Lei Chen

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

182 3,879 48 34 h-index g-index citations papers 6.19 193 5,741 3.5 L-index avg, IF ext. citations ext. papers

#	Paper	IF	Citations
182	Predicting Anatomical Therapeutic Chemical (ATC) classification of drugs by integrating chemical-chemical interactions and similarities. <i>PLoS ONE</i> , 2012 , 7, e35254	3.7	133
181	Identify Key Sequence Features to Improve CRISPR sgRNA Efficacy. <i>IEEE Access</i> , 2017 , 5, 26582-26590	3.5	91
180	A similarity-based method for prediction of drug side effects with heterogeneous information. <i>Mathematical Biosciences</i> , 2018 , 306, 136-144	3.9	89
179	Classification and analysis of regulatory pathways using graph property, biochemical and physicochemical property, and functional property. <i>PLoS ONE</i> , 2011 , 6, e25297	3.7	77
178	Analysis of cancer-related lncRNAs using gene ontology and KEGG pathways. <i>Artificial Intelligence in Medicine</i> , 2017 , 76, 27-36	7.4	74
177	Prediction and analysis of essential genes using the enrichments of gene ontology and KEGG pathways. <i>PLoS ONE</i> , 2017 , 12, e0184129	3.7	74
176	Prediction of protein domain with mRMR feature selection and analysis. <i>PLoS ONE</i> , 2012 , 7, e39308	3.7	74
175	Prediction of protein-protein interaction sites by random forest algorithm with mRMR and IFS. <i>PLoS ONE</i> , 2012 , 7, e43927	3.7	69
174	Gene expression differences among different MSI statuses in colorectal cancer. <i>International Journal of Cancer</i> , 2018 , 143, 1731-1740	7.5	68
173	Predicting Drug Side Effects with Compact Integration of Heterogeneous Networks. <i>Current Bioinformatics</i> , 2019 , 14, 709-720	4.7	64
172	Analysis and prediction of drug-drug interaction by minimum redundancy maximum relevance and incremental feature selection. <i>Journal of Biomolecular Structure and Dynamics</i> , 2017 , 35, 312-329	3.6	61
171	Prediction and analysis of cell-penetrating peptides using pseudo-amino acid composition and random forest models. <i>Amino Acids</i> , 2015 , 47, 1485-93	3.5	59
170	A hybrid method for prediction and repositioning of drug Anatomical Therapeutic Chemical classes. <i>Molecular BioSystems</i> , 2014 , 10, 868-77		58
169	Gene expression profiling gut microbiota in different races of humans. Scientific Reports, 2016 , 6, 2307	5 4.9	55
168	Deciphering the effects of gene deletion on yeast longevity using network and machine learning approaches. <i>Biochimie</i> , 2012 , 94, 1017-25	4.6	54
167	Identification of Drug-Drug Interactions Using Chemical Interactions. <i>Current Bioinformatics</i> , 2017 , 12,	4.7	54
166	Predicting the network of substrate-enzyme-product triads by combining compound similarity and functional domain composition. <i>BMC Bioinformatics</i> , 2010 , 11, 293	3.6	50

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165	Identification of compound-protein interactions through the analysis of gene ontology, KEGG enrichment for proteins and molecular fragments of compounds. <i>Molecular Genetics and Genomics</i> , 2016 , 291, 2065-2079	3.1	48	
164	Discriminating cirRNAs from other lncRNAs using a hierarchical extreme learning machine (H-ELM) algorithm with feature selection. <i>Molecular Genetics and Genomics</i> , 2018 , 293, 137-149	3.1	45	
163	Identification of gene expression signatures across different types of neural stem cells with the Monte-Carlo feature selection method. <i>Journal of Cellular Biochemistry</i> , 2018 , 119, 3394-3403	4.7	45	
162	iATC-NRAKEL: an efficient multi-label classifier for recognizing anatomical therapeutic chemical classes of drugs. <i>Bioinformatics</i> , 2020 , 36, 1391-1396	7.2	44	
