# Edwin Chacko

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## EDUCATION

### University of Toronto

Sep. 2022 – May 2027

BASc. in Engineering Science, Machine Intelligence Option.

Toronto, Canada

Relevant Coursework: Machine Learning, Data structures, Algorithms, Matrix Algebra and Optimization, Computational Linguistics, Natural Language Computing, Computer Architecture, Probability & Statistics

# TECHNICAL SKILLS

Languages: Python, C/C++, CUDA, SQL, JS, Assembly, Verilog, MATLAB, React, Django

Libraries: TensorFlow, PyTorch, scikit-learn, keras, NumPy, pandas, Matplotlib, Hugging Face Transformers

Developer Tools: Git, Docker, VS Code, Visual Studio, Linux (Ubuntu), Shell/Bash, HDF5

## EXPERIENCE

# Machine Learning Researcher

May 2024 - Present

 $McMaster\ University\ -\ ChemAI\ Lab$ 

Hamilton, ON

- Founded and leading the Spectro project at McMaster University, exploring novel applications of AI in chemistry.
- Designed a multi-modal molecule prediction model, **Spectro**, achieving an **accuracy of 93**% (see projects).
- Collaborated with Dr. Kylie Luksa and other domain experts to inform model development and behaviour.
- Co-authored and submitted a paper to AI4Mat-NeurIPS, currently pending approval.

# Calibrations Engineering Intern

May 2023 – Sep. 2023

VACS Calibrations Toronto, ON

- Calibrated electronic and mechanical equipment, following the IEE and ISO17025 standards.
- Performed **statistical analysis**, including standard deviation, uncertainty propagation, and regression, to validate calibration accuracy and assess reliability.
- Developed insights from calibration data, improving process efficiency by 10% and measurement precision by 15%.

## **PROJECTS**

**Spectro** | TensorFlow, PyTorch, Hugging Face Transformers, JS, Linux

May 2024 – Present

- $\bullet \ \ \mathbf{Led} \ \ \mathbf{development} \ \ \mathbf{of} \ \mathbf{a} \ \mathbf{multimodal} \ \ \mathbf{molecule} \ \ \mathbf{elucidation} \ \ \mathbf{model} \ \ \mathbf{using} \ \ \mathbf{IR} \ \ \mathbf{and} \ \ \mathbf{NMR} \ \ \mathbf{data}, \ \mathbf{achieving} \ \ \mathbf{93\%} \ \ \mathbf{accuracy} \ \ \mathbf{a$
- Preprocessed and augmented datasets, employing data normalization and, reducing class imbalance by 30%.
- Fine-tuned a modified ResNet for functional group prediction from IR images, achieving an f1 score of 91%.
- Designed and tuned a decoder (RNN with LSTM), achieving 93% accuracy in molecule prediction.
- Implemented a **complete ML pipeline** in TensorFlow with: data augmentation and oversampling, **SMOTE**, cross-validation, a custom learning rate scheduler, **distributed training**, joint training, on-the-fly dataloading.

#### Chess Engine | C++, CUDA, Docker, Postman, JavaScript

Jun. 2023 – Apr. 2024

- C++ chess engine (rated 1800) with hardware optimizations like BitBoards and compile-time optimizations.
- Implemented multithreaded Negamax search with alpha-beta pruning, reducing search time by 45%.
- Integrated Zobrist hashing, a transposition table, and quiescence search, to improve search accuracy and eliminate redundant calculations.
- Utilized custom **CUDA** kernels for move generation and evaluations, significantly accelerating parallelizable tasks to achieve around **35 million nodes per second** in performance test.

Chess NNUE (Efficiently Updatable Neural Network) | PyTorch, C++, SQL, HDF5 May 2024 - Present

- Developing chess static evaluation with NNUE in PyTorch, currently achieving 80% accuracy.
- Augmented and preprocessed 83 million data points, exploring utilizing a custom C dataloader.
- Bridging the int\_8 quantized variant into my C++ chess engine, reducing inference time by 70%.

#### Remote ML Server | Linux (Ubuntu), Bash, Docker, Git

Jun. 2024

• Configured a retired GPU cryptocurrency mining rig into a remote server using docker, achieving 10x faster training time for ML models.

#### Publications

[1] (Pending) Chacko, Sondhi, et al. "A Multi-modal Approach for Molecule Elucidation Using IR and NMR Data." AI4Mat-NeurIPS 2024. December 2024.