Summer Intern

2018

- Synthesized over forty compounds for the inhibition of kinases implicated in oncology
- Interpreted structural biology data and assay results to inform future work

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## TEACHING EXPERIENCE

**Harvard University**, Cambridge, MA Spring 2021  
Head Teaching Fellow

- Courses taught: Organic Chemistry of Life (CHEM27) with Prof. Emily Balskus
- Responsibilities include setting and implementing course policy, managing fifteen junior teaching fellows, interfacing with other course staff, giving review lectures, and holding office hours.

**Harvard University**, Cambridge, MA 2020  
Teaching Fellow

- Courses taught: Organic Chemistry of Life (CHEM 27) with Prof. Emily Balskus and Principles of Organic Chemistry (CHEM17) with Prof. Christina Woo
- Responsibilities include holding office hours, leading discussion sections, and proctoring exams.

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## PUBLICATIONS

Anderson, C.L., ... Liang, J., *et al.* **Solution-processable and functionalizable ultra-high molecular weight polymers via topochemical synthesis.** *Nat Commun* **12**, 6818 (2021).  
<https://doi.org/10.1038/s41467-021-27090-1>

Anderson, C.L., Liang, J., *et al.* **A highly substituted pyrazinophane generated from a quinoidal system: Via a cascade reaction.** *Chemical Commun* **56**, 4472–4475 (2020).  
<https://doi.org/10.1039/d0cc00916d>

Jones, H.T., ... Liang, J., *et al.* **A Dataset for Distilling Knowledge Priors from Literature for Scientific Discovery.** *NeurIPS.* (2025).  
<https://doi.org/10.48550/arXiv.2508.10899>

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## PRESENTATIONS

**Harvard Chemistry ChemTalks Series**  
• Speaker September 2021

**Harvard Chemistry and Chemical Biology Symposium**  
• Poster presentation March 2022, 2023

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## SKILLS AND EXPERIENCE

- Planning and executing multi-step chemical syntheses, with focus in heterocyclic systems
- Analysis of structural biology data on platforms such as Molecular Operating Environment (MOE) and Pymol, including generation of homology models based on both published structures and computationally generated AlphaFold models
- Coding in Python, Java, and R, including scripting workflows for processing, analyzing, and visualizing biological data from RNASeq