161	Identification of Differentially Expressed Genes between Original Breast Cancer and Xenograft Using Machine Learning Algorithms. <i>Genes</i> , 2018 , 9,	4.2	42	
160	Identification of synthetic lethality based on a functional network by using machine learning algorithms. <i>Journal of Cellular Biochemistry</i> , 2019 , 120, 405-416	4.7	42	
159	Signal propagation in protein interaction network during colorectal cancer progression. <i>BioMed Research International</i> , 2013 , 2013, 287019	3	41	
158	Identification of the copy number variant biomarkers for breast cancer subtypes. <i>Molecular Genetics and Genomics</i> , 2019 , 294, 95-110	3.1	41	
157	iATC-FRAKEL: a simple multi-label web server for recognizing anatomical therapeutic chemical classes of drugs with their fingerprints only. <i>Bioinformatics</i> , 2020 , 36, 3568-3569	7.2	40	
156	Gene Ontology and KEGG Pathway Enrichment Analysis of a Drug Target-Based Classification System. <i>PLoS ONE</i> , 2015 , 10, e0126492	3.7	39	
155	Prediction of Drug Side Effects with a Refined Negative Sample Selection Strategy. <i>Computational and Mathematical Methods in Medicine</i> , 2020 , 2020, 1573543	2.8	38	
154	Identification of hepatocellular carcinoma related genes with k-th shortest paths in a protein-protein interaction network. <i>Molecular BioSystems</i> , 2013 , 9, 2720-8		37	
153	Analysis of tumor suppressor genes based on gene ontology and the KEGG pathway. <i>PLoS ONE</i> , 2014 , 9, e107202	3.7	37	
152	Application of the Shortest Path Algorithm for the Discovery of Breast Cancer-Related Genes. <i>Current Bioinformatics</i> , 2016 , 11, 51-58	4.7	37	
151	Similarity-Based Machine Learning Model for Predicting the Metabolic Pathways of Compounds. <i>IEEE Access</i> , 2020 , 8, 130687-130696	3.5	36	
150	The use of Gene Ontology terms and KEGG pathways for analysis and prediction of oncogenes. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2016 , 1860, 2725-34	4	35	
149	Tissue Expression Difference between mRNAs and lncRNAs. <i>International Journal of Molecular Sciences</i> , 2018 , 19,	6.3	35	
148	Classifying ten types of major cancers based on reverse phase protein array profiles. <i>PLoS ONE</i> , 2015 , 10, e0123147	3.7	33	

147	Predicting metabolic pathways of small molecules and enzymes based on interaction information of chemicals and proteins. <i>PLoS ONE</i> , 2012 , 7, e45944	3.7	33
146	Predicting the types of metabolic pathway of compounds using molecular fragments and sequential minimal optimization. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2016 , 19, 136	-43	33
145	Identifying and analyzing different cancer subtypes using RNA-seq data of blood platelets. <i>Oncotarget</i> , 2017 , 8, 87494-87511	3.3	32
144	Identifying Patients with Atrioventricular Septal Defect in Down Syndrome Populations by Using Self-Normalizing Neural Networks and Feature Selection. <i>Genes</i> , 2018 , 9,	4.2	31
143	Multiple classifier integration for the prediction of protein structural classes. <i>Journal of Computational Chemistry</i> , 2009 , 30, 2248-54	3.5	31
142	Labelling algorithms for paired-domination problems in block and interval graphs. <i>Journal of Combinatorial Optimization</i> , 2010 , 19, 457-470	0.9	31
141	Determining protein-protein functional associations by functional rules based on gene ontology and KEGG pathway. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2021 , 1869, 140621	4	31
140	Identifying Transcriptomic Signatures and Rules for SARS-CoV-2 Infection. <i>Frontiers in Cell and Developmental Biology</i> , 2020 , 8, 627302	5.7	31
