

# Juan Liu

## List of Publications by Year in descending order

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Version: 2024-02-01

81  
papers

1,296  
citations

394421

19  
h-index

377865

34  
g-index

81  
all docs

81  
docs citations

81  
times ranked

1754  
citing authors

#	ARTICLE	IF	CITATIONS
1	CSCD2: an integrated interactional database of cancer-specific circular RNAs. <i>Nucleic Acids Research</i> , 2022, 50, D1179-D1183.	14.5	35
2	Interpretable Machine Learning for Early Prediction of Prognosis in Sepsis: A Discovery and Validation Study. <i>Infectious Diseases and Therapy</i> , 2022, 11, 1117-1132.	4.0	37
3	DeepRF: A deep learning method for predicting metabolic pathways in organisms based on annotated genomes. <i>Computers in Biology and Medicine</i> , 2022, 147, 105756.	7.0	4
4	CytoBrain: Cervical Cancer Screening System Based on Deep Learning Technology. <i>Journal of Computer Science and Technology</i> , 2021, 36, 347-360.	1.5	38
5	Review of Machine Learning Methods for the Prediction and Reconstruction of Metabolic Pathways. <i>Frontiers in Molecular Biosciences</i> , 2021, 8, 634141.	3.5	15
6	Improved Diagnostic Accuracy of Ameloblastoma and Odontogenic Keratocyst on Cone-Beam CT by Artificial Intelligence. <i>Frontiers in Oncology</i> , 2021, 11, 793417.	2.8	11
7	Network-based cancer genomic data integration for pattern discovery. <i>BMC Genomic Data</i> , 2021, 22, 54.	1.7	2
8	COMNA: Core-attachment based protein complex detection via multiple network alignment. , 2021, , .		1
9	TransMixNet: An Attention Based Double-Branch Model for White Blood Cell Classification and Its Training with the Fuzzified Training Data. , 2021, , .		5
10	Link Prediction Only With Interaction Data and its Application on Drug Repositioning. <i>IEEE Transactions on Nanobioscience</i> , 2020, 19, 547-555.	3.3	10
11	GPPIAL: A New Global PPI Network Aligner Based on Orthologs. , 2020, , .		0
12	Group-sparse SVD Models via $L_1$ - and $L_0$ -norm Penalties and Their Applications in Biological Data. <i>IEEE Transactions on Knowledge and Data Engineering</i> , 2019, , 1-1.	5.7	6
13	Prediction of drug-disease associations based on ensemble meta paths and singular value decomposition. <i>BMC Bioinformatics</i> , 2019, 20, 134.	2.6	24
14	A Robust Approach to Locate HER2 and CEN17 Signals in Varied FISH Images. , 2019, , .		0
15	Predicting Drug-Disease Treatment Associations Based on Topological Similarity and Singular Value Decomposition. , 2019, , .		4
16	Gene Functional Module Discovery via Integrating Gene Expression and PPI Network Data. <i>Lecture Notes in Computer Science</i> , 2019, , 116-126.	1.3	1
17	Edge-group sparse PCA for network-guided high dimensional data analysis. <i>Bioinformatics</i> , 2018, 34, 3479-3487.	4.1	43
18	Discovering large conserved functional components in global network alignment by graph matching. <i>BMC Genomics</i> , 2018, 19, 670.	2.8	6

