

Nagasuma Chandra

List of Publications by Year in descending order

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106
papers

2,835
citations

159358

30
h-index

223531

46
g-index

109
all docs

109
docs citations

109
times ranked

4050
citing authors

#	ARTICLE	IF	CITATIONS
1	A new blood-based RNA signature (R9), for monitoring effectiveness of tuberculosis treatment in a South Indian longitudinal cohort. <i>IScience</i> , 2022, 25, 103745.	1.9	1
2	SiteMotif: A graph-based algorithm for deriving structural motifs in Protein Ligand binding sites. <i>PLoS Computational Biology</i> , 2022, 18, e1009901.	1.5	7
3	Functional and Biochemical Characterization of the MazEF6 Toxin-Antitoxin System of <i>Mycobacterium tuberculosis</i> . <i>Journal of Bacteriology</i> , 2022, 204, e0005822.	1.0	8
4	PathExt: a general framework for path-based mining of omics-integrated biological networks. <i>Bioinformatics</i> , 2021, 37, 1254-1262.	1.8	11
5	VB10, a new blood biomarker for differential diagnosis and recovery monitoring of acute viral and bacterial infections. <i>EBioMedicine</i> , 2021, 67, 103352.	2.7	15
6	A Strategic Target Rescues Trimethoprim Sensitivity in <i>Escherichia coli</i> . <i>IScience</i> , 2020, 23, 100986.	1.9	15
7	Î©76: A designed antimicrobial peptide to combat carbapenem- and tigecycline-resistant <i>Acinetobacter baumannii</i> . <i>Science Advances</i> , 2019, 5, eaax1946.	4.7	64
8	Interrogation of genome-wide networks in biology: comparison of knowledge-based and statistical methods. <i>International Journal of Advances in Engineering Sciences and Applied Mathematics</i> , 2019, 11, 119-137.	0.7	5
9	An Augmented Pocketome: Detection and Analysis of Small-Molecule Binding Pockets in Proteins of Known 3D Structure. <i>Structure</i> , 2018, 26, 499-512.e2.	1.6	38
10	Computational antimicrobial peptide design and evaluation against multidrug-resistant clinical isolates of bacteria. <i>Journal of Biological Chemistry</i> , 2018, 293, 3492-3509.	1.6	93
11	Design of a heme-binding peptide motif adopting a Î²-hairpin conformation. <i>Journal of Biological Chemistry</i> , 2018, 293, 9412-9422.	1.6	8
12	Identification of a coâ€target for enhancing efficacy of sorafenib in HCC through a quantitative modeling approach. <i>FEBS Journal</i> , 2018, 285, 3977-3992.	2.2	8
13	Role of genetic heterogeneity in determining the epidemiological severity of H1N1 influenza. <i>PLoS Computational Biology</i> , 2018, 14, e1006069.	1.5	14
14	Unbiased Identification of Blood-based Biomarkers for Pulmonary Tuberculosis by Modeling and Mining Molecular Interaction Networks. <i>EBioMedicine</i> , 2017, 15, 112-126.	2.7	75
15	A Systems Perspective of Signalling Networks in Hostâ€Pathogen Interactions. <i>Journal of the Indian Institute of Science</i> , 2017, 97, 41-57.	0.9	0
16	Deciphering common recognition principles of nucleoside mono/di and tri-phosphates binding in diverse proteins via structural matching of their binding sites. <i>Proteins: Structure, Function and Bioinformatics</i> , 2017, 85, 1699-1712.	1.5	3
17	Resolving protein structureâ€functionâ€binding site relationships from a binding site similarity network perspective. <i>Proteins: Structure, Function and Bioinformatics</i> , 2017, 85, 1319-1335.	1.5	14
18	A genome-wide structure-based survey of nucleotide binding proteins in <i>M. tuberculosis</i> . <i>Scientific Reports</i> , 2017, 7, 12489.	1.6	5