139	Identification of the Gene Expression Rules That Define the Subtypes in Glioma. <i>Journal of Clinical Medicine</i> , 2018 , 7,	5.1	30
138	Finding candidate drugs for hepatitis C based on chemical-chemical and chemical-protein interactions. <i>PLoS ONE</i> , 2014 , 9, e107767	3.7	29
137	Identification of Protein Subcellular Localization With Network and Functional Embeddings. <i>Frontiers in Genetics</i> , 2020 , 11, 626500	4.5	29
136	Identification of Human Membrane Protein Types by Incorporating Network Embedding Methods. <i>IEEE Access</i> , 2019 , 7, 140794-140805	3.5	28
135	Prediction of human genesTregulatory functions based on proteinprotein interaction network. <i>Protein and Peptide Letters</i> , 2012 , 19, 910-6	1.9	28
134	Prediction of Metabolic Pathway Using Graph Property, Chemical Functional Group and Chemical Structural Set. <i>Current Bioinformatics</i> , 2013 , 8, 200-207	4.7	27
133	A Binary Classifier for the Prediction of EC Numbers of Enzymes. <i>Current Proteomics</i> , 2019 , 16, 383-391	0.7	27
132	Identification of leukemia stem cell expression signatures through Monte Carlo feature selection strategy and support vector machine. <i>Cancer Gene Therapy</i> , 2020 , 27, 56-69	5.4	27
131	Identification of Genes Associated with Breast Cancer Metastasis to Bone on a Protein-Protein Interaction Network with a Shortest Path Algorithm. <i>Journal of Proteome Research</i> , 2017 , 16, 1027-1038	3 ^{5.6}	26
130	Identification of novel proliferative diabetic retinopathy related genes on proteinprotein interaction network. <i>Neurocomputing</i> , 2016 , 217, 63-72	5.4	26

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129	Inferring Novel Tumor Suppressor Genes with a Protein-Protein Interaction Network and Network Diffusion Algorithms. <i>Molecular Therapy - Methods and Clinical Development</i> , 2018 , 10, 57-67	6.4	26
128	Prediction of gene phenotypes based on GO and KEGG pathway enrichment scores. <i>BioMed Research International</i> , 2013 , 2013, 870795	3	26
127	Classification of Widely and Rarely Expressed Genes with Recurrent Neural Network. <i>Computational and Structural Biotechnology Journal</i> , 2019 , 17, 49-60	6.8	26
126	Analysis of Expression Pattern of snoRNAs in Different Cancer Types with Machine Learning Algorithms. <i>International Journal of Molecular Sciences</i> , 2019 , 20,	6.3	25
125	Analysis of protein pathway networks using hybrid properties. <i>Molecules</i> , 2010 , 15, 8177-92	4.8	25
124	Computational analysis of HIV-1 resistance based on gene expression profiles and the virus-host interaction network. <i>PLoS ONE</i> , 2011 , 6, e17291	3.7	25
123	Mining for novel tumor suppressor genes using a shortest path approach. <i>Journal of Biomolecular Structure and Dynamics</i> , 2016 , 34, 664-75	3.6	24
122	SACICA: a sparse approximation coefficient-based ICA model for functional magnetic resonance imaging data analysis. <i>Journal of Neuroscience Methods</i> , 2013 , 216, 49-61	3	24
121	Prediction of effective drug combinations by chemical interaction, protein interaction and target enrichment of KEGG pathways. <i>BioMed Research International</i> , 2013 , 2013, 723780	3	24
120	Identifying Methylation Pattern and Genes Associated with Breast Cancer Subtypes. <i>International Journal of Molecular Sciences</i> , 2019 , 20,	6.3	23
119	A computational method using the random walk with restart algorithm for identifying novel epigenetic factors. <i>Molecular Genetics and Genomics</i> , 2018 , 293, 293-301	3.1	23
118	Inferring anatomical therapeutic chemical (ATC) class of drugs using shortest path and random walk with restart algorithms. <i>Biochimica Et Biophysica Acta - Molecular Basis of Disease</i> , 2018 , 1864, 2228	-2240	22
