

# Raphaël A G Chaleil

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/8502160/publications.pdf>

Version: 2024-02-01

17  
papers

1,635  
citations

623734

14  
h-index

888059

17  
g-index

17  
all docs

17  
docs citations

17  
times ranked

2479  
citing authors

#	ARTICLE	IF	CITATIONS
1	Updates to the Integrated Protein-Protein Interaction Benchmarks: Docking Benchmark Version 5 and Affinity Benchmark Version 2. <i>Journal of Molecular Biology</i> , 2015, 427, 3031-3041.	4.2	348
2	SwarmDock: a server for flexible protein-protein docking. <i>Bioinformatics</i> , 2013, 29, 807-809.	4.1	259
3	Butyrophilin-2A1 Directly Binds Germline-Encoded Regions of the V $\beta$ 9V $\beta$ 2 TCR and Is Essential for Phosphoantigen Sensing. <i>Immunity</i> , 2020, 52, 487-498.e6.	14.3	164
4	The $\beta$ TCR combines innate immunity with adaptive immunity by utilizing spatially distinct regions for agonist selection and antigen responsiveness. <i>Nature Immunology</i> , 2018, 19, 1352-1365.	14.5	163
5	Prediction of homoprotein and heteroprotein complexes by protein docking and template-based modeling: A CASP-CAPRI experiment. <i>Proteins: Structure, Function and Bioinformatics</i> , 2016, 84, 323-348.	2.6	148
6	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
7	A simple biophysical model emulates budding yeast chromosome condensation. <i>ELife</i> , 2015, 4, e05565.	6.0	87
8	Evolution of Enzymes in Metabolism: A Network Perspective. <i>Journal of Molecular Biology</i> , 2002, 320, 751-770.	4.2	73
9	Prediction of protein assemblies, the next frontier: The CASP14-CAPRI experiment. <i>Proteins: Structure, Function and Bioinformatics</i> , 2021, 89, 1800-1823.	2.6	73
10	Application of abductive ILP to learning metabolic network inhibition from temporal data. <i>Machine Learning</i> , 2006, 64, 209-230.	5.4	71
11	Implicit flexibility in protein docking: Cross-docking and local refinement. <i>Proteins: Structure, Function and Bioinformatics</i> , 2007, 69, 750-757.	2.6	53
12	A Markov-chain model description of binding funnels to enhance the ranking of docked solutions. <i>Proteins: Structure, Function and Bioinformatics</i> , 2013, 81, 2143-2149.	2.6	32
13	Flexible Protein-Protein Docking with SwarmDock. <i>Methods in Molecular Biology</i> , 2018, 1764, 413-428.	0.9	20
14	A machine learning approach for ranking clusters of docked protein-protein complexes by pairwise cluster comparison. <i>Proteins: Structure, Function and Bioinformatics</i> , 2017, 85, 528-543.	2.6	18
15	Enhanced sampling of protein conformational states for dynamic cross-docking within the protein-protein docking server SwarmDock. <i>Proteins: Structure, Function and Bioinformatics</i> , 2020, 88, 962-972.	2.6	16
16	Modeling the effects of toxins in metabolic networks. <i>IEEE Engineering in Medicine and Biology Magazine</i> , 2007, 26, 37-46.	0.8	10
17	A Guide for Protein-Protein Docking Using SwarmDock. <i>Methods in Molecular Biology</i> , 2020, 2165, 199-216.	0.9	1