

Alexey STROKACH

PHD STUDENT • UNIVERSITY OF TORONTO • TORONTO, ON

✉ alex.strokach@utoronto.ca | ☎ (647) 226-4521 | 🌐 [ostrokach](https://ostrokach.github.io) | 📁 ostrokach.gitlab.io

EDUCATION

DEC. 2023 (Expected)	PhD in COMPUTER SCIENCE Supervisor: Dr. Philip M. KIM GPA: 3.93 / 4.0 Awards: • NSERC Postgraduate Scholarship (2018-2021, \$63000)	University of Toronto, Toronto, ON
DEC. 2016	MSc in COMPUTER SCIENCE Supervisor: Dr. Philip M. KIM GPA: 4.0 / 4.0	University of Toronto, Toronto, ON
AUG. 2013	MSc in PHARMACEUTICAL SCIENCES Supervisor: Dr. James W. WELLS GPA: 4.0 / 4.0 Awards: • Ontario Graduate Scholarship (2011-2012, \$15000)	University of Toronto, Toronto, ON
AUG. 2010	BMSc in Medical Sciences (Honors) GPA: 88.7 / 100 Awards: • Dean's Honor List (2006-2010) • CIS Academic All-Canadian in Cross Country and Track and Field (2009) • Queen Elisabeth II Scholarship (2006-2009, \$3500 x 4) • Reinhard Konrad Memorial Awards in Science (2008, \$1000) • Laurene Paterson Estate Scholarship (2007, \$2000) • The Western Scholarship of Distinction (2006, \$1000)	University of Western Ontario, London, ON

WORK EXPERIENCE

JAN. 2022– PRESENT	Software Developer Working on machine learning models for detecting fraud in advertising.	Google, Toronto, ON
MAY–AUG. 2021	Software Engineering Intern Implemented differentially private mechanisms and trained ML models using anonymized data.	Google, Remote, ON
MAY–AUG. 2019	Software Engineering Intern Designed and evaluated APIs for interacting with Apache Beam batch and streaming pipelines.	Google, Sunnyvale, CA
SEP. 2010– PRESENT	Teaching Assistant Developed content, held tutorials, graded assignments and exams for the following courses: <ul style="list-style-type: none">CSC413H – Neural Networks and Deep Learning 1 termCSC411H – Introduction to Machine Learning 1 termCSC384H – Introduction to Artificial Intelligence 3 termsCSC343H – Introduction to Databases 3 termsCSC311H – Introduction to Machine Learning 1 termCSC209H – Software Tools and Systems Programming 1 termCSC148H – Introduction to Computer Science 6 termsCSC128H – Content Developer for PythonTA 1 termCSC120H – Computer Science for the Sciences 1 termCSC108H – Introduction to Computer Programming 2 termsC4M – Computing for Medicine 1 termCS Help Center 1 termPHM140H / PHC301H – Molecular Pharmacology 2 termsPHM225H – Introduction to Biochemistry and Molecular Biology 1 term	University of Toronto, Toronto, ON
MAY–AUG. 2009	Summer Research Assistant Supervisors: Dr. David W. SHOESMITH and Dr. J. Clara WREN Performed computational modeling of gas-phase radiolysis and carbon steel corrosion.	University of Western Ontario, London, ON

MISCELLANEOUS

- Reviewer for the *Machine Learning in Computational Biology (MLCB)* workshop and the *Computational and Structural Biotechnology Journal*, *Biotechnology Letters* and *BMC Supplements* journals.
- Competed on Varsity Cross Country and Track and Field teams for five years.