117	Identifying novel protein phenotype annotations by hybridizing protein-protein interactions and protein sequence similarities. <i>Molecular Genetics and Genomics</i> , 2016 , 291, 913-34	3.1	22
116	A fast-FENICA method on resting state fMRI data. <i>Journal of Neuroscience Methods</i> , 2012 , 209, 1-12	3	22
115	Identifying protein complexes using hybrid properties. <i>Journal of Proteome Research</i> , 2009 , 8, 5212-8	5.6	22
114	Drug Target Group Prediction with Multiple Drug Networks. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2020 , 23, 274-284	1.3	22
113	Tissue differences revealed by gene expression profiles of various cell lines. <i>Journal of Cellular Biochemistry</i> , 2018 , 120, 7068	4.7	22
112	Identification of new candidate drugs for lung cancer using chemical-chemical interactions, chemical-protein interactions and a K-means clustering algorithm. <i>Journal of Biomolecular Structure and Dynamics</i> , 2016 , 34, 906-17	3.6	21

111	Predicting chemical toxicity effects based on chemical-chemical interactions. <i>PLoS ONE</i> , 2013 , 8, e5651	73.7	21
110	A Feature and Algorithm Selection Method for Improving the Prediction of Protein Structural Class. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2017 , 20, 612-621	1.3	21
109	Detecting the Multiomics Signatures of Factor-Specific Inflammatory Effects on Airway Smooth Muscles. <i>Frontiers in Genetics</i> , 2020 , 11, 599970	4.5	21
108	A computational method for the identification of new candidate carcinogenic and non-carcinogenic chemicals. <i>Molecular BioSystems</i> , 2015 , 11, 2541-50		20
107	Predicting drugs side effects based on chemical-chemical interactions and protein-chemical interactions. <i>BioMed Research International</i> , 2013 , 2013, 485034	3	20
106	A linear-time algorithm for paired-domination problem in strongly chordal graphs. <i>Information Processing Letters</i> , 2009 , 110, 20-23	0.8	20
105	Prediction of Nitrated Tyrosine Residues in Protein Sequences by Extreme Learning Machine and Feature Selection Methods. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2018 , 21, 393-402	1.3	20
104	Identifying protein subcellular location with embedding features learned from networks. <i>Current Proteomics</i> , 2020 , 17,	0.7	20
103	Identification of novel candidate drivers connecting different dysfunctional levels for lung adenocarcinoma using protein-protein interactions and a shortest path approach. <i>Scientific Reports</i> , 2016 , 6, 29849	4.9	20
102	iMPTCE-Hnetwork: A Multilabel Classifier for Identifying Metabolic Pathway Types of Chemicals and Enzymes with a Heterogeneous Network. <i>Computational and Mathematical Methods in Medicine</i> , 2021, 6683051	2.8	20
101	Identification of drugdisease associations by using multiple drug and disease networks. <i>Current Bioinformatics</i> , 2021 , 16,	4.7	20
100	Computationally identifying virulence factors based on KEGG pathways. <i>Molecular BioSystems</i> , 2013 , 9, 1447-52		19
99	Gene ontology and KEGG enrichment analyses of genes related to age-related macular degeneration. <i>BioMed Research International</i> , 2014 , 2014, 450386	3	19
98	Identification of lung-cancer-related genes with the shortest path approach in a protein-protein interaction network. <i>BioMed Research International</i> , 2013 , 2013, 267375	3	19
97	HIV infection alters the human epigenetic landscape. <i>Gene Therapy</i> , 2019 , 26, 29-39	4	19
96	A Network Integration Method for Deciphering the Types of Metabolic Pathway of Chemicals with Heterogeneous Information. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2018 , 21, 670-680	1.3	16
