## **Ognjen** Perisic

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Single-molecule force spectroscopy reveals a mechanically stable protein fold and the rational tuning of its mechanical stability. Proceedings of the National Academy of Sciences of the United States of America, 2007, 104, 9278-9283.	7.1	117
2	Ligand efficiency indices for an effective mapping of chemico-biological space: the concept of an atlas-like representation. Drug Discovery Today, 2010, 15, 804-811.	6.4	102
3	Modeling Studies of Chromatin Fiber Structure as a Function of DNA Linker Length. Journal of Molecular Biology, 2010, 403, 777-802.	4.2	98
4	Mesoscale simulations of two nucleosome-repeat length oligonucleosomes. Physical Chemistry Chemical Physics, 2009, 11, 10729.	2.8	45
5	Sensitive effect of linker histone binding mode and subtype on chromatin condensation. Nucleic Acids Research, 2019, 47, 4948-4957.	14.5	43
6	Gene Mutation Profiles in Primary Diffuse Large B Cell Lymphoma of Central Nervous System: Next Generation Sequencing Analyses. International Journal of Molecular Sciences, 2016, 17, 683.	4.1	29
7	Mechanical Signaling on the Single Protein Level Studied Using Steered Molecular Dynamics. Cell Biochemistry and Biophysics, 2009, 55, 141-152.	1.8	27
8	Dependence of the Linker Histone and Chromatin Condensation on the Nucleosome Environment. Journal of Physical Chemistry B, 2017, 121, 7823-7832.	2.6	25
9	Computational strategies to address chromatin structure problems. Physical Biology, 2016, 13, 035006.	1.8	14
10	On the Improvement of Free-Energy Calculation from Steered Molecular Dynamics Simulations Using Adaptive Stochastic Perturbation Protocols. PLoS ONE, 2014, 9, e101810.	2.5	13
11	Parallel targeted next generation sequencing of childhood and adult acute myeloid leukemia patients reveals uniform genomic profile of the disease. Tumor Biology, 2016, 37, 13391-13401.	1.8	13
12	Recognition of Potential COVID-19 Drug Treatments through the Study of Existing Protein–Drug and Protein–Protein Structures: An Analysis of Kinetically Active Residues. Biomolecules, 2020, 10, 1346.	4.0	13
13	Improved protein fold assignment using support vector machines. International Journal of Bioinformatics Research and Applications, 2005, 1, 319.	0.2	8
14	Heterodimer Binding Scaffolds Recognition via the Analysis of Kinetically Hot Residues. Pharmaceuticals, 2018, 11, 29.	3.8	5
15	Efficient free-energy-profile reconstruction using adaptive stochastic perturbation protocols. Physical Review E, 2011, 84, 056705.	2.1	4
16	Pulling-spring modulation as a method for improving the potential-of-mean-force reconstruction in single-molecule manipulation experiments. Physical Review E, 2013, 87, 013303.	2.1	4
17	Mechanism for the Unfolding of the TOP7 Protein in Steered Molecular Dynamics Simulations as Revealed by Mutual Information Analysis. Frontiers in Molecular Biosciences, 2021, 8, 696609.	3.5	1