

Indira Ghosh

List of Publications by Year in descending order

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28
papers

484
citations

840776

11
h-index

677142

22
g-index

30
all docs

30
docs citations

30
times ranked

593
citing authors

#	ARTICLE	IF	CITATIONS
1	Application and Comprehensive Analysis of Neighbor Approximated Information Theoretic Configurational Entropy Methods to Protein-Ligand Binding Cases. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 7581-7600.	5.3	2
2	Coupling Supervised Molecular Dynamics (SuMD) with Entropy Estimations To Shine Light on the Stability of Multiple Binding Sites. <i>ACS Medicinal Chemistry Letters</i> , 2019, 10, 444-449.	2.8	5
3	In silico identification and characterization of stress and virulence associated repeats in <i>Salmonella</i> . <i>Genomics</i> , 2018, 110, 23-34.	2.9	5
4	ILP-assisted de novo drug design. <i>Machine Learning</i> , 2016, 103, 309-341.	5.4	11
5	Modularity in protein structures: study on all-alpha proteins. <i>Journal of Biomolecular Structure and Dynamics</i> , 2015, 33, 2667-2681.	3.5	16
6	PfalDB: An Integrated Drug Target and Chemical Database for <i>Plasmodium falciparum</i> . <i>Current Drug Targets</i> , 2014, 15, 1089-1093.	2.1	1
7	Methylerythritol phosphate pathway to isoprenoids: Kinetic modeling and <i>in silico</i> enzyme inhibitions in <i>Plasmodium falciparum</i> . <i>FEBS Letters</i> , 2013, 587, 2806-2817.	2.8	15
8	Concepts on the protein folding problem. <i>Journal of Biomolecular Structure and Dynamics</i> , 2013, 31, 1020-1023.	3.5	8
9	Modeling and Experimental Analyses Reveals Signaling Plasticity in a Bi-Modular Assembly of CD40 Receptor Activated Kinases. <i>PLoS ONE</i> , 2012, 7, e39898.	2.5	17
10	Proteins will Fold Anyway!!. <i>Journal of Biomolecular Structure and Dynamics</i> , 2011, 28, 627-628.	3.5	2
11	Mathematical criteria to observe mesoscopic emergence of protein biochemical properties. <i>Journal of Mathematical Chemistry</i> , 2011, 49, 643-665.	1.5	5
12	Fractal symmetry of protein interior: what have we learned?. <i>Cellular and Molecular Life Sciences</i> , 2011, 68, 2711-2737.	5.4	33
13	Characterization of <i>Leishmania donovani</i> Aquaporins Shows Presence of Subcellular Aquaporins Similar to Tonoplast Intrinsic Proteins of Plants. <i>PLoS ONE</i> , 2011, 6, e24820.	2.5	16
14	Analysis of physico-chemical properties of substrates of ABC and MFS multidrug transporters of pathogenic <i>Candida albicans</i> . <i>European Journal of Medicinal Chemistry</i> , 2010, 45, 4813-4826.	5.5	31
15	Structure and Dynamics of Double Helical DNA in Torsion Angle Hyperspace: A Molecular Mechanics Approach. <i>Journal of Biomolecular Structure and Dynamics</i> , 2010, 27, 695-712.	3.5	11
16	Revisiting the Myths of Protein Interior: Studying Proteins with Mass-Fractal Hydrophobicity-Fractal and Polarizability-Fractal Dimensions. <i>PLoS ONE</i> , 2009, 4, e7361.	2.5	25
17	A new computational model to study mass inhomogeneity and hydrophobicity inhomogeneity in proteins. <i>European Biophysics Journal</i> , 2009, 38, 577-587.	2.2	11
18	Determination of Phosphorylation Sites for NADP-specific Isocitrate Dehydrogenase from <i>Mycobacterium tuberculosis</i> . <i>Journal of Biomolecular Structure and Dynamics</i> , 2009, 26, 741-754.	3.5	15

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19	Simple sequence repeats in different genome sequences of Shigella and comparison with high GC and AT-rich genomes. DNA Sequence, 2008, 19, 167-176.	0.7	3
20	Mapping Selectivity and Specificity of Active Site of Plasmepsins from Plasmodium falciparum Using Molecular Interaction Field Approach. Protein and Peptide Letters, 2007, 14, 569-574.	0.9	7
21	Developing an Antituberculosis Compounds Database and Data Mining in the Search of a Motif Responsible for the Activity of a Diverse Class of Antituberculosis Agents. Journal of Chemical Information and Modeling, 2006, 46, 17-23.	5.4	23
22	Biochemical characterization of Plasmodium falciparum hypoxanthine-guanine-xanthine phosphoribosyltransferase: role of histidine residue in substrate selectivity. Molecular and Biochemical Parasitology, 2004, 137, 267-276.	1.1	11
23	Analysis of E.coli promoter structures using neural networks. Nucleic Acids Research, 1994, 22, 2158-2165.	14.5	51
24	Molecular dynamics study of the binding of phenylalanine stereoisomers to thermolysin. Biophysical Chemistry, 1994, 50, 237-248.	2.8	8
25	Hysteresis and Statistical Errors in Free Energy Perturbation L to D Amino Acid Conversion. Molecular Simulation, 1993, 10, 241-253.	2.0	6
26	Hydration of chloride and bromide anions: determination of relative free energy by computer simulation. Journal of the American Chemical Society, 1985, 107, 7793-7794.	13.7	135
27	Effect of Configuration of the Inhibitors on the Mode of Binding to the Enzyme, Thermolysin. Journal of Biomolecular Structure and Dynamics, 1984, 2, 29-40.	3.5	2
28	A conformational approach to the study of the dynamics of enzyme inhibition: studies on thermolysin. International Journal of Biological Macromolecules, 1982, 4, 130-136.	7.5	7