95	Identification of age-related macular degeneration related genes by applying shortest path algorithm in protein-protein interaction network. <i>BioMed Research International</i> , 2013 , 2013, 523415	3	16
94	Predicting subcellular location of proteins using integrated-algorithm method. <i>Molecular Diversity</i> , 2010 , 14, 551-8	3.1	16

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93	Recognizing Novel Tumor Suppressor Genes Using a Network Machine Learning Strategy. <i>IEEE Access</i> , 2019 , 7, 155002-155013	3.5	16	
92	An integrated method for the identification of novel genes related to oral cancer. <i>PLoS ONE</i> , 2017 , 12, e0175185	3.7	15	
91	Prediction of drug target groups based on chemical-chemical similarities and chemical-chemical/protein connections. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2014 , 1844, 207-13	4	15	
90	Joint Access Selection and Resource Allocation in Cache-Enabled HCNs with D2D Communications 2017 ,		14	
89	Determination of Genes Related to Uveitis by Utilization of the Random Walk with Restart Algorithm on a Protein-Protein Interaction Network. <i>International Journal of Molecular Sciences</i> , 2017 , 18,	6.3	14	
88	A novel approach for fMRI data analysis based on the combination of sparse approximation and affinity propagation clustering. <i>Magnetic Resonance Imaging</i> , 2014 , 32, 736-46	3.3	14	
87	Prediction of multi-type membrane proteins in human by an integrated approach. <i>PLoS ONE</i> , 2014 , 9, e93553	3.7	14	
86	Prediction and analysis of retinoblastoma related genes through gene ontology and KEGG. <i>BioMed Research International</i> , 2013 , 2013, 304029	3	13	
85	Identifying chemicals with potential therapy of HIV based on protein-protein and protein-chemical interaction network. <i>PLoS ONE</i> , 2013 , 8, e65207	3.7	13	
84	Screening of Methylation Signature and Gene Functions Associated With the Subtypes of Isocitrate Dehydrogenase-Mutation Gliomas. <i>Frontiers in Bioengineering and Biotechnology</i> , 2019 , 7, 339	5.8	13	
83	Investigating the gene expression profiles of cells in seven embryonic stages with machine learning algorithms. <i>Genomics</i> , 2020 , 112, 2524-2534	4.3	12	
82	Hardness results and approximation algorithms for (weighted) paired-domination in graphs. <i>Theoretical Computer Science</i> , 2009 , 410, 5063-5071	1.1	12	
81	Discriminating between lysine sumoylation and lysine acetylation using mRMR feature selection and analysis. <i>PLoS ONE</i> , 2014 , 9, e107464	3.7	12	
80	A hybrid computational method for the discovery of novel reproduction-related genes. <i>PLoS ONE</i> , 2015 , 10, e0117090	3.7	12	
79	OPMSP: A Computational Method Integrating Protein Interaction and Sequence Information for the Identification of Novel Putative Oncogenes. <i>Protein and Peptide Letters</i> , 2016 , 23, 1081-1094	1.9	12	
78	Identification of COVID-19 Infection-Related Human Genes Based on a Random Walk Model in a Virus-Human Protein Interaction Network. <i>BioMed Research International</i> , 2020 , 2020, 4256301	3	12	
77	Prediction of cancer drugs by chemical-chemical interactions. <i>PLoS ONE</i> , 2014 , 9, e87791	3.7	11	
76	Prediction of interactiveness between small molecules and enzymes by combining gene ontology and compound similarity. <i>Journal of Computational Chemistry</i> , 2010 , 31, 1766-76	3.5	11	

75	Analysis of Gene Expression Profiles in the Human Brain Stem, Cerebellum and Cerebral Cortex. <i>PLoS ONE</i> , 2016 , 11, e0159395	3.7	11
