

Ursula Pieper

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/3912953/publications.pdf>

Version: 2024-02-01

49
papers

8,305
citations

147801

31
h-index

214800

47
g-index

49
all docs

49
docs citations

49
times ranked

18194
citing authors

#	ARTICLE	IF	CITATIONS
1	Atlas of the Radical SAM Superfamily: Divergent Evolution of Function Using a "Plug and Play" Domain. <i>Methods in Enzymology</i> , 2018, 606, 1-71.	1.0	99
2	Prediction of Functionally Important Phospho-Regulatory Events in <i>Xenopus laevis</i> Oocytes. <i>PLoS Computational Biology</i> , 2015, 11, e1004362.	3.2	14
3	ModBase, a database of annotated comparative protein structure models and associated resources. <i>Nucleic Acids Research</i> , 2014, 42, D336-D346.	14.5	275
4	Coordinating the impact of structural genomics on the human α -helical transmembrane proteome. <i>Nature Structural and Molecular Biology</i> , 2013, 20, 135-138.	8.2	64
5	A Role for Matrix Metalloproteinases in Regulating Mammary Stem Cell Function via the Wnt Signaling Pathway. <i>Cell Stem Cell</i> , 2013, 13, 300-313.	11.1	123
6	Structure, Dynamics, Evolution, and Function of a Major Scaffold Component in the Nuclear Pore Complex. <i>Structure</i> , 2013, 21, 560-571.	3.3	53
7	Target Prediction for an Open Access Set of Compounds Active against <i>Mycobacterium tuberculosis</i> . <i>PLoS Computational Biology</i> , 2013, 9, e1003253.	3.2	51
8	Consequences of domain insertion on sequence-structure divergence in a superfold. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013, 110, E3381-7.	7.1	25
9	Biochemical characterization and structural modeling of human cathepsin E variant 2 in comparison to the wild-type protein. <i>Biological Chemistry</i> , 2012, 393, 177-186.	2.5	3
10	SALIGN: a web server for alignment of multiple protein sequences and structures. <i>Bioinformatics</i> , 2012, 28, 2072-2073.	4.1	72
11	Atomic structure of the nuclear pore complex targeting domain of a Nup116 homologue from the yeast, <i>Candida glabrata</i> . <i>Proteins: Structure, Function and Bioinformatics</i> , 2012, 80, 2110-2116.	2.6	7
12	Facile backbone structure determination of human membrane proteins by NMR spectroscopy. <i>Nature Methods</i> , 2012, 9, 834-839.	19.0	83
13	Structure of the C-terminal domain of <i>Saccharomyces cerevisiae</i> Nup133, a component of the nuclear pore complex. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 1672-1677.	2.6	16
14	Response to "Predictable difficulty or difficulty to predict". <i>Protein Science</i> , 2011, 20, 4-5.	7.6	0
15	A Conserved Coatmer-related Complex Containing Sec13 and Seh1 Dynamically Associates With the Vacuole in <i>Saccharomyces cerevisiae</i> . <i>Molecular and Cellular Proteomics</i> , 2011, 10, M110.006478.	3.8	115
16	ModBase, a database of annotated comparative protein structure models, and associated resources. <i>Nucleic Acids Research</i> , 2011, 39, D465-D474.	14.5	506
17	Comparison of human solute carriers. <i>Protein Science</i> , 2010, 19, 412-428.	7.6	99
18	Functional hot spots in human ATP-binding cassette transporter nucleotide binding domains. <i>Protein Science</i> , 2010, 19, 2110-2121.	7.6	19