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19	Methionine synthase is localized to the nucleus in <i>Pichia pastoris</i> and <i>Candida albicans</i> and to the cytoplasm in <i>Saccharomyces cerevisiae</i> . <i>Journal of Biological Chemistry</i> , 2017, 292, 14730-14746.	1.6	9
20	DenHunt - A Comprehensive Database of the Intricate Network of Dengue-Human Interactions. <i>PLoS Neglected Tropical Diseases</i> , 2016, 10, e0004965.	1.3	23
21	Structural analysis of dihydrofolate reductases enables rationalization of antifolate binding affinities and suggests repurposing possibilities. <i>FEBS Journal</i> , 2016, 283, 1139-1167.	2.2	19
22	Dissecting the structural and functional features of the Luteinizing hormone receptor using receptor specific single chain fragment variables. <i>Molecular and Cellular Endocrinology</i> , 2016, 427, 1-12.	1.6	5
23	A comparative analysis of the DNA recombination repair pathway in mycobacterial genomes. <i>Tuberculosis</i> , 2016, 99, 109-119.	0.8	27
24	Curcumin Reduces the Motility of <i>Salmonella enterica</i> Serovar Typhimurium by Binding to the Flagella, Thereby Leading to Flagellar Fragility and Shedding. <i>Journal of Bacteriology</i> , 2016, 198, 1798-1811.	1.0	27
25	Editorial overview: Carbohydrate-protein interactions and glycosylation: integrating structural biology, informatics and systems modelling to understand glycan structure and glycan-protein interactions. <i>Current Opinion in Structural Biology</i> , 2016, 40, v-viii.	2.6	1
26	Identifying and Tackling Emergent Vulnerability in Drug-Resistant Mycobacteria. <i>ACS Infectious Diseases</i> , 2016, 2, 592-607.	1.8	34
27	EpiTracer - an algorithm for identifying epicenters in condition-specific biological networks. <i>BMC Genomics</i> , 2016, 17, 543.	1.2	17
28	Hypothetical protein Rv3423.1 of <i>Mycobacterium tuberculosis</i> is a histone acetyltransferase. <i>FEBS Journal</i> , 2016, 283, 265-281.	2.2	51
29	High IL-6 and low IL-15 levels mark the presence of TB infection: A preliminary study. <i>Cytokine</i> , 2016, 81, 57-62.	1.4	22
30	Probing the Druggability Limits for Enzymes of the NAD Biosynthetic Network in Glioma. <i>Journal of Chemical Information and Modeling</i> , 2016, 56, 843-853.	2.5	4
31	Complete Genome Sequences of an <i>Escherichia coli</i> Laboratory Strain and Trimethoprim-Resistant (TMP32XR) Mutant Strains. <i>Genome Announcements</i> , 2015, 3, .	0.8	4
32	Mechanism of Iron-Dependent Repressor (IdeR) Activation and DNA Binding: A Molecular Dynamics and Protein Structure Network Study. <i>PLoS Computational Biology</i> , 2015, 11, e1004500.	1.5	13
33	NrichD database: sequence databases enriched with computationally designed protein-like sequences aid in remote homology detection. <i>Nucleic Acids Research</i> , 2015, 43, D300-D305.	6.5	12
34	Enriching the annotation of <i>Mycobacterium tuberculosis</i> H37Rv proteome using remote homology detection approaches: Insights into structure and function. <i>Tuberculosis</i> , 2015, 95, 14-25.	0.8	9
35	Architectural plan of transcriptional regulation in <i>Mycobacterium tuberculosis</i> . <i>Trends in Microbiology</i> , 2015, 23, 123-125.	3.5	1
36	A molecular systems approach to modelling human skin pigmentation: identifying underlying pathways and critical components. <i>BMC Research Notes</i> , 2015, 8, 170.	0.6	9