74	A Binary Classifier for Prediction of the Types of Metabolic Pathway of Chemicals. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2017 , 20, 140-146	1.3	11
73	Analysis and Prediction of Myristoylation Sites Using the mRMR Method, the IFS Method and an Extreme Learning Machine Algorithm. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2017 , 20, 96-106	1.3	11
72	Predicting Biological Functions of Protein Complexes Using Graphic and Functional Features <i>Current Bioinformatics</i> , 2013 , 8, 545-551	4.7	11
71	iMPT-FDNPL: Identification of Membrane Protein Types with Functional Domains and a Natural Language Processing Approach. <i>Computational and Mathematical Methods in Medicine</i> , 2021 , 2021, 768	1497	11
70	Using compound similarity and functional domain composition for prediction of drug-target interaction networks. <i>Medicinal Chemistry</i> , 2010 , 6, 388-95	1.8	10
69	Predicting RNA 5-Methylcytosine Sites by Using Essential Sequence Features and Distributions BioMed Research International, 2022 , 2022, 4035462	3	10
68	Predicting Heart Cell Types by Using Transcriptome Profiles and a Machine Learning Method <i>Life</i> , 2022 , 12,	3	10
67	Prediction of Drug Combinations with a Network Embedding Method. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2018 , 21, 789-797	1.3	10
66	Investigation and Prediction of Human Interactome Based on Quantitative Features. <i>Frontiers in Bioengineering and Biotechnology</i> , 2020 , 8, 730	5.8	10
65	Prediction of Linear B-Cell Epitopes with mRMR Feature Selection and Analysis. <i>Current Bioinformatics</i> , 2016 , 11, 22-31	4.7	10
64	Analysis of the chemical toxicity effects using the enrichment of Gene Ontology terms and KEGG pathways. <i>Biochimica Et Biophysica Acta - General Subjects</i> , 2016 , 1860, 2619-26	4	10
63	A Novel Brain Networks Enhancement Model (BNEM) for BOLD fMRI Data Analysis With Highly Spatial Reproducibility. <i>IEEE Journal of Biomedical and Health Informatics</i> , 2016 , 20, 1107-19	7.2	9
62	Identifying Essential Signature Genes and Expression Rules Associated With Distinctive Development Stages of Early Embryonic Cells. <i>IEEE Access</i> , 2019 , 7, 128570-128578	3.5	9
61	(k)-Power domination in block graphs. <i>Journal of Combinatorial Optimization</i> , 2016 , 31, 865-873	0.9	9
60	Validating the performance of one-time decomposition for fMRI analysis using ICA with automatic target generation process. <i>Magnetic Resonance Imaging</i> , 2013 , 31, 970-5	3.3	9
59	A Computational Method for Classifying Different Human Tissues with Quantitatively Tissue-Specific Expressed Genes. <i>Genes</i> , 2018 , 9,	4.2	9
58	Discovery of new candidate genes related to brain development using protein interaction information. <i>PLoS ONE</i> , 2015 , 10, e0118003	3.7	8

57	Distinguishing Glioblastoma Subtypes by Methylation Signatures. Frontiers in Genetics, 2020, 11, 604330	54.5	8
56	Feature Classification and Analysis of Lung Cancer Related Genes through Gene Ontology and KEGG Pathways. <i>Current Bioinformatics</i> , 2016 , 11, 40-50	4.7	8
55	The Use of Gene Ontology Term and KEGG Pathway Enrichment for Analysis of Drug Half-Life. <i>PLoS ONE</i> , 2016 , 11, e0165496	3.7	8
54	Primary Tumor Site Specificity is Preserved in Patient-Derived Tumor Xenograft Models. <i>Frontiers in Genetics</i> , 2019 , 10, 738	4.5	7
53	Analysis of Protein-Protein Functional Associations by Using Gene Ontology and KEGG Pathway. <i>BioMed Research International</i> , 2019 , 2019, 4963289	3	7
52	Predicting triplet of transcription factor Imediating enzyme Itarget gene by functional profiles. <i>Neurocomputing</i> , 2011 , 74, 3677-3681	5.4	7
