

Divya Karade

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Professional Summary

- AI researcher in drug discovery with 10+ years of experience in Biological Sciences (Biotech, Cheminformatics & Comp. Bio.)
- National-award-winning innovator for “AIDrugApp: Artificial Intelligence-based Web-Application for Drug Discovery”
- Collaborated and Managing a team of 3 for upcoming ‘AIDrugApp’ projects
- Proficient knowledge in Drug Designing, Machine/Deep Learning, Data Analytics, Algorithms and metabolomics datasets
- Excellent record of inter-disciplinary collaboration with researchers from biology, chemistry & computational science fields
- Solid publication record with 4 first-author publications in peer-reviewed journals and many in the pipeline
- Well versed in popular data science python libraries such as TensorFlow, Keras, sklearn, NumPy, Pandas, RDKit, PubChemPy

Technical Skills

- **Programming Languages:** Python, R
- **Machine learning:** Building and deploying ML/ DL algorithms, Auto-ML, Auto-DL, Web-app development - streamlit, Data analysis -TensorFlow, Keras, scikit-learn, NumPy, Pandas, Data visualization - Matplotlib, seaborn, yellowbrick & ggplot2
- **Data Science:** Metabolomics and Chemical Data analysis, Data science pipeline (cleansing, wrangling, optimization, EDA, modeling, interpretation), Feature engineering, Statistics, Hypothesis testing, GitHub, Data collection & compiling, Excel
- **Chemoinformatics:** Structure & ligand-based drug design, Text mining, Network analysis, Chemical structure analysis, QSAR, virtual screening, virtual libraries, ADME, scaffold analysis, MOE, ChemAxon, ChemDraw, Cytoscape, RDKit, PubChemPy
- **Bioinformatics:** Molecular docking and analysis using computational chemistry tools like Schrodinger, MOE and AutoDock
- **Metabolomics:** Untargeted metabolomics, Mass spectrometric data analysis- PUTMEDID-LCMS in Taverna workbench, ProbMetab R package, Univariate and Multivariate statistical analysis
- **Genomics & Proteomics:** DNA/RNA/protein extraction, PCR, Phylogenetic statistical analysis, SDS-PAGE, gel electrophoresis
- **Miscellaneous:** MySQL, HTML, Scientific and technical writing, cloud platform - Heroku, AWS

Innovation Project

AIDrugApp: Artificial Intelligence-based Virtual Screening Web-App for Drug Discovery

Sep 2020 - Present

<https://sars-covid-app.herokuapp.com/>

(Python3 | TensorFlow | Keras | scikit-learn | Matplotlib | Streamlit | Heroku)

- Innovated and founded a user-friendly and novel web application platform called ‘AIDrugApp’ for drug discovery
- **Bioactivity prediction module:** Originated and Implemented QSAR based DL algorithms to AIDrugApp for bioactivity prediction of user-defined molecules against SARS-CoV-2
- **Custom ML module:** ML tools integrated to AIDrugApp for user-defined model building, calculations and predictions on user data
 1. Auto-Multi-ML - Created for building and comparing user-defined multiple ML models and making predictions on users data
 2. Auto-DL - Developed to build user-defined deep learning models with neural networks and making predictions on users data
 3. DesCal - Developed for calculating user-defined 2-D descriptors and fingerprints using molecular SMILES of users data

Education

Ph.D. in Science (Chemoinformatics, Metabolomics and Computational Biology)

Jan 2015 - Thesis submitted

Academy of Scientific and Innovative Research (AcSIR)

Pune, Maharashtra, India

- Thesis: Chemoinformatics based investigation of plant metabolites for their medicinal and crop protection values

Master of Science (Plant Biotechnology) - Cumulative GPA: 8.14 /10.0

Sep 2008 - Nov 2010

A.N.G.R.A. University, Dept. of Biotechnology

Hyderabad, India

- Thesis: Characterization of genetic variability in cowpea using morphological, biochemical and molecular markers.

Bachelor of Science (Agricultural Science) - Cumulative GPA: 7.68 /10.0

Aug 2004 - Aug 2008

Dr. P.D.K.V. University, College of Agriculture

Nagpur, Maharashtra, India

- Relevant Coursework: Statistics, Maths, Plant Biotechnology, Plant Biochemistry, Molecular Biology, Genetics, Microbiology

Experience

Senior Research Fellow

Jan 2017 - Present

CSIR National Chemical Laboratory, Chemical Engineering and Process Development Division

Pune, India

- Bridging In-Silico and Experimental: Chemoinformatics Analysis of Mass Spectrometry-Based Metabolomics Study of Soybean
 - Designed and executed laboratory biochemical experiments to optimize soybean tissue samples for LC/MS studies
 - Discovered and reported the presence of 14 novel small molecules in soybean by mass spectrometry data analysis
 - Developed virtual library of 815 novel drug-like molecules from soybean small molecules through chemoinformatics analysis
- Chemoinformatics Investigation on Chemical Defense in Plants
 - Investigated 280 allelochemicals from 162 plants for their pesticide properties by chemoinformatics and QSAR approaches
 - Designed virtual libraries of allelochemicals that were then quantitatively evaluated to screen 169 pesticide-like molecules

Junior Research Fellow

Jan 2015 - Dec 2016

CSIR National Chemical Laboratory, Chemical Engineering and Process Development Division