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37	Deciphering complex patterns of class I HLA peptide cross-reactivity via hierarchical grouping. <i>Immunology and Cell Biology</i> , 2015, 93, 522-532.	1.0	9
38	Recognizing drug targets using evolutionary information: implications for repurposing FDA-approved drugs against <i>Mycobacterium tuberculosis</i> H37Rv. <i>Molecular BioSystems</i> , 2015, 11, 3316-3331.	2.9	20
39	De-DUFing the DUFs: Deciphering distant evolutionary relationships of Domains of Unknown Function using sensitive homology detection methods. <i>Biology Direct</i> , 2015, 10, 38.	1.9	34
40	Complete Genome Sequences of a <i>Mycobacterium smegmatis</i> Laboratory Strain (MC ² 155) and Isoniazid-Resistant (4XR1/R2) Mutant Strains. <i>Genome Announcements</i> , 2015, 3, .	0.8	54
41	Gene expression profiles of wild-type and isoniazid-resistant strains of <i>Mycobacterium smegmatis</i> . <i>Data in Brief</i> , 2015, 4, 186-189.	0.5	5
42	CHEXVIS: a tool for molecular channel extraction and visualization. <i>BMC Bioinformatics</i> , 2015, 16, 119.	1.2	74
43	Common recognition principles across diverse sequence and structural families of sialic acid binding proteins. <i>Glycobiology</i> , 2014, 24, 5-16.	1.3	7
44	Mapping Post-translational Modifications of Mammalian Testicular Specific Histone Variant TH2B in Tetraploid and Haploid Germ Cells and Their Implications on the Dynamics of Nucleosome Structure. <i>Journal of Proteome Research</i> , 2014, 13, 5603-5617.	1.8	20
45	Identifying feasible metabolic routes in <i>Mycobacterium smegmatis</i> and possible alterations under diverse nutrient conditions. <i>BMC Microbiology</i> , 2014, 14, 276.	1.3	22
46	Weighting schemes in metabolic graphs for identifying biochemical routes. <i>Systems and Synthetic Biology</i> , 2014, 8, 47-57.	1.0	2
47	Filling-in Void and Sparse Regions in Protein Sequence Space by Protein-Like Artificial Sequences Enables Remarkable Enhancement in Remote Homology Detection Capability. <i>Journal of Molecular Biology</i> , 2014, 426, 962-979.	2.0	15
48	Rationalization and prediction of drug resistant mutations in targets for clinical anti-tubercular drugs. <i>Journal of Biomolecular Structure and Dynamics</i> , 2013, 31, 44-58.	2.0	3
49	A multi-level multi-scale approach to study essential genes in <i>Mycobacterium tuberculosis</i> . <i>BMC Systems Biology</i> , 2013, 7, 132.	3.0	17
50	Network approaches to drug discovery. <i>Expert Opinion on Drug Discovery</i> , 2013, 8, 7-20.	2.5	40
51	Current trends in modeling host-pathogen interactions. <i>Wiley Interdisciplinary Reviews: Data Mining and Knowledge Discovery</i> , 2013, 3, 109-128.	4.6	8
52	PocketAnnotate: towards site-based function annotation. <i>Nucleic Acids Research</i> , 2012, 40, W400-W408.	6.5	18
53	Cloning, expression, purification, and biochemical characterisation of the FIC motif containing protein of <i>Mycobacterium tuberculosis</i> . <i>Protein Expression and Purification</i> , 2012, 86, 58-67.	0.6	9
54	Crowd Sourcing a New Paradigm for Interactome Driven Drug Target Identification in <i>Mycobacterium tuberculosis</i> . <i>PLoS ONE</i> , 2012, 7, e39808.	1.1	36