51	Identifying Robust Microbiota Signatures and Interpretable Rules to Distinguish Cancer Subtypes. <i>Frontiers in Molecular Biosciences</i> , 2020 , 7, 604794	5.6	7
50	Identifying the RNA signatures of coronary artery disease from combined lncRNA and mRNA expression profiles. <i>Genomics</i> , 2020 , 112, 4945-4958	4.3	7
49	Immunosignature Screening for Multiple Cancer Subtypes Based on Expression Rule. <i>Frontiers in Bioengineering and Biotechnology</i> , 2019 , 7, 370	5.8	7
48	Network-Based Method for Identifying Co- Regeneration Genes in Bone, Dentin, Nerve and Vessel Tissues. <i>Genes</i> , 2017 , 8,	4.2	6
47	Extremal problems on consecutive L(2,1)-labelling. <i>Discrete Applied Mathematics</i> , 2007 , 155, 1302-1313	1	6
46	The Use of Chemical-Chemical Interaction and Chemical Structure to Identify New Candidate Chemicals Related to Lung Cancer. <i>PLoS ONE</i> , 2015 , 10, e0128696	3.7	6
45	A Shortest-Path-Based Method for the Analysis and Prediction of Fruit-Related Genes in Arabidopsis thaliana. <i>PLoS ONE</i> , 2016 , 11, e0159519	3.7	6
44	Identifying the Signatures and Rules of Circulating Extracellular MicroRNA for Distinguishing Cancer Subtypes. <i>Frontiers in Genetics</i> , 2021 , 12, 651610	4.5	6
43	Analysis of Gene Expression Differences between Different Pancreatic Cells. ACS Omega, 2019, 4, 6421	6,4335	5
42	Recognizing and Predicting Thioether Bridges Formed by Lanthionine and EMethyllanthionine in Lantibiotics Using a Random Forest Approach with Feature Selection. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2017 , 20, 582-593	1.3	5
41	A path-based method for identification of protein phenotypic annotations. <i>Current Bioinformatics</i> , 2021 , 16,	4.7	5
40	iATC-NFMLP: Identifying classes of anatomical therapeutic chemicals based on drug networks, fingerprints and multilayer perceptron. <i>Current Bioinformatics</i> , 2022 , 17,	4.7	5

39	Inferring novel genes related to oral cancer with a network embedding method and one-class learning algorithms. <i>Gene Therapy</i> , 2019 , 26, 465-478	4	4
38	Alternative Polyadenylation Modification Patterns Reveal Essential Posttranscription Regulatory Mechanisms of Tumorigenesis in Multiple Tumor Types. <i>BioMed Research International</i> , 2020 , 2020, 638	43120	4
37	Deciphering the Relationship between Obesity and Various Diseases from a Network Perspective. <i>Genes</i> , 2017 , 8,	4.2	4
36	Selection of reprogramming factors of induced pluripotent stem cells based on the protein interaction network and functional profiles. <i>Protein and Peptide Letters</i> , 2012 , 19, 113-9	1.9	4
35	Prediction of drugs target groups based on ChEBI ontology. <i>BioMed Research International</i> , 2013 , 2013, 132724	3	4
34	A two-step similarity-based method for prediction of drug's target group. <i>Protein and Peptide Letters</i> , 2013 , 20, 364-70	1.9	4
33	Predicting Protein Ligand Binding Sites with Structure Alignment Method on Hadoop. <i>Current Proteomics</i> , 2016 , 13, 113-121	0.7	4
32	A Fast Approximate Hypervolume Calculation Method by a Novel Decomposition Strategy. <i>Lecture Notes in Computer Science</i> , 2017 , 14-25	0.9	4
31	Predicting gene phenotype by multi-label multi-class model based on essential functional features. <i>Molecular Genetics and Genomics</i> , 2021 , 296, 905-918	3.1	4
30	Identification of Cell Markers and Their Expression Patterns in Skin Based on Single-Cell RNA-Sequencing Profiles <i>Life</i> , 2022 , 12,	3	4
29	Distance paired-domination problems on subclasses of chordal graphs. <i>Theoretical Computer Science</i> , 2009 , 410, 5072-5081	1.1	3
28	Analysis of metabolic pathway using hybrid properties. <i>Protein and Peptide Letters</i> , 2012 , 19, 99-107	1.9	3