#	ARTICLE	IF	CITATIONS
19	Structures of the autoproteolytic domain from the <i>Saccharomyces cerevisiae</i> nuclear pore complex component, Nup145. <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 1992-1998.	2.6	13
20	Structure of a putative BenF-like porin from <i>Pseudomonas fluorescens</i> Pf5 at 2.6 Å... resolution. <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 3056-3062.	2.6	17
21	Prediction of protease substrates using sequence and structure features. <i>Bioinformatics</i> , 2010, 26, 1714-1722.	4.1	61
22	MODBASE, a database of annotated comparative protein structure models and associated resources. <i>Nucleic Acids Research</i> , 2009, 37, D347-D354.	14.5	154
23	A Kernel for Open Source Drug Discovery in Tropical Diseases. <i>PLoS Neglected Tropical Diseases</i> , 2009, 3, e418.	3.0	23
24	Carbanion or Amide? First Charge Density Study of Parent 2-picolyllithium. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 2978-2982.	13.8	51
25	Target selection and annotation for the structural genomics of the amidohydrolase and enolase superfamilies. <i>Journal of Structural and Functional Genomics</i> , 2009, 10, 107-125.	1.2	25
26	A survey of integral α -helical membrane proteins. <i>Journal of Structural and Functional Genomics</i> , 2009, 10, 269-280.	1.2	12
27	A kernel for the Tropical Disease Initiative. <i>Nature Biotechnology</i> , 2009, 27, 320-321.	17.5	7
28	Selecting Optimum Eukaryotic Integral Membrane Proteins for Structure Determination by Rapid Expression and Solubilization Screening. <i>Journal of Molecular Biology</i> , 2009, 385, 820-830.	4.2	53
29	Genomic-scale prioritization of drug targets: the TDR Targets database. <i>Nature Reviews Drug Discovery</i> , 2008, 7, 900-907.	46.4	282
30	DBAli tools: mining the protein structure space. <i>Nucleic Acids Research</i> , 2007, 35, W393-W397.	14.5	25
31	The AnnoLite and AnnoLyze programs for comparative annotation of protein structures. <i>BMC Bioinformatics</i> , 2007, 8, S4.	2.6	36
32	Comparative Protein Structure Modeling Using MODELLER. <i>Current Protocols in Protein Science</i> , 2007, 50, Unit 2.9.	2.8	1,056
33	Comparative Protein Structure Modeling Using Modeller. <i>Current Protocols in Bioinformatics</i> , 2006, 15, Unit-5.6.	25.8	2,858
34	Protein complex compositions predicted by structural similarity. <i>Nucleic Acids Research</i> , 2006, 34, 2943-2952.	14.5	56
35	MODBASE: a database of annotated comparative protein structure models and associated resources. <i>Nucleic Acids Research</i> , 2006, 34, D291-D295.	14.5	265
36	Comparative Protein Structure Modeling. , 2005, , 831-860.		15

#	ARTICLE	IF	CITATIONS
37	LS-SNP: large-scale annotation of coding non-synonymous SNPs based on multiple information sources. <i>Bioinformatics</i> , 2005, 21, 2814-2820.	4.1	202
38	High-Throughput Computational and Experimental Techniques in Structural Genomics. <i>Genome Research</i> , 2004, 14, 2145-2154.	5.5	59
39	MODBASE, a database of annotated comparative protein structure models, and associated resources. <i>Nucleic Acids Research</i> , 2004, 32, 217D-222.	14.5	256
40	Modeling Protein Structure from its Sequence. <i>Current Protocols in Bioinformatics</i> , 2003, 3, 5.1.1.	25.8	6
41	Tools for comparative protein structure modeling and analysis. <i>Nucleic Acids Research</i> , 2003, 31, 3375-3380.	14.5	406
42	ModView, visualization of multiple protein sequences and structures. <i>Bioinformatics</i> , 2003, 19, 165-166.	4.1	18
43	MODBASE, a database of annotated comparative protein structure models. <i>Nucleic Acids Research</i> , 2002, 30, 255-259.	14.5	114
44	Structural genomics: A pipeline for providing structures for the biologist. <i>Protein Science</i> , 2002, 11, 723-738.	7.6	168
45	Homology-based annotation yields 1,042 new candidate genes in the <i>Drosophila melanogaster</i> genome. <i>Nature Genetics</i> , 2001, 27, 337-340.	21.4	58
46	Protein structure modeling for structural genomics. <i>Nature Structural Biology</i> , 2000, 7, 986-990.	9.7	199
47	Structural features of halophilicity derived from the crystal structure of dihydrofolate reductase from the Dead Sea halophilic archaeon, <i>Haloferax volcanii</i> . <i>Structure</i> , 1998, 6, 75-88.	3.3	96
48	Structural evidence for the evolutionary divergence of mycoplasma from Gram-positive bacteria: the histidine-containing phosphocarrier protein. <i>Structure</i> , 1995, 3, 781-790.	3.3	22
49	Syntheses and x-ray structures of (diphenylpyridylmethyl)lithium, -sodium, and -potassium in comparison with the triphenylmethyl derivatives. <i>Organometallics</i> , 1993, 12, 1201-1206.	2.3	68