NIDADAVOLU KAPIL CHANDRA

University of Texas at Austin

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EDUCATION

Univeristy of Texas at Austin

Master's in Computational Science, Engineering, and Mathematics

Austin, Texas 2019 – 2021

Indian Institute of Technology Madras

Dual Degree in Metallurgical and Materials Engineering; CGPA: 8.80/10

Chennai, India 2014 – 2019

o Dual Degree Class Rank: 4

o Minor: Physics

PUBLICATIONS

• "Modeling macroscopic material behavior with machine learning algorithms trained by micromechanical simulations" [PDF]

Authors: Denise Reimann, Kapil Nidadavolu, Hamad ul Hassan, Napat Vajragupta, Tobias Glasmachers, Philipp Junker, Alexander Hartmaier

SCHOLASTIC ACHIEVEMENTS

- One among 400 candidates all over India to be selected for the SRFP¹, 2017 to pursue a research internship in Indian Institute of Science, Bengaluru
- One among 4 candidates in my institute to be awarded a pre-placement interview for working in American Express on account of outstanding performance in Amex Analyze This, 2018 Comptetition [Link]
- Achieved All India Rank 3384 in Joint Entrance Examination 2014 (98th percentile)

PROJECTS

* RESEARCH PROJECTS

Development of Atomistic Potentials for Ag Nanoclusters using ML

• Master's Thesis, IIT Madras

Aug '18 - May'19

Guide: Dr. Satyesh Kumar Yadav

- o Performed first principle calculations for Ag nanoclusters using VASP over OpenMPI framework
- Developed a complete pipeline involving unsupervised and supervised learning techniques using scikit-learn, implementing Kernel Ridge Regression in conjunction with an atomic fingerprinting system to fit the force fields of Ag

Structure-Property linkages using Machine Learning

• Research intern at ICAMS Lab, Germany

May '18 - March '19

Guides: Prof. Alexander Hartmaier, Dr. Napat Vajragupta

- Developed machine learning models to replace a non-local crystal plasticity model
- Achieved an accuracy of 97.9% on grain size distribution predictions from flow curve data
- Performed feature engineering using Modified Voce Law to improve accuracy of the model

Development of memory efficient phase field model [Link]

• Research Intern at Indian Institute of Science, Bengaluru

May '17 - July '17

Guides: Prof. T.A. Abinandanan and Dr. Abhik Choudhury

- Implemented a phase field model for grain growth during recrystallization based on the paper Computer simulations of 2-D and 3-D ideal grain growth by Kim, Kim and Suzuki, in C
- Developed a memory efficient code by using **dynamic memory allocation** and **linked lists** which reduced the space complexity of algorithm from $\mathcal{O}(no.ofgrains \times x \times y)$ to $\mathcal{O}(1 \times x \times y)$

¹Summer Research Fellowship Program

Machine Learning for Engineering and Scientific Applications

• Course Project, IIT Madras

Jan '18 - April '18

Instructors: Dr. Balaji Srinivasan, Dr. Ganapathy Krishnamurthi

- Implemented CNN² in **TensorFlow** to classify MNIST dataset and obtained an accuracy of 98.4%
- \circ Developed codes for Linear Regression, Logistic Regression and ANN³ from scratch in Python without using already available frameworks
- Implemented high performance gradient descent optimization algorithms like **Adam**, **Adagrad** and **RMSprop** from scratch in Python

Computational Materials Engineering Lab

• Course Project, IIT Madras

Aug '17 - Dec '17

Instructors: Prof. G Phanikumar, Dr. Kanjarla Anand, Dr. Parasuraman Swaminathan

- Implemented **isotropic and anisotropic dendritic growth** model based on the paper "Modeling and numerical simulations of dendritic crystal growth" by Ryo Kobayashi, in **Matlab**
- o Devised algorithms for image processing tools like Object Counting and Edge Detection in MATLAB
- Implemented algorithms for solving ordinary partial differential equations in MATLAB

* SELF MOTIVATED PROJECTS

Online Automation Tools [Link]

- Python Script
 - Developed scripts in Python which can automatically book train tickets online
 - \circ Developed automatic mass SMS sender using **Selenium** and **Beautiful Soup** module in Python

PUBLICATIONS

• Reimann, Denise; Nidadavolu, Kapil; Hassan, Hamad ul; Vajragupta, Napat; Glasmachers, Tobias; Junker, Philipp; et al. (2019): Modeling Macroscopic Material Behavior With Machine Learning Algorithms Trained by Micromechanical Simulations. Frontiers. Collection.

TEACHING ASSISTANTSHIP

Computational Materials Engineering Lab

Guides: Prof. G Phanikumar, Dr. Kanjarla Anand, Dr. Parasuraman Swaminathan Aug '18 – Dec '18

 $\circ\,$ Assisted undergraduate students to understand concepts and clarified their doubts inside and outside the class

COURSEWORK

- Algorithms & Computing: Data Structures and Algorithms, Statistical Models for Big Data, Bayesian Statistical Methods
- Machine Learning and Deep Learning: Machine Learning, Neural Networks and Deep Learning, Improving Deep Neural Networks, Introduction to Data Science In Python, Reinforcement Learning
- Materials Science & Physics: Computational Materials Engineering Lab, Atomistic Modeling of Materials, Quantum Physics

TECHNOLOGY SUMMARY

- Programming & Scripting Languages: Python, C, R, SQL, Fortran, BASH Shell
- Machine Learning Frameworks: TensorFlow, Keras, Scikit-Learn
- Miscellaneous Tools: MATLAB, LATEX, Git

²Convolutional Neural Networks

³Artificial Neural Networks

CO CURRICULAR & EXTRA CURRICULAR ACTIVITIES

• Data Science Competitions

- \circ Ranked 12thout of 1500+ teams all over India in **Amex Analyze This** competition, 2018
- Awarded a **Pre-Placement Interview** offer for working in American Express on account of outstanding performance in the competition
- Regular participant in data science competitions on the campus and on online platforms like Kaggle