

Cang-Zhi Jia

List of Publications by Year in descending order

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Version: 2024-02-01

37
papers

1,459
citations

361413

20
h-index

330143

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37
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docs citations

37
times ranked

969
citing authors

#	ARTICLE	IF	CITATIONS
1	4mCPred: machine learning methods for DNA N4-methylcytosine sites prediction. <i>Bioinformatics</i> , 2019, 35, 593-601.	4.1	146
2	O-GlcNAcPRED-II: an integrated classification algorithm for identifying O-GlcNAcylation sites based on fuzzy undersampling and a <i>K</i> -means PCA oversampling technique. <i>Bioinformatics</i> , 2018, 34, 2029-2036.	4.1	128
3	MULTiPly: a novel multi-layer predictor for discovering general and specific types of promoters. <i>Bioinformatics</i> , 2019, 35, 2957-2965.	4.1	109
4	Prediction of Protein S-Nitrosylation Sites Based on Adapted Normal Distribution Bi-Profile Bayes and Chou's Pseudo Amino Acid Composition. <i>International Journal of Molecular Sciences</i> , 2014, 15, 10410-10423.	4.1	94
5	An Interpretable Prediction Model for Identifying N7-Methylguanosine Sites Based on XGBoost and SHAP. <i>Molecular Therapy - Nucleic Acids</i> , 2020, 22, 362-372.	5.1	93
6	DeepTorrent: a deep learning-based approach for predicting DNA N4-methylcytosine sites. <i>Briefings in Bioinformatics</i> , 2021, 22, .	6.5	84
7	EnhancerPred: a predictor for discovering enhancers based on the combination and selection of multiple features. <i>Scientific Reports</i> , 2016, 6, 38741.	3.3	80
8	70ProPred: a predictor for discovering sigma70 promoters based on combining multiple features. <i>BMC Systems Biology</i> , 2018, 12, 44.	3.0	72
9	PASSION: an ensemble neural network approach for identifying the binding sites of RBPs on circRNAs. <i>Bioinformatics</i> , 2020, 36, 4276-4282.	4.1	58
10	RNA-MethylPred: A high-accuracy predictor to identify N6-methyladenosine in RNA. <i>Analytical Biochemistry</i> , 2016, 510, 72-75.	2.4	56
11	S-SulfPred: A sensitive predictor to capture S-sulfenylation sites based on a resampling one-sided selection undersampling-synthetic minority oversampling technique. <i>Journal of Theoretical Biology</i> , 2017, 422, 84-89.	1.7	52
12	O-GlcNAcPRED: a sensitive predictor to capture protein O-GlcNAcylation sites. <i>Molecular BioSystems</i> , 2013, 9, 2909.	2.9	45
13	Computational identification of eukaryotic promoters based on cascaded deep capsule neural networks. <i>Briefings in Bioinformatics</i> , 2021, 22, .	6.5	44
14	EnhancerPred2.0: predicting enhancers and their strength based on position-specific trinucleotide propensity and electron-ion interaction potential feature selection. <i>Molecular BioSystems</i> , 2017, 13, 767-774.	2.9	42
15	Inspector: a lysine succinylation predictor based on edited nearest-neighbor undersampling and adaptive synthetic oversampling. <i>Analytical Biochemistry</i> , 2020, 593, 113592.	2.4	40
16	PRISMOID: a comprehensive 3D structure database for post-translational modifications and mutations with functional impact. <i>Briefings in Bioinformatics</i> , 2020, 21, 1069-1079.	6.5	38
17	NucPosPred: Predicting species-specific genomic nucleosome positioning via four different modes of general PseKNC. <i>Journal of Theoretical Biology</i> , 2018, 450, 15-21.	1.7	35
18	Prediction of mitochondrial proteins of malaria parasite using bi-profile Bayes feature extraction. <i>Biochimie</i> , 2011, 93, 778-782.	2.6	34

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19	Harmonic number identities via the Newton–Andrews method. <i>Ramanujan Journal</i> , 2014, 35, 263-285.	0.7	32
20	Positive-unlabeled learning in bioinformatics and computational biology: a brief review. <i>Briefings in Bioinformatics</i> , 2022, 23, .	6.5	26
21	Prediction of aptamer–protein interacting pairs based on sparse autoencoder feature extraction and an ensemble classifier. <i>Mathematical Biosciences</i> , 2019, 311, 103-108.	1.9	22
22	Staem5: A novel computational approach for accurate prediction of m5C site. <i>Molecular Therapy - Nucleic Acids</i> , 2021, 26, 1027-1034.	5.1	20
23	Formator: Predicting Lysine Formylation Sites Based on the Most Distant Undersampling and Safe-Level Synthetic Minority Oversampling. <i>IEEE/ACM Transactions on Computational Biology and Bioinformatics</i> , 2021, 18, 1937-1945.	3.0	17
24	OH-PRED: prediction of protein hydroxylation sites by incorporating adapted normal distribution bi-profile Bayes feature extraction and physicochemical properties of amino acids. <i>Journal of Biomolecular Structure and Dynamics</i> , 2017, 35, 829-835.	3.5	16
25	EnsemPseU: Identifying Pseudouridine Sites With an Ensemble Approach. <i>IEEE Access</i> , 2020, 8, 79376-79382.	4.2	15
26	CarSite: identifying carbonylated sites of human proteins based on a one-sided selection resampling method. <i>Molecular BioSystems</i> , 2017, 13, 2362-2369.	2.9	13
27	Critical assessment of computational tools for prokaryotic and eukaryotic promoter prediction. <i>Briefings in Bioinformatics</i> , 2022, 23, .	6.5	11
28	Alignment-free Comparison of Protein Sequences Based on Reduced Amino Acid Alphabets. <i>Journal of Biomolecular Structure and Dynamics</i> , 2009, 26, 763-769.	3.5	10
29	DephosSitePred: A High Accuracy Predictor for Protein Dephosphorylation Sites. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2017, 20, 153-157.	1.1	6
30	Computational Prediction of Protein O-GlcNAc Modification. <i>Methods in Molecular Biology</i> , 2018, 1754, 235-246.	0.9	5
31	TRANSFORMATION AND REDUCTION FORMULAE FOR DOUBLE q -SERIES OF TYPE $\hat{1} \uparrow \langle \sup \rangle 2:1; \hat{1} \langle \sub \rangle 2:0; \hat{1} \uparrow \langle \sub \rangle 4$. <i>Glasgow Mathematical Journal</i> , 2010, 52, 195-204.	0.3	3
32	Quartic theta hypergeometric series. <i>Ramanujan Journal</i> , 2013, 32, 23-81.	0.7	3
33	Recognition of Protein Pupylation Sites by Adopting Resampling Approach. <i>Molecules</i> , 2018, 23, 3097.	3.8	3
34	O-glycosylation site prediction for <i>Homo sapiens</i> by combining properties and sequence features with support vector machine. <i>Journal of Bioinformatics and Computational Biology</i> , 2022, 20, 2150029.	0.8	3
35	Protein secondary structure class assignment on the basis of a new graphic representation. <i>International Journal of Quantum Chemistry</i> , 2009, 109, 819-825.	2.0	2
36	ABEL'S METHOD ON SUMMATION BY PARTS FOR ELLIPTIC HYPERGEOMETRIC SERIES. <i>Communications in Contemporary Mathematics</i> , 2009, 11, 337-353.	1.2	1

#	ARTICLE	IF	CITATIONS
37	Pippin: A random forest-based method for identifying presynaptic and postsynaptic neurotoxins. Journal of Bioinformatics and Computational Biology, 2020, 18, 2050008.	0.8	1