27	Study of drug-drug combinations based on molecular descriptors and physicochemical properties. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2016 , 19, 153-60	1.3	3
26	miR-29 mediates exercise-induced skeletal muscle angiogenesis by targeting VEGFA, COL4A1 and COL4A2 via the PI3K/Akt signaling pathway. <i>Molecular Medicine Reports</i> , 2020 , 22, 661-670	2.9	3
25	An Integrated Multi-Label Classifier with Chemical-Chemical Interactions for Prediction of Chemical Toxicity Effects. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2018 , 21, 403-410	1.3	3
24	Prediction of membrane protein types by fusing protein-protein interaction and protein sequence information. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2020 , 1868, 140524	4	3
23	Recognizing novel chemicals/drugs for anatomical therapeutic chemical classes with a heat diffusion algorithm. <i>Biochimica Et Biophysica Acta - Molecular Basis of Disease</i> , 2020 , 1866, 165910	6.9	3
22	Drug-Drug Interactions Prediction Using Fingerprint Only <i>Computational and Mathematical Methods in Medicine</i> , 2022 , 2022, 7818480	2.8	3

21	NP-completeness and APX-completeness of restrained domination in graphs. <i>Theoretical Computer Science</i> , 2012 , 448, 1-8	1.1	2
20	Identification of Pan-Cancer Biomarkers Based on the Gene Expression Profiles of Cancer Cell Lines <i>Frontiers in Cell and Developmental Biology</i> , 2021 , 9, 781285	5.7	2
19	Computational method for distinguishing lysine acetylation, sumoylation, and ubiquitination using the random forest algorithm with a feature selection procedure. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2017 ,	1.3	2
18	iMPT-FRAKEL: A Simple Multi-label Web-server that Only Uses Fingerprints to Identify which Metabolic Pathway Types Compounds can Participate In. <i>Open Bioinformatics Journal</i> , 2020 , 13, 83-91	0.8	2
17	Identification of Latent Oncogenes with a Network Embedding Method and Random Forest. <i>BioMed Research International</i> , 2020 , 2020, 5160396	3	2
16	Identifying COVID-19-Specific Transcriptomic Biomarkers with Machine Learning Methods. <i>BioMed Research International</i> , 2021 , 2021, 9939134	3	2
15	Identification of Microbiota Biomarkers With Orthologous Gene Annotation for Type 2 Diabetes. <i>Frontiers in Microbiology</i> , 2021 , 12, 711244	5.7	2
14	Identification of Gene Signatures and Expression Patterns During Epithelial-to-Mesenchymal Transition From Single-Cell Expression Atlas. <i>Frontiers in Genetics</i> , 2020 , 11, 605012	4.5	2
13	Identification of Carcinogenic Chemicals with Network Embedding and Deep Learning Methods. <i>Current Bioinformatics</i> , 2021 , 15, 1017-1026	4.7	2
12	Similarity-Based Method with Multiple-Feature Sampling for Predicting Drug Side Effects <i>Computational and Mathematical Methods in Medicine</i> , 2022 , 2022, 9547317	2.8	2
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10	Vertices in all minimum paired-dominating sets of block graphs. <i>Journal of Combinatorial Optimization</i> , 2012 , 24, 176-191	0.9	1
9	A computational method for the identification of candidate drugs for non-small cell lung cancer. <i>PLoS ONE</i> , 2017 , 12, e0183411	3.7	1
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5	Recognition of Immune Cell Markers of COVID-19 Severity with Machine Learning Methods <i>BioMed Research International</i> , 2022 , 2022, 6089242	3	О
4	A Two-step Similarity-based Method for Prediction of Drug's Target Group. <i>Protein and Peptide Letters</i> , 2013 , 20, 364-370	1.9	

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