

Nurit Haspel

List of Publications by Year in descending order

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Version: 2024-02-01

51
papers

958
citations

516215

16
h-index

454577

30
g-index

55
all docs

55
docs citations

55
times ranked

1156
citing authors

#	ARTICLE	IF	CITATIONS
1	Taking geometry to its edge: Fast unbound rigid (and hinge-bent) docking. <i>Proteins: Structure, Function and Bioinformatics</i> , 2003, 52, 107-112.	1.5	238
2	A Comparative Study of Amyloid Fibril Formation by Residues 15â€“19 of the Human Calcitonin Hormone: A Single Î²-Sheet Model with a Small Hydrophobic Core. <i>Journal of Molecular Biology</i> , 2005, 345, 1213-1227.	2.0	71
3	Tracing conformational changes in proteins. <i>BMC Structural Biology</i> , 2010, 10, S1.	2.3	61
4	Sequence dependence of C-end rule peptides in binding and activation of neuropilin-1 receptor. <i>Journal of Structural Biology</i> , 2013, 182, 78-86.	1.3	58
5	Principles of nanostructure design with protein building blocks. <i>Proteins: Structure, Function and Bioinformatics</i> , 2007, 68, 1-12.	1.5	51
6	Reducing the computational complexity of protein folding via fragment folding and assembly. <i>Protein Science</i> , 2003, 12, 1177-1187.	3.1	48
7	The Stability and Dynamics of the Human Calcitonin Amyloid Peptide DFNKF. <i>Biophysical Journal</i> , 2004, 87, 146-158.	0.2	46
8	De Novo Tubular Nanostructure Design Based on Self-Assembly of Î²-Helical Protein Motifs. <i>Structure</i> , 2006, 14, 1137-1148.	1.6	41
9	Electrostatic contributions drive the interaction between <i>Staphylococcus aureus</i> protein Efbâ€“ and its complement target C3d. <i>Protein Science</i> , 2008, 17, 1894-1906.	3.1	34
10	Predicting the Effect of Single and Multiple Mutations on Protein Structural Stability. <i>Molecules</i> , 2018, 23, 251.	1.7	31
11	Hierarchical protein folding pathways: A computational study of protein fragments. <i>Proteins: Structure, Function and Bioinformatics</i> , 2003, 51, 203-215.	1.5	29
12	Concepts and schemes for the re-engineering of physical protein modules: generating nanodevices via targeted replacements with constrained amino acids. <i>Physical Biology</i> , 2006, 3, S54-S62.	0.8	20
13	Active and Inactive Cdc42 Differ in Their Insert Region Conformational Dynamics. <i>Biophysical Journal</i> , 2021, 120, 306-318.	0.2	20
14	Side chain interactions determine the amyloid organization: a single layer Å-sheet molecular structure of the calcitonin peptide segment 15â€“19. <i>Physical Biology</i> , 2004, 1, 89-99.	0.8	19
15	AN EVOLUTIONARY CONSERVATION-BASED METHOD FOR REFINING AND RERANKING PROTEIN COMPLEX STRUCTURES. <i>Journal of Bioinformatics and Computational Biology</i> , 2012, 10, 1242002.	0.3	19
16	Changing the Charge Distribution of Î²-Helical-Based Nanostructures Can Provide the Conditions for Charge Transfer. <i>Biophysical Journal</i> , 2007, 93, 245-253.	0.2	18
17	Accurate refinement of docked protein complexes using evolutionary information and deep learning. <i>Journal of Bioinformatics and Computational Biology</i> , 2016, 14, 1642002.	0.3	16
18	Protein Segments with Conformationally Restricted Amino Acids Can Control Supramolecular Organization at the Nanoscale. <i>Journal of Chemical Information and Modeling</i> , 2009, 49, 1623-1629.	2.5	14