SELECT RESEARCH PROJECTS

2018–PRESENT	Fast and Flexible Design of Novel Proteins (ProteinSolver) Supervisor: Dr. Philip M. KIM Technologies: PYTHON, PYTORCH, PYTORCH GEOMETRIC, VOILA, DOCKER, KUBERNETES <ul style="list-style-type: none">• Pre-trained a graph convolutional network to reconstruct masked amino acids in proteins.• Evaluated the quality of the learned embeddings using a variety of downstream applications.• Generated new protein sequences which underwent successful experimental validation.• Currently extending this approach to generate proteins with entirely “hand-drawn” topologies.
2017–2018	Predicting Synergistic and Beneficial Drug Combinations Supervisor: Dr. Philip M. KIM Technologies: PYTHON, SCIKIT-LEARN, XGBOOST, TENSORFLOW, MYSQL <ul style="list-style-type: none">• Created a data warehouse with information about drugs (rdkit), drug targets (STITCH), protein interactions (Mentha), protein functional associations (STRING), and cell line- and drug-dependent gene expression levels (LINCS), mutation profiles (CCLE), and essentiality scores (Achilles).• Compared decision tree- and neural network-based models for predicting drug combinations.• Participated in the AstraZeneca-Sanger Drug Combination Prediction DREAM Challenge.
2013–2016	Predicting the Effect of Mutations on a Genome-Wide Scale Supervisor: Dr. Philip M. KIM Technologies: PYTHON, SCIKIT-LEARN, XGBOOST, DJANGO, MYSQL, AIOHTTP <ul style="list-style-type: none">• Created > 100k homology models of proteins and protein-protein interactions.• Trained a gradient boosted decision tree model to predict the $\Delta\Delta G$ associated with mutations.• Performed feature elimination while optimizing performance on external validation tasks.• Implemented a web server backend which executes user-submitted jobs on a SLURM cluster.
2010–2013	Probing the Oligomeric Status of G Protein-Coupled Receptors using Förster Resonance Energy Transfer (FRET) and Single-Particle Tracking (SPT) Supervisor: Dr. James W. WELLS Technologies: SPT MICROSCOPY, FRET MICROSCOPY, MATLAB, CELL CULTURE, MOLECULAR CLONING, PCR <ul style="list-style-type: none">• Created DNA constructs for proteins tagged with fluorescent proteins or other labels.• Expressed fluorescently-tagged proteins in cells and collected thousands of microscopy images.• Derived 4 mathematical models describing expected patterns of FRET for different interactions.• Fit single particle fluorescence to Poisson distributions using custom CUDA kernels.

PUBLICATIONS

- Strokach A, Kim PM. *Deep generative modeling for protein design*. **Current Opinion in Structural Biology**. 2021 ([submitted](#)).
- Strokach A, Becerra D, Corbi-Verge C, Perez-Riba A, Kim PM. *Computational generation of proteins with predetermined three-dimensional shapes using ProteinSolver*. **STAR Protocols**. 2021; doi: [10.1016/j.xpro.2021.100505](#).
- Strokach A, Lu TY, Kim PM. *ELASPIC2 (EL2): Combining contextualized language models and graph neural networks to predict effects of mutations*. **Journal of Molecular Biology**. 2021; doi: [10.1016/j.jmb.2021.166810](#).
- Strokach A, Becerra D, Corbi-Verge C, Perez-Riba A, Kim PM. *Fast and flexible protein design using deep graph neural networks*. **Cell Systems**. 2020; 11: 1–10. doi: [10.1016/j.cels.2020.08.016](#).
- Strokach A, Corbi-Verge C, Kim PM. *Predicting changes in protein stability caused by mutations using sequence- and structure-based methods in a CAGI5 blind challenge*. **Human Mutation**. 2019; 40: 1414–1423. doi: [10.1002/humu.23852](#).
- Savojardo C, Petrosino M, Babbi G, Bovo S, Corbi-Verge C, et al.. *Evaluating the predictions of the protein stability change upon single amino acid substitutions for the FXN CAGI5 challenge*. **Human Mutation**. 2019; 0. doi: [10.1002/humu.23843](#).
- Strokach A, Corbi-Verge C, Teyra J, Kim PM. *Predicting the Effect of Mutations on Protein Folding and Protein-Protein Interactions*. **Methods in Molecular Biology**. 2019; 1851: 1–17. doi: [10.1007/978-1-4939-8736-8_1](#).
- Witvliet D*, Strokach A*, Giraldo-Forero AF, Teyra J, Colak R, and Kim PM. *ELASPIC web-server: proteome-wide structure based prediction of mutation effects on protein stability and binding affinity*. **Bioinformatics** 32 (10): 1589–1591. doi: [10.1093/bioinformatics/btw031](#).

CONFERENCE POSTERS

- Strokach A, Becerra D, Corbi-Verge C, Perez-Riba A, Kim PM *Designing real novel proteins using deep graph neural networks*. Machine Learning in Computational Biology (MLCB); **NeurIPS Workshop on Learning Meaningful Representation of Life (LMRL)**. Vancouver, Canada, 2019.
- Strokach A, Wells JW. *Using Förster resonance energy transfer to distinguish between proteins that exist as monomers, transient oligomers and stable oligomers in the plasma membrane of living cells*. **Chemical Biophysics Symposium**. Toronto, Canada, 2013.
- Strokach A, Wells JW. *Probing the Oligomeric Status of G Protein-Coupled Receptors by Single-Molecule Fluorescence*. **Great Lakes GPCR Retreat**. London, Canada, 2012.