

Kathryn A Porter

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/11772834/publications.pdf>

Version: 2024-02-01

12
papers

2,780
citations

1039406

9
h-index

1199166

12
g-index

13
all docs

13
docs citations

13
times ranked

4504
citing authors

#	ARTICLE	IF	CITATIONS
1	The ClusPro web server for protein-protein docking. <i>Nature Protocols</i> , 2017, 12, 255-278.	5.5	1,959
2	Performance and Its Limits in Rigid Body Protein-Protein Docking. <i>Structure</i> , 2020, 28, 1071-1081.e3.	1.6	306
3	High-resolution global peptide-protein docking using fragments-based PIPER-FlexPepDock. <i>PLoS Computational Biology</i> , 2017, 13, e1005905.	1.5	119
4	ClusPro PeptiDock: efficient global docking of peptide recognition motifs using FFT. <i>Bioinformatics</i> , 2017, 33, 3299-3301.	1.8	102
5	What method to use for protein-protein docking?. <i>Current Opinion in Structural Biology</i> , 2019, 55, 1-7.	2.6	83
6	Prediction of protein assemblies, the next frontier: The CASP14-CAPRI experiment. <i>Proteins: Structure, Function and Bioinformatics</i> , 2021, 89, 1800-1823.	1.5	73
7	Kinase Atlas: Druggability Analysis of Potential Allosteric Sites in Kinases. <i>Journal of Medicinal Chemistry</i> , 2019, 62, 6512-6524.	2.9	52
8	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.	3.3	43
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.	1.5	15
10	Progress toward improved understanding of antibody maturation. <i>Current Opinion in Structural Biology</i> , 2021, 67, 226-231.	2.6	12
11	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.	1.5	5
12	Conservation of binding properties in protein models. <i>Computational and Structural Biotechnology Journal</i> , 2021, 19, 2549-2566.	1.9	2