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55	Rule-based modelling of iron homeostasis in tuberculosis. <i>Molecular BioSystems</i> , 2011, 7, 2750.	2.9	13
56	Computational approaches for drug target identification in pathogenic diseases. <i>Expert Opinion on Drug Discovery</i> , 2011, 6, 975-979.	2.5	9
57	PocketAlign A Novel Algorithm for Aligning Binding Sites in Protein Structures. <i>Journal of Chemical Information and Modeling</i> , 2011, 51, 1725-1736.	2.5	36
58	Systems biology of tuberculosis. <i>Tuberculosis</i> , 2011, 91, 487-496.	0.8	20
59	Structural biology of Mycobacterium tuberculosis proteins: The Indian efforts. <i>Tuberculosis</i> , 2011, 91, 456-468.	0.8	22
60	Open source drug discovery— A new paradigm of collaborative research in tuberculosis drug development. <i>Tuberculosis</i> , 2011, 91, 479-86.	0.8	42
61	Structural Annotation of Mycobacterium tuberculosis Proteome. <i>PLoS ONE</i> , 2011, 6, e27044.	1.1	33
62	Structural bioinformatics: Deriving biological insights from protein structures. <i>Interdisciplinary Sciences, Computational Life Sciences</i> , 2010, 2, 347-366.	2.2	9
63	Modeling metabolic adjustment in Mycobacterium tuberculosis upon treatment with isoniazid. <i>Systems and Synthetic Biology</i> , 2010, 4, 299-309.	1.0	11
64	Protein—protein interaction networks suggest different targets have different propensities for triggering drug resistance. <i>Systems and Synthetic Biology</i> , 2010, 4, 311-322.	1.0	21
65	The Multifunctional PE_PGRS11 Protein from Mycobacterium tuberculosis Plays a Role in Regulating Resistance to Oxidative Stress. <i>Journal of Biological Chemistry</i> , 2010, 285, 30389-30403.	1.6	59
66	Mycobacterium tuberculosis FtsZ requires at least one arginine residue at the C-terminal end for polymerization in vitro. <i>Acta Biochimica Et Biophysica Sinica</i> , 2010, 42, 58-69.	0.9	4
67	A systems perspective of host—pathogen interactions: predicting disease outcome in tuberculosis. <i>Molecular BioSystems</i> , 2010, 6, 516-530.	2.9	47
68	Localisation of Plasmodium falciparum uroporphyrinogen III decarboxylase of the heme-biosynthetic pathway in the apicoplast and characterisation of its catalytic properties. <i>International Journal for Parasitology</i> , 2009, 39, 559-568.	1.3	44
69	Structural basis for the function of anti-idiotypic antibody in immune memory. <i>Molecular Immunology</i> , 2009, 46, 1250-1255.	1.0	2
70	Computational systems approach for drug target discovery. <i>Expert Opinion on Drug Discovery</i> , 2009, 4, 1221-1236.	2.5	21
71	Flux balance analysis of biological systems: applications and challenges. <i>Briefings in Bioinformatics</i> , 2009, 10, 435-449.	3.2	354
72	Strategies for efficient disruption of metabolism in Mycobacterium tuberculosis from network analysis. <i>Molecular BioSystems</i> , 2009, 5, 1740.	2.9	35

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73	Functionally important movements in RecA molecules and filaments: studies involving mutation and environmental changes. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2008, 64, 1146-1157.	2.5	11
74	HLA-A*0201-restricted Cytotoxic T-cell Epitopes in Three PE/PPE Family Proteins of <i>Mycobacterium tuberculosis</i> . <i>Scandinavian Journal of Immunology</i> , 2008, 67, 411-417.	1.3	16
75	Scanning the Genome of <i>Mycobacterium tuberculosis</i> to Identify Potential Lectins. <i>Protein and Peptide Letters</i> , 2007, 14, 683-691.	0.4	9
76	Perpetuation of immunological memory through common MHC-I binding modes of peptidomimic and antigenic peptides. <i>Biochemical and Biophysical Research Communications</i> , 2007, 364, 308-312.	1.0	2
77	Snapshots of RecA Protein Involving Movement of the C-domain and Different Conformations of the DNA-binding Loops: Crystallographic and Comparative Analysis of 11 Structures of <i>Mycobacterium smegmatis</i> RecA. <i>Journal of Molecular Biology</i> , 2007, 367, 1130-1144.	2.0	40
78	Diagnosis and therapy of oral squamous cell carcinoma. <i>Expert Review of Anticancer Therapy</i> , 2007, 7, 317-329.	1.1	19
79	Conformationally locked thiosugars as potent α -mannosidase inhibitors: Synthesis, biochemical and docking studies. <i>Bioorganic and Medicinal Chemistry</i> , 2007, 15, 5659-5665.	1.4	17
80	Parallel implementation of AutoDock. <i>Journal of Applied Crystallography</i> , 2007, 40, 598-599.	1.9	31
81	Hallmarks of mycolic acid biosynthesis: A comparative genomics study. <i>Proteins: Structure, Function and Bioinformatics</i> , 2007, 69, 358-368.	1.5	11
82	Functional interaction of diphenols with polyphenol oxidase. <i>FEBS Journal</i> , 2007, 274, 4177-4187.	2.2	25
83	Metabolome Based Reaction Graphs of <i>M. tuberculosis</i> and <i>M. leprae</i> : A Comparative Network Analysis. <i>PLoS ONE</i> , 2007, 2, e881.	1.1	38
84	In-Silico Pharmacodynamics. <i>Applied Bioinformatics</i> , 2006, 5, 141-150.	1.7	4
85	A K52Q substitution in the globular domain of histone H1t modulates its nucleosome binding properties. <i>FEBS Letters</i> , 2006, 580, 5999-6006.	1.3	9
86	A combined immuno-informatics and structure-based modeling approach for prediction of T cell epitopes of secretory proteins of <i>Mycobacterium tuberculosis</i> . <i>Microbes and Infection</i> , 2006, 8, 738-746.	1.0	33
87	Common Scaffolds, Diverse Recognition Profiles. <i>Structure</i> , 2006, 14, 1093-1094.	1.6	3
88	Crystallographic identification of an ordered C-terminal domain and a second nucleotide-binding site in RecA: new insights into allostery. <i>Nucleic Acids Research</i> , 2006, 34, 2186-2195.	6.5	38
89	LectinDB: a plant lectin database. <i>Glycobiology</i> , 2006, 16, 938-946.	1.3	53
90	Molecular Mechanism of Dimerization of Bowman-Birk Inhibitors. <i>Journal of Biological Chemistry</i> , 2004, 279, 30425-30432.	1.6	46