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19	Prediction of drug-disease treatment relations based on positive and unlabeled samples. Journal of Intelligent and Fuzzy Systems, 2018, 35, 1363-1373.	1.4	3
20	Sparse Weighted Canonical Correlation Analysis. Chinese Journal of Electronics, 2018, 27, 459-466.	1.5	6
21	Differential function analysis: identifying structure and activation variations in dysregulated pathways. Science China Information Sciences, 2017, 60, 1.	4.3	3
22	Comparative network stratification analysis for identifying functional interpretable network biomarkers. BMC Bioinformatics, 2017, 18, 48.	2.6	7
23	Local network component analysis for quantifying transcription factor activities. Methods, 2017, 124, 25-35.	3.8	14
24	A Novel Sparse Penalty for Singular Value Decomposition. Chinese Journal of Electronics, 2017, 26, 306-312.	1.5	2
25	GMAAlign: A new network aligner for revealing large conserved functional components. , 2017, , .		0
26	Predicting drug-disease interactions by semi-supervised graph cut algorithm and three-layer data integration. BMC Medical Genomics, 2017, 10, 79.	1.5	23
27	Surface shapes and surrounding environment analysis of single- and double-stranded DNA-binding proteins in protein-DNA interface. Proteins: Structure, Function and Bioinformatics, 2016, 84, 979-989.	2.6	10
28	Quantum dots-based double imaging combined with organic dye imaging to establish an automatic computerized method for cancer Ki67 measurement. Scientific Reports, 2016, 6, 20564.	3.3	20
29	Semi-supervised graph cut algorithm for drug repositioning by integrating drug, disease and genomic associations. , 2016, , .		2
30	Integration of multiple heterogeneous omics data. , 2016, , .		4
31	A Two-Stage Method to Identify Joint Modules From Matched MicroRNA and mRNA Expression Data. IEEE Transactions on Nanobioscience, 2016, 15, 362-370.	3.3	10
32	Network stratification analysis for identifying function-specific network layers. Molecular BioSystems, 2016, 12, 1232-1240.	2.9	1
33	Integration of a prognostic gene module with a drug sensitivity module to identify drugs that could be repurposed for breast cancer therapy. Computers in Biology and Medicine, 2015, 61, 163-171.	7.0	4
34	Segmentation of Hematoxylin-Eosin stained breast cancer histopathological images based on pixel-wise SVM classifier. Science China Information Sciences, 2015, 58, 1-13.	4.3	24
35	Measuring the quality of linear patterns in biclusters. Methods, 2015, 83, 18-27.	3.8	8
36	Predicting Response to Preoperative Chemotherapy Agents by Identifying Drug Action on Modeled MicroRNA Regulation Networks. PLoS ONE, 2014, 9, e98140.	2.5	3

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37	The convergence rate of a three-term HS method with restart strategy for unconstrained optimization problems. <i>Optimization</i> , 2014, 63, 1387-1400.	1.7	0
38	Construction and investigation of breast cancer-specific ceRNA network based on the mRNA and miRNA expression data. <i>IET Systems Biology</i> , 2014, 8, 96-103.	1.5	110
39	Analysis and classification of DNA-binding sites in single-stranded and double-stranded DNA-binding proteins using protein information. <i>IET Systems Biology</i> , 2014, 8, 176-183.	1.5	11
40	Two-step segmentation of Hematoxylin-Eosin stained histopathological images for prognosis of breast cancer. , 2014, , .		14
41	A modified Polak-Ribiere-Polyak descent method for unconstrained optimization. <i>Optimization Methods and Software</i> , 2014, 29, 177-188.	2.4	3
42	A computational model to predict bone metastasis in breast cancer by integrating the dysregulated pathways. <i>BMC Cancer</i> , 2014, 14, 618.	2.6	14
43	Deciphering early development of complex diseases by progressive module network. <i>Methods</i> , 2014, 67, 334-343.	3.8	42
44	Identification of single-stranded and double-stranded dna binding proteins based on protein structure. <i>BMC Bioinformatics</i> , 2014, 15, S4.	2.6	18
45	Databases for B-Cell Epitopes. <i>Methods in Molecular Biology</i> , 2014, 1184, 135-148.	0.9	5
46	Inferring Gene Dependency Network Specific to Phenotypic Alteration Based on Gene Expression Data and Clinical Information of Breast Cancer. <i>PLoS ONE</i> , 2014, 9, e92023.	2.5	9
47	Integrating peptides' sequence and energy of contact residues information improves prediction of peptide and HLA-I binding with unknown alleles. <i>BMC Bioinformatics</i> , 2013, 14, S1.	2.6	2
48	Predicting immunogenic T-cell epitopes by combining various sequence-derived features. , 2013, , .		10
49	Computer-Based Image Studies on Tumor Nests Mathematical Features of Breast Cancer and Their Clinical Prognostic Value. <i>PLoS ONE</i> , 2013, 8, e82314.	2.5	25
50	Predicting linear B-cell epitopes by using sequence-derived structural and physicochemical features. <i>International Journal of Data Mining and Bioinformatics</i> , 2012, 6, 557.	0.1	22
51	Predicting distant metastasis in breast cancer using ensemble classifier based on context-specific miRNA regulation modules. , 2012, , .		2
52	Computational Prediction of Conformational B-Cell Epitopes from Antigen Primary Structures by Ensemble Learning. <i>PLoS ONE</i> , 2012, 7, e43575.	2.5	52
53	Prediction of heme binding residues from protein sequences with integrative sequence profiles. <i>Proteome Science</i> , 2012, 10, S20.	1.7	36
54	Predicting Binding-Peptide of HLA-I on Unknown Alleles by Integrating Sequence Information and Energies of Contact Residues. <i>Communications in Computer and Information Science</i> , 2012, , 160-165.	0.5	0

