

Joseph Rebehmed

List of Publications by Year in descending order

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Version: 2024-02-01

23
papers

470
citations

840776

11
h-index

713466

21
g-index

23
all docs

23
docs citations

23
times ranked

890
citing authors

#	ARTICLE	IF	CITATIONS
1	Protein flexibility in the light of structural alphabets. <i>Frontiers in Molecular Biosciences</i> , 2015, 2, 20.	3.5	71
2	Molecular modeling studies of N-substituted pyrrole derivativesâ€”Potential HIV-1 gp41 inhibitors. <i>Bioorganic and Medicinal Chemistry</i> , 2008, 16, 3039-3048.	3.0	68
3	Cisâ€”trans isomerization of omega dihedrals in proteins. <i>Amino Acids</i> , 2013, 45, 279-289.	2.7	60
4	PTM-SD: a database of structurally resolved and annotated posttranslational modifications in proteins. <i>Database: the Journal of Biological Databases and Curation</i> , 2014, 2014, bau041-bau041.	3.0	44
5	Î²â€”Bulges: Extensive structural analyses of Î²â€”sheets irregularities. <i>Protein Science</i> , 2013, 22, 1366-1378.	7.6	27
6	Flexible computational docking studies of new aminoglycosides targeting RNA 16S bacterial ribosome site. <i>European Journal of Medicinal Chemistry</i> , 2008, 43, 1648-1656.	5.5	26
7	Discrete analyses of protein dynamics. <i>Journal of Biomolecular Structure and Dynamics</i> , 2020, 38, 2988-3002.	3.5	20
8	Evolution study of the Baeyerâ€”Villiger monooxygenases enzyme family: Functional importance of the highly conserved residues. <i>Biochimie</i> , 2013, 95, 1394-1402.	2.6	19
9	Expanding the SRI domain family: A common scaffold for binding the phosphorylated Câ€”terminal domain of RNA polymerase II. <i>FEBS Letters</i> , 2014, 588, 4431-4437.	2.8	19
10	Investigation of the impact of PTMs on the protein backbone conformation. <i>Amino Acids</i> , 2019, 51, 1065-1079.	2.7	17
11	In silico prediction of protein flexibility with local structure approach. <i>Biochimie</i> , 2019, 165, 150-155.	2.6	15
12	iPBAvizu: a PyMOL plugin for an efficient 3D protein structure superimposition approach. <i>Source Code for Biology and Medicine</i> , 2019, 14, 5.	1.7	14
13	The respective roles of polar/nonpolar binary patterns and amino acid composition in protein regular secondary structures explored exhaustively using hydrophobic cluster analysis. <i>Proteins: Structure, Function and Bioinformatics</i> , 2016, 84, 624-638.	2.6	12
14	A topology-based investigation of protein interaction sites using Hydrophobic Cluster Analysis. <i>Biochimie</i> , 2019, 167, 68-80.	2.6	12
15	2D and 3D QSAR studies of diarylpyrimidine HIV-1 reverse transcriptase inhibitors. <i>Journal of Computer-Aided Molecular Design</i> , 2008, 22, 831-841.	2.9	10
16	Dynamics and deformability of Î±-, 310- and Î³-helices. <i>Archives of Biological Sciences</i> , 2018, 70, 21-31.	0.5	10
17	Current status of PTMs structural databases: applications, limitations and prospects. <i>Amino Acids</i> , 2022, 54, 575-590.	2.7	9
18	Enhancing the Activity of a <i>Dietzia</i> sp. D5 Baeyerâ€”Villiger Monooxygenase towards Cyclohexanone by Saturation Mutagenesis. <i>ChemistrySelect</i> , 2017, 2, 7169-7177.	1.5	7

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19	Enzymatic Synthesis of Galactosylated Serine/Threonine Derivatives by β -Galactosidase from <i>Escherichia coli</i> . <i>International Journal of Molecular Sciences</i> , 2015, 16, 13714-13728.	4.1	4
20	Shaking the β -Bulges. <i>IEEE/ACM Transactions on Computational Biology and Bioinformatics</i> , 2022, 19, 14-18.	3.0	3
21	Editorial: Advances in Molecular Docking and Structure-Based Modelling. <i>Frontiers in Molecular Biosciences</i> , 2022, 9, 839826.	3.5	2
22	Combining In Silico Phylogenetic and Threading Approaches to Assist the in vitro Protein Engineering of BVMO Enzymes. <i>Biophysical Journal</i> , 2020, 118, 45a.	0.5	1
23	Efficient Protein Structure Alignment Methods Based on a Structural Alphabet. <i>Biophysical Journal</i> , 2014, 106, 54a.	0.5	0