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91	A database analysis of jacalin-like lectins: sequence-structure-function relationships. <i>Glycobiology</i> , 2004, 14, 1247-1263.	1.3	70
92	Î-Aminolevulinic Acid Dehydratase from <i>Plasmodium falciparum</i> . <i>Journal of Biological Chemistry</i> , 2004, 279, 6934-6942.	1.6	62
93	Extracting Hydrogen-Bond Signature Patterns from Protein Structure Data. <i>Applied Bioinformatics</i> , 2004, 3, 125-135.	1.7	2
94	Structural studies on MtRecA-nucleotide complexes: Insights into DNA and nucleotide binding and the structural signature of NTP recognition. <i>Proteins: Structure, Function and Bioinformatics</i> , 2003, 50, 474-485.	1.5	58
95	Determinants of histamine recognition: implications for the design of antihistamines. <i>Biochemical and Biophysical Research Communications</i> , 2003, 309, 425-431.	1.0	6
96	Conformation of gramicidin-A in CTAB micellar media. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2003, 70, 117-124.	1.7	5
97	Computational analysis of multivalency in lectins: structures of garlic lectin-oligosaccharide complexes and their aggregates. <i>Glycobiology</i> , 2003, 13, 765-775.	1.3	25
98	Molecular modeling of the chromatosome particle. <i>Nucleic Acids Research</i> , 2003, 31, 4264-4274.	6.5	68
99	Crystal Structures of <i>Mycobacterium smegmatis</i> RecA and Its Nucleotide Complexes. <i>Journal of Bacteriology</i> , 2003, 185, 4280-4284.	1.0	61
100	Re-refinement using reprocessed data to improve the quality of the structure: a case study involving garlic lectin. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2002, 58, 414-420.	2.5	18
101	Prediction of an HMG-box fold in the C-terminal domain of histone H1: Insights into its role in DNA condensation. <i>Proteins: Structure, Function and Bioinformatics</i> , 2002, 49, 71-81.	1.5	39
102	Crystal Structure of a Dimeric Mannose-specific Agglutinin from Garlic: Quaternary Association and Carbohydrate Specificity. <i>Journal of Molecular Biology</i> , 1999, 285, 1157-1168.	2.0	95
103	X-ray Studies on Crystalline Complexes Involving Amino Acids and Peptides. XXXIII. Crystal Structures of L- and DL-Arginine Complexed with Oxalic Acid and a Comparative Study of Amino Acidâ€“Oxalic Acid Complexes. <i>Acta Crystallographica Section B: Structural Science</i> , 1998, 54, 257-263.	1.8	19
104	Xray Studies on Crystalline Complexes Involving Amino Acids and Peptides XXXIV. Novel Mode of Aggregation, Interaction Patterns and Chiral Effects in the Maleic Acid Complexes of DL- and L-Arginine. <i>Journal of Biomolecular Structure and Dynamics</i> , 1998, 15, 1093-1100.	2.0	22
105	Crystallization and preliminary crystallographic studies on the mannose-specific lectin from garlic. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 1997, 53, 787-788.	2.5	6
106	Water-dependent domain motion and flexibility in ribonuclease A and the invariant features in its hydration shell. An X-ray study of two low-humidity crystal forms of the enzyme. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 1995, 51, 703-710.	2.5	28