

Dzmitry Padhorny

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6147681/publications.pdf>

Version: 2024-02-01

16
papers

2,709
citations

1040056

9
h-index

996975

15
g-index

17
all docs

17
docs citations

17
times ranked

4942
citing authors

#	ARTICLE	IF	CITATIONS
1	Elucidation of protein function using computational docking and hotspot analysis by <i>ClusPro</i> and <i>FTMap</i> . <i>Acta Crystallographica Section D: Structural Biology</i> , 2022, 78, 690-697.	2.3	3
2	Assessing the binding properties of <i>CASP14</i> targets and models. <i>Proteins: Structure, Function and Bioinformatics</i> , 2021, 89, 1922-1939.	2.6	11
3	Prediction of protein assemblies, the next frontier: The <i>CASP14</i> CAPRI experiment. <i>Proteins: Structure, Function and Bioinformatics</i> , 2021, 89, 1800-1823.	2.6	73
4	Side-chain Packing Using SE(3)-Transformer. , 2021, , .		2
5	Sampling and refinement protocols for template-based macrocycle docking: 2018 D3R Grand Challenge 4. <i>Journal of Computer-Aided Molecular Design</i> , 2020, 34, 179-189.	2.9	8
6	Modeling beta-sheet peptide-protein interactions: Rosetta FlexPepDock in CAPRI rounds 38-45. <i>Proteins: Structure, Function and Bioinformatics</i> , 2020, 88, 1037-1049.	2.6	10
7	Actionable Cytopathogenic Host Responses of Human Alveolar Type 2 Cells to SARS-CoV-2. <i>Molecular Cell</i> , 2020, 80, 1104-1122.e9.	9.7	94
8	ClusPro in rounds 38 to 45 of CAPRI: Toward combining template-based methods with free docking. <i>Proteins: Structure, Function and Bioinformatics</i> , 2020, 88, 1082-1090.	2.6	5
9	Template-based modeling by ClusPro in CASP13 and the potential for using co-evolutionary information in docking. <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 1241-1248.	2.6	15
10	Blind prediction of homo- and hetero-protein complexes: The CASP13 CAPRI experiment. <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 1200-1221.	2.6	99
11	Allostery in Its Many Disguises: From Theory to Applications. <i>Structure</i> , 2019, 27, 566-578.	3.3	285
12	Monte Carlo on the manifold and MD refinement for binding pose prediction of protein-ligand complexes: 2017 D3R Grand Challenge. <i>Journal of Computer-Aided Molecular Design</i> , 2019, 33, 119-127.	2.9	12
13	Protein-ligand docking using FFT based sampling: D3R case study. <i>Journal of Computer-Aided Molecular Design</i> , 2018, 32, 225-230.	2.9	9
14	The ClusPro web server for protein-protein docking. <i>Nature Protocols</i> , 2017, 12, 255-278.	12.0	1,959
15	ClusPro-DC: Dimer Classification by the Cluspro Server for Protein-Protein Docking. <i>Journal of Molecular Biology</i> , 2017, 429, 372-381.	4.2	36
16	Protein-protein docking by fast generalized Fourier transforms on 5D rotational manifolds. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, E4286-93.	7.1	43