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55	Dynamic remodeling of context-specific miRNAs regulation networks facilitate in silico cancer drug screening. , 2011, , .		0
56	Context-specific miRNA regulation network predicts cancer prognosis. , 2011, , .		2
57	A novel computational framework for simultaneous integration of multiple types of genomic data to identify microRNA-gene regulatory modules. <i>Bioinformatics</i> , 2011, 27, i401-i409.	4.1	212
58	Exploiting a Reduced Set of Weighted Average Features to Improve Prediction of DNA-Binding Residues from 3D Structures. <i>PLoS ONE</i> , 2011, 6, e28440.	2.5	30
59	Prediction of conformational B-cell epitopes from 3D structures by random forests with a distance-based feature. <i>BMC Bioinformatics</i> , 2011, 12, 341.	2.6	92
60	An accurate feature-based method for identifying DNA-binding residues on protein surfaces. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 509-517.	2.6	68
61	Distinct interfacial biclique patterns between ssDNA-binding proteins and those with dsDNAs. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 598-610.	2.6	4
62	Predicting cleavage sites in exogenous antigen using weighted SVM. , 2010, , .		0
63	A novel locally linear embedding and wavelet transform based encoding method for prediction of MHC-II binding affinity. <i>Interdisciplinary Sciences, Computational Life Sciences</i> , 2010, 2, 145-150.	3.6	1
64	Mixture classification model based on clinical markers for breast cancer prognosis. <i>Artificial Intelligence in Medicine</i> , 2010, 48, 129-137.	6.5	33
65	Quantitative prediction of MHC-II binding affinity using particle swarm optimization. <i>Artificial Intelligence in Medicine</i> , 2010, 50, 127-132.	6.5	16
66	In silico Genetic Network Models for *Pre-clinical Drug Prioritization. <i>Nature Precedings</i> , 2010, , .	0.1	0
67	Pre-Clinical Drug Prioritization via Prognosis-Guided Genetic Interaction Networks. <i>PLoS ONE</i> , 2010, 5, e13937.	2.5	9
68	Quantitative prediction of MHC-II peptide binding affinity using relevance vector machine. <i>Applied Intelligence</i> , 2009, 31, 180-187.	5.3	10
69	Prediction of $\beta$ -turns using double BP network with novel coding schemes of amino acids. <i>Wuhan University Journal of Natural Sciences</i> , 2009, 14, 119-124.	0.4	1
70	MDAS: An integrated system for metabonomic data analysis. <i>Interdisciplinary Sciences, Computational Life Sciences</i> , 2009, 1, 61-71.	3.6	1
71	Prediction of Linear B-Cell Epitopes Using AAT Scale. , 2009, , .		3
72	BiodMHC: an online server for the prediction of MHC class II-peptide binding affinity. <i>Journal of Genetics and Genomics</i> , 2009, 36, 289-296.	3.9	2

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73	A Bayesian regression approach to the prediction of MHC-II binding affinity. Computer Methods and Programs in Biomedicine, 2008, 92, 1-7.	4.7	33
74	Web-Based Parallel Corpora for Statistical Machine Translation. , 2007, , .		0
75	Web-based parallel corpora for statistical machine translation. , 2007, , .		0
76	A simple method of inferring pairwise gene interactions from microarray time series data. , 2005, , .		1
77	A concept learning method based on a hybrid genetic algorithm. Science in China Series D: Earth Sciences, 1998, 41, 488-495.	0.9	3
78	Selecting informative genes using a multiobjective evolutionary algorithm. , 0, , .		11
79	Tumor classification based on gene microarray data and hybrid learning method. , 0, , .		2
80	The simulated annealing method applied to protein structure prediction. , 0, , .		0
81	A genetic algorithm applied to optimal gene subset selection. , 0, , .		1