Pune, India

- Design of Novel Drug-like Molecules using Informatics Rich Secondary Metabolites Analysis of Indian Medicinal Plants
 - Analyzed molecular data (n=3459) by text mining 91206 PubMed literature related to 104 Indian medicinal plants
 - Created and Contributed a virtual library of 4147 novel drug-like molecules for building a database of Indian medicinal plants
 - Formulated 4 computational protocols for designing drug-like novel molecules from metabolites of medicinal & food crops

Patent Analyst (Life Science)

Mar 2013 - Dec 2014

CSIR Unit of Research & Development of Information Products, Life Science Division

Pune, India

- Examined current technology alerts and patent updates in the field of life sciences for 5 projects
- Interpreted comparative and competitive patent portfolio after novelty, patentability and prior art search
- Mapped data for constructing a technological patent landscape

Reviewer for the “Journal of Experimental and Theoretical Artificial Intelligence” (Taylor & Francis)

Feb 2021- Present

Honors & Awards

National & International

- National-level AWS DeepRacer (**Selected in top 100**) at Woman's league for community car races driven by reinforcement learning codes in python through AWS cloud-based 3D racing simulator organized by AWS Educate Apr 2021
- National Award (Consolation - **Selected in top 10 out of 377 shortlisted applicants**) for “AIDrugApp” at the open-source “Drug Discovery Hackathon-2020” against Covid-19 in phase-1 under moonshot category organized by Govt. of India. Mar 2021
- National Award (Certificate of Merit - **Selected in top 4 out of 31 finalists**) for “AIDrugApp” in student engineering model competition under “Digital India” theme held at “Indian International Science Festival” organized by Govt. of India Dec 2020

Academic

- Awarded Senior Research Fellowship for continuing Ph.D., CSIR-UGC (National Eligibility Test) NET exam, India 2017
- All India rank **90** for Junior Research Fellowship for Ph.D. admission in Life-Science, CSIR-UGC NET exam, India 2013
- Awarded for Graduate Aptitude Test in Engineering (Biotechnology) for Ph.D. admission, MHRD, India 2011, 2012
- All India rank **64** in National Eligibility for Lectureship in Life-Science, CSIR-UGC NET, India 2011
- Awarded National Eligibility for Lectureship in Plant biotechnology by ASRB NET, India 2011
- All India rank **31** in Plant Biotechnology common entrance exam for P.G. admission, JNU, India 2008
- All India rank **14** in Plant Biotechnology entrance exam for P.G. admission, ICAR, India 2008

Publications

- **Divya Karade**, D. V., N. Kadoo, R. Vyas, P. K. Ingle and M. Karthikeyan. Design of Novel Drug-like Molecules using Informatics Rich Secondary Metabolites Analysis of Indian Medicinal and Aromatic Plants. *Combinatorial Chemistry & High Throughput Screening* **2020**, 23 (10), pp. 1113-1131. (Impact factor: 1.2, [click here](#))
- **Divya Karade**, Vikas Karade. AIDrugApp: Artificial Intelligence-based Web-App for Virtual Screening of Inhibitors against SARS-COV-2. *ChemRxiv. Preprint* **2020**. (Web-App: [click here](#))
(Accepted in peer-reviewed *Journal of Experimental and Theoretical Artificial Intelligence*, Impact factor: 2.11)
- **Divya Karade**. AutoDL: Automated Deep Learning (Machine learning module of AIDrugApp - Artificial Intelligence Based Virtual Screening Web-App for Drug Discovery) (Version 1.0.0). *Zenodo*. **2021**. (Software: [click here](#))
- **Divya Karade**, S. Sivaramakrishnan, K. Venkateswaran, R.S. Reddy. Characterization of genetic variability among cowpea (*Vigna Unguiculata* L. Walp) germplasm using morphological, biochemical and molecular markers. *Jour. Of Med. Sci. & Tech.* **2012**, 1 (2), pp. 43-61. ([click here](#))
- **Divya Karade**, S. Mundhe, N. Kadoo, M. Ratnaparkhe, R. Vyas and M. Karthikeyan. Bridging In-Silico and Experimental: Chemoinformatics Analysis for Mass Spectrometry-Based Metabolomics Study of Soybean. *Metabolomics*. **2021**. (Impact factor: 3.167, Under review)

Research Presentations

Posters

- **Divya Karade**, Vikas Karade (**2021-2021**). ‘AIDrugApp: Artificial Intelligence Based Virtual Screening Web-App for Drug Discovery’ presented at “Indian International Science Festival” and “Drug Discovery Hackathon-2020” organized by GOI.
- **Divya Karade**, N. Kadoo, M. Karthikeyan (**2019**) ‘Antidiabetic Drug Designing Based on Organic Metabolites from Fenugreek: A Chemo-and Bioinformatics Approach’ presented at an international symposium conducted by C-DAC, Pune
- **Divya Karade**, N. Kadoo, M. Karthikeyan (**2017**) ‘Chemoinformatics Investigation of Organic Metabolites from Soybean Reveals Common Drug Scaffolds presented at CSIR-National Chemical Laboratory, Pune on the “National Science Day”
- **Divya Karade**, N. Kadoo, M. Karthikeyan (**2017**) ‘Drug Design based on Metabolomics of Indian Medicinal and Aromatic Plants: A Chemoinformatics Approach’ presented at CSIR-National Chemical Laboratory, Pune on the “National Science Day”

Workshop

Instructor (10 people) and participated in “Chemoinformatics-2018” for skill development program by CSIR-NCL, Pune