

# Tunca Dogan

## List of Publications by Year in descending order

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Version: 2024-02-01

34  
papers

22,397  
citations

430754

18  
h-index

454834

30  
g-index

44  
all docs

44  
docs citations

44  
times ranked

40846  
citing authors

#	ARTICLE	IF	CITATIONS
1	Learning functional properties of proteins with language models. <i>Nature Machine Intelligence</i> , 2022, 4, 227-245.	8.3	72
2	SLPred: a multi-view subcellular localization prediction tool for multi-location human proteins. <i>Bioinformatics</i> , 2022, 38, 4226-4229.	1.8	3
3	MDeePred: novel multi-channel protein featurization for deep learning-based binding affinity prediction in drug discovery. <i>Bioinformatics</i> , 2021, 37, 693-704.	1.8	61
4	UniProt: the universal protein knowledgebase in 2021. <i>Nucleic Acids Research</i> , 2021, 49, D480-D489.	6.5	4,709
5	Crowdsourced mapping of unexplored target space of kinase inhibitors. <i>Nature Communications</i> , 2021, 12, 3307.	5.8	41
6	CROsSBAR: comprehensive resource of biomedical relations with knowledge graph representations. <i>Nucleic Acids Research</i> , 2021, 49, e96-e96.	6.5	19
7	Protein domain-based prediction of drug/compound–target interactions and experimental validation on LIM kinases. <i>PLoS Computational Biology</i> , 2021, 17, e1009171.	1.5	13
8	Editorial: Machine Learning Methodologies to Study Molecular Interactions. <i>Frontiers in Molecular Biosciences</i> , 2021, 8, 806474.	1.6	1
9	A crowdsourcing open platform for literature curation in UniProt. <i>PLoS Biology</i> , 2021, 19, e3001464.	2.6	74
10	Data Centric Molecular Analysis and Evaluation of Hepatocellular Carcinoma Therapeutics Using Machine Intelligence-Based Tools. <i>Journal of Gastrointestinal Cancer</i> , 2021, 52, 1266-1276.	0.6	0
11	DEEPScreen: high performance drug–target interaction prediction with convolutional neural networks using 2-D structural compound representations. <i>Chemical Science</i> , 2020, 11, 2531-2557.	3.7	131
12	iBioProVis: interactive visualization and analysis of compound bioactivity space. <i>Bioinformatics</i> , 2020, 36, 4227-4230.	1.8	7
13	Abstract 5235: In vitro validation of drug-target interactions revealed in silico by Comprehensive Resource of Biomedical Relations with Network Representations and Deep Learning (CROsSBAR) in HCC. , 2020, , .		0
14	Recent applications of deep learning and machine intelligence on in silico drug discovery: methods, tools and databases. <i>Briefings in Bioinformatics</i> , 2019, 20, 1878-1912.	3.2	310
15	FAIR adoption, assessment and challenges at UniProt. <i>Scientific Data</i> , 2019, 6, 175.	2.4	11
16	DEEPred: Automated Protein Function Prediction with Multi-task Feed-forward Deep Neural Networks. <i>Scientific Reports</i> , 2019, 9, 7344.	1.6	80
17	The CAFA challenge reports improved protein function prediction and new functional annotations for hundreds of genes through experimental screens. <i>Genome Biology</i> , 2019, 20, 244.	3.8	261
18	UniProt: a worldwide hub of protein knowledge. <i>Nucleic Acids Research</i> , 2019, 47, D506-D515.	6.5	6,185

#	ARTICLE	IF	CITATIONS
19	Phylogenetic and Other Conservation-Based Approaches to Predict Protein Functional Sites. <i>Methods in Molecular Biology</i> , 2018, 1762, 51-69.	0.4	4
20	Large-scale automated function prediction of protein sequences and an experimental case study validation on PTEN transcript variants. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018, 86, 135-151.	1.5	13
21	HPO2GO: prediction of human phenotype ontology term associations for proteins using cross ontology annotation co-occurrences. <i>PeerJ</i> , 2018, 6, e5298.	0.9	27
22	ECPred: a tool for the prediction of the enzymatic functions of protein sequences based on the EC nomenclature. <i>BMC Bioinformatics</i> , 2018, 19, 334.	1.2	99
23	A Structural Perspective on the Modulation of Protein-Protein Interactions with Small Molecules. <i>Current Topics in Medicinal Chemistry</i> , 2018, 18, 700-713.	1.0	6
24	UniProt: the universal protein knowledgebase. <i>Nucleic Acids Research</i> , 2017, 45, D158-D169.	6.5	4,240
25	From the research laboratory to the database: the <i>Caenorhabditis elegans</i> kinome in UniProtKB. <i>Biochemical Journal</i> , 2017, 474, 493-515.	1.7	9
26	An expanded evaluation of protein function prediction methods shows an improvement in accuracy. <i>Genome Biology</i> , 2016, 17, 184.	3.8	308
27	UniProt-DAAC: domain architecture alignment and classification, a new method for automatic functional annotation in UniProtKB. <i>Bioinformatics</i> , 2016, 32, 2264-2271.	1.8	37
28	Tools and data services registry: a community effort to document bioinformatics resources. <i>Nucleic Acids Research</i> , 2016, 44, D38-D47.	6.5	113
29	Unsupervised identification of redundant domain entries in InterPro database using clustering techniques. , 2015, , .		0
30	UniProt: a hub for protein information. <i>Nucleic Acids Research</i> , 2015, 43, D204-D212.	6.5	4,370
31	Activities at the Universal Protein Resource (UniProt). <i>Nucleic Acids Research</i> , 2014, 42, D191-D198.	6.5	1,162
32	Automatic Identification of Highly Conserved Family Regions and Relationships in Genome Wide Datasets Including Remote Protein Sequences. <i>PLoS ONE</i> , 2013, 8, e75458.	1.1	7
33	2-D Thresholding of the Connectivity Map Following the Multiple Sequence Alignments of Diverse Datasets. , 2013, , .		0
34	Evolutionary relationships between gene sequences via nonlinear embedding. , 2010, , .		0