

Ke Chen

List of Publications by Year in descending order

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Version: 2024-02-01

27
papers

1,193
citations

687220

13
h-index

677027

22
g-index

27
all docs

27
docs citations

27
times ranked

1217
citing authors

#	ARTICLE	IF	CITATIONS
1	Computational Prediction of Secondary and Supersecondary Structures from Protein Sequences. <i>Methods in Molecular Biology</i> , 2019, 1958, 73-100.	0.4	11
2	Accurate prediction of secondary structure of tRNAs. <i>Biochemical and Biophysical Research Communications</i> , 2019, 509, 64-68.	1.0	1
3	A statistical analysis on transcriptome sequences: The enrichment of Alu-element is associated with subcellular location. <i>Biochemical and Biophysical Research Communications</i> , 2018, 499, 397-402.	1.0	9
4	Pairwise alignment for very long nucleic acid sequences. <i>Biochemical and Biophysical Research Communications</i> , 2018, 502, 313-317.	1.0	9
5	NSiteMatch: Prediction of Binding Sites of Nucleotides by Identifying the Structure Similarity of Local Surface Patches. <i>Computational and Mathematical Methods in Medicine</i> , 2017, 2017, 1-16.	0.7	8
6	Systematic investigation of sequence and structural motifs that recognize ATP. <i>Computational Biology and Chemistry</i> , 2015, 56, 131-141.	1.1	3
7	Quad-PRE: A Hybrid Method to Predict Protein Quaternary Structure Attributes. <i>Computational and Mathematical Methods in Medicine</i> , 2014, 2014, 1-9.	0.7	3
8	newDNA-Prot: Prediction of DNA-binding proteins by employing support vector machine and a comprehensive sequence representation. <i>Computational Biology and Chemistry</i> , 2014, 52, 51-59.	1.1	20
9	A matrix grammar approach for automatic distributed network resource management. <i>Frontiers of Computer Science</i> , 2013, 7, 583-594.	1.6	1
10	Prediction and analysis of nucleotide-binding residues using sequence and sequence-derived structural descriptors. <i>Bioinformatics</i> , 2012, 28, 331-341.	1.8	106
11	Computational Prediction of Secondary and Supersecondary Structures. <i>Methods in Molecular Biology</i> , 2012, 932, 63-86.	0.4	13
12	A Critical Comparative Assessment of Predictions of Protein-Binding Sites for Biologically Relevant Organic Compounds. <i>Structure</i> , 2011, 19, 613-621.	1.6	59
13	iFC2: an integrated web-server for improved prediction of protein structural class, fold type, and secondary structure content. <i>Amino Acids</i> , 2011, 40, 963-973.	1.2	13
14	ATPsite: sequence-based prediction of ATP-binding residues. <i>Proteome Science</i> , 2011, 9, S4.	0.7	57
15	Accurate prediction of ATP-binding residues using sequence and sequence-derived structural descriptors. , 2010, , .		0
16	Prediction of integral membrane protein type by collocated hydrophobic amino acid pairs. <i>Journal of Computational Chemistry</i> , 2009, 30, 163-172.	1.5	66
17	Fuzzy Filtering of Nonlinear Systems With Intermittent Measurements. <i>IEEE Transactions on Fuzzy Systems</i> , 2009, 17, 291-300.	6.5	267
18	Investigation of Atomic Level Patterns in Protein-Small Ligand Interactions. <i>PLoS ONE</i> , 2009, 4, e4473.	1.1	55

#	ARTICLE	IF	CITATIONS
19	Prediction of protein structural class using novel evolutionary collocation-based sequence representation. <i>Journal of Computational Chemistry</i> , 2008, 29, 1596-1604.	1.5	148
20	Identification of tubulin drug binding sites and prediction of relative differences in binding affinities to tubulin isotypes using digital signal processing. <i>Journal of Molecular Graphics and Modelling</i> , 2008, 27, 497-505.	1.3	15
21	PFRES: protein fold classification by using evolutionary information and predicted secondary structure. <i>Bioinformatics</i> , 2007, 23, 2843-2850.	1.8	108
22	Prediction of protein crystallization using collocation of amino acid pairs. <i>Biochemical and Biophysical Research Communications</i> , 2007, 355, 764-769.	1.0	102
23	Improved Prediction of Relative Solvent Accessibility Using Two-stage Support Vector Regression. , 2007, , .		0
24	Prediction of Protein Structural Class Using PSI-BLAST Profile Based Collocation of Amino Acid Pairs. , 2007, , .		1
25	Prediction of flexible/rigid regions from protein sequences using k-spaced amino acid pairs. <i>BMC Structural Biology</i> , 2007, 7, 25.	2.3	100
26	Optimization of the Sliding Window Size for Protein Structure Prediction. , 2006, , .		18
27	Prediction of the Number of Helices for the Twilight Zone Proteins. , 2006, , .		0