

# Oliver Carrillo

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/3287310/publications.pdf>

Version: 2024-02-01

10  
papers

387  
citations

1040056

9  
h-index

1474206

9  
g-index

11  
all docs

11  
docs citations

11  
times ranked

642  
citing authors

#	ARTICLE	IF	CITATIONS
1	MoDEL (Molecular Dynamics Extended Library): A Database of Atomistic Molecular Dynamics Trajectories. <i>Structure</i> , 2010, 18, 1399-1409.	3.3	123
2	FlexServ: an integrated tool for the analysis of protein flexibility. <i>Bioinformatics</i> , 2009, 25, 1709-1710.	4.1	72
3	Prediction and validation of protein intermediate states from structurally rich ensembles and coarse-grained simulations. <i>Nature Communications</i> , 2016, 7, 12575.	12.8	62
4	Exploring the Suitability of Coarse-Grained Techniques for the Representation of Protein Dynamics. <i>Biophysical Journal</i> , 2008, 95, 2127-2138.	0.5	42
5	Coarse-grained Representation of Protein Flexibility. Foundations, Successes, and Shortcomings. <i>Advances in Protein Chemistry and Structural Biology</i> , 2011, 85, 183-215.	2.3	33
6	Structural Plasticity and Functional Implications of Internal Cavities in Distal Mutants of Type 1 Non-Symbiotic Hemoglobin AHb1 from <i>Arabidopsis thaliana</i> . <i>Journal of Physical Chemistry B</i> , 2009, 113, 16028-16038.	2.6	20
7	GRID-MD: A tool for massive simulation of protein channels. <i>Proteins: Structure, Function and Bioinformatics</i> , 2008, 70, 892-899.	2.6	13
8	Application of Drug-Perturbed Essential Dynamics/Molecular Dynamics (ED/MD) to Virtual Screening and Rational Drug Design. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 2204-2214.	5.3	11
9	Fast Atomistic Molecular Dynamics Simulations from Essential Dynamics Samplings. <i>Journal of Chemical Theory and Computation</i> , 2012, 8, 792-799.	5.3	10
10	Multiplicative white-noise-induced phase transitions: the role of the stochastic interpretation. , 2004, , ,		0