

Claudia Milln

List of Publications by Citations

Source: <https://exaly.com/author-pdf/6837428/claudia-millan-publications-by-citations.pdf>

Version: 2024-04-27

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

26

papers

962

citations

13

h-index

28

g-index

28

ext. papers

2,286

ext. citations

7.8

avg, IF

4.21

L-index

#	Paper	IF	Citations
26	Accurate prediction of protein structures and interactions using a three-track neural network. <i>Science</i> , 2021 , 373, 871-876	33.3	522
25	Structure and mechanism of a bacterial host-protein citrullinating virulence factor, <i>Porphyromonas gingivalis</i> peptidylarginine deiminase. <i>Scientific Reports</i> , 2015 , 5, 11969	4.9	57
24	Exploiting tertiary structure through local folds for crystallographic phasing. <i>Nature Methods</i> , 2013 , 10, 1099-101	21.6	51
23	Structural basis of meiotic chromosome synapsis through SYCP1 self-assembly. <i>Nature Structural and Molecular Biology</i> , 2018 , 25, 557-569	17.6	38
22	ARCIMBOLDO_LITE: single-workstation implementation and use. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2015 , 71, 1921-30		37
21	Macromolecular ab initio phasing enforcing secondary and tertiary structure. <i>IUCrJ</i> , 2015 , 2, 95-105	4.7	35
20	Structure solution with ARCIMBOLDO using fragments derived from distant homology models. <i>FEBS Journal</i> , 2014 , 281, 4029-45	5.7	33
19	On the application of the expected log-likelihood gain to decision making in molecular replacement. <i>Acta Crystallographica Section D: Structural Biology</i> , 2018 , 74, 245-255	5.5	22
18	ARCIMBOLDO on coiled coils. <i>Acta Crystallographica Section D: Structural Biology</i> , 2018 , 74, 194-204	5.5	20
17	Exploiting distant homologues for phasing through the generation of compact fragments, local fold refinement and partial solution combination. <i>Acta Crystallographica Section D: Structural Biology</i> , 2018 , 74, 290-304	5.5	19
16	Exolytic and endolytic turnover of peptidoglycan by lytic transglycosylase Slt of. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018 , 115, 4393-4398	11.5	18
15	Ega3 from the fungal pathogen is an endo- β -1,4-galactosaminidase that disrupts microbial biofilms. <i>Journal of Biological Chemistry</i> , 2019 , 294, 13833-13849	5.4	18
14	Assessing the utility of CASP14 models for molecular replacement. <i>Proteins: Structure, Function and Bioinformatics</i> , 2021 , 89, 1752-1769	4.2	18
13	Gyre and gimble: a maximum-likelihood replacement for Patterson correlation refinement. <i>Acta Crystallographica Section D: Structural Biology</i> , 2018 , 74, 279-289	5.5	13
12	Combining phase information in reciprocal space for molecular replacement with partial models. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2015 , 71, 1931-45		10
11	A highly conserved region essential for NMD in the Upf2 N-terminal domain. <i>Journal of Molecular Biology</i> , 2014 , 426, 3689-3702	6.5	10
10	ALEPH: a network-oriented approach for the generation of fragment-based libraries and for structure interpretation. <i>Acta Crystallographica Section D: Structural Biology</i> , 2020 , 76, 193-208	5.5	9

9	Accurate prediction of protein structures and interactions using a 3-track network		9
8	ALIXE: a phase-combination tool for fragment-based molecular replacement. <i>Acta Crystallographica Section D: Structural Biology</i> , 2020 , 76, 209-220	5.5	7
7	Fragment-based determination of a proteinase K structure from MicroED data using ARCIMBOLDO_SHREDDER. <i>Acta Crystallographica Section D: Structural Biology</i> , 2020 , 76, 703-712	5.5	6
6	SEQUENCE SLIDER: expanding polyalanine fragments for phasing with multiple side-chain hypotheses. <i>Acta Crystallographica Section D: Structural Biology</i> , 2020 , 76, 221-237	5.5	5
5	Structure of a 13-fold superhelix (almost) determined from first principles. <i>IUCrJ</i> , 2015 , 2, 177-87	4.7	2
4	Improving AlphaFold modeling using implicit information from experimental density maps		2
3	Assessing the utility of CASP14 models for molecular replacement		1
2	Crystallographic structure solution. <i>Arbor</i> , 2015 , 191, a218	0.2	
1	Structure of the class XI myosin globular tail reveals evolutionary hallmarks for cargo recognition in plants. <i>Acta Crystallographica Section D: Structural Biology</i> , 2021 , 77, 522-533	5.5	