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19	GUIDING PROTEIN DOCKING WITH GEOMETRIC AND EVOLUTIONARY INFORMATION. Journal of Bioinformatics and Computational Biology, 2012, 10, 1242008.	0.3	11
20	Molecular characterization of l-phenylalanine terminated poly(l-lactide) conjugates. RSC Advances, 2014, 4, 23231.	1.7	11
21	Stability of Tubular Structures Based on α -Helical Proteins: Self-Assembled versus Polymerized Nanoconstructs and Wild-Type versus Mutated Sequences. Biomacromolecules, 2007, 8, 3135-3146.	2.6	10
22	Detecting intermediate protein conformations using algebraic topology. BMC Bioinformatics, 2017, 18, 502.	1.2	10
23	Patch-DCA: improved protein interface prediction by utilizing structural information and clustering DCA scores. Bioinformatics, 2020, 36, 1460-1467.	1.8	10
24	Multi-Resolution Rigidity-Based Sampling of Protein Conformational Paths. , 2013, , .		8
25	Conformational Exploration of Two Peptides and Their Hybrid Polymer Conjugates: Potentialities As Self-Aggregating Materials. Journal of Physical Chemistry B, 2012, 116, 13941-13952.	1.2	7
26	An Evolutionary Conservation & Rigidity Analysis Machine Learning Approach for Detecting Critical Protein Residues. , 2013, , .		6
27	Multi-scale characterization of the energy landscape of proteins with application to the C3D/Efb complex. Proteins: Structure, Function and Bioinformatics, 2010, 78, 1004-1014.	1.5	5
28	Accurate Prediction of Docked Protein Structure Similarity. Journal of Computational Biology, 2015, 22, 892-904.	0.8	5
29	Protein hormone fragmentation in intercellular signaling: hormones as nested information systems. Biology of Reproduction, 2021, 104, 887-901.	1.2	5
30	Preface: Selected Articles from 2015 Computational Structural Bioinformatics Workshop. Journal of Computational Biology, 2017, 24, 1-1.	0.8	4
31	Detecting chromosomal structural variation using jaccard distance and parallel architecture. , 2017, , .		4
32	TIDE: Inter-chromosomal translocation and insertion detection using embeddings. , 2014, , .		3
33	Methods for Detecting Critical Residues in Proteins. Methods in Molecular Biology, 2017, 1498, 227-242.	0.4	3
34	Integrating Co-Evolutionary Information in Monte Carlo Based Method for Proteins Trajectory Simulation. , 2019, , .		3
35	Efficient Exploration of Protein Conformational Pathways using RRT* and MC. , 2020, , .		3
36	Fusion Transcript Detection from RNA-Seq using Jaccard Distance. , 2020, , .		3

#	ARTICLE	IF	CITATIONS
37	Refining multimeric protein complexes using conservation, electrostatics and probabilistic selection. , 2012, , .		2
38	A Protocol for the Design of Protein and Peptide Nanostructure Self-Assemblies Exploiting Synthetic Amino Acids. Methods in Molecular Biology, 2017, 1529, 323-352.	0.4	2
39	Tracing conformational changes in proteins. , 2009, , .		1
40	Towards a hybrid method for detecting critical protein residues. , 2012, , .		1
41	Machine Learning Approaches for Predicting Protein Complex Similarity. Journal of Computational Biology, 2017, 24, 40-51.	0.8	1
42	Integrating Rigidity Analysis into the Exploration of Protein Conformational Pathways Using RRT* and MC. Molecules, 2021, 26, 2329.	1.7	1
43	Novel insights into target specificities and molecular mechanisms for two potent complement evasion proteins from Staphylococcus aureus. Molecular Immunology, 2008, 45, 4114-4115.	1.0	0
44	THE 5TH INTERNATIONAL CONFERENCE ON BIO-INSPIRED MODELS OF NETWORK, INFORMATION AND COMPUTING SYSTEMS (BIONETICS 2010) SPECIAL TRACK ON BIOINFORMATICS. Journal of Bioinformatics and Computational Biology, 2011, 09, v-vii.	0.3	0
45	Combining conservation and rigidity methods to detect critical residues in proteins. , 2012, , .		0
46	An evolutionary-guided iterative refinement approach for protein multimers. , 2012, , .		0
47	A new DP algorithm for comparing gene expression data using geometric similarity. , 2015, , .		0
48	The 7th Computational Structural Bioinformatics Workshop. Journal of Computational Biology, 2015, 22, 785-786.	0.8	0
49	Detecting intermediate protein conformations using algebraic topology. , 2016, , .		0
50	Introduction to the selected papers from the 7th International Conference on Bioinformatics and Computational Biology (BICoB 2015). Journal of Bioinformatics and Computational Biology, 2016, 14, 1602002.	0.3	0
51	Introduction to selected papers from the 8th International Conference on Bioinformatics and Computational Biology (BICOB 2016). Journal of Bioinformatics and Computational Biology, 2017, 15, 1702002.	0.3	0