

# Harianto Tjong

## List of Publications by Year in descending order

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31  
papers

2,351  
citations

331670

21  
h-index

454955

30  
g-index

34  
all docs

34  
docs citations

34  
times ranked

3520  
citing authors

#	ARTICLE	IF	CITATIONS
1	Oncogenic extrachromosomal DNA functions as mobile enhancers to globally amplify chromosomal transcription. <i>Cancer Cell</i> , 2021, 39, 694-707.e7.	16.8	115
2	ChIA-PIPE: A fully automated pipeline for comprehensive ChIA-PET data analysis and visualization. <i>Science Advances</i> , 2020, 6, eaay2078.	10.3	22
3	Chromatin interaction analyses elucidate the roles of PRC2-bound silencers in mouse development. <i>Nature Genetics</i> , 2020, 52, 264-272.	21.4	104
4	Mapping the Global Chromatin Connectivity Network for Sox2 Function in Neural Stem Cell Maintenance. <i>Cell Stem Cell</i> , 2019, 24, 462-476.e6.	11.1	72
5	Producing genome structure populations with the dynamic and automated PGS software. <i>Nature Protocols</i> , 2018, 13, 915-926.	12.0	67
6	Picky comprehensively detects high-resolution structural variants in nanopore long reads. <i>Nature Methods</i> , 2018, 15, 455-460.	19.0	80
7	Quantitative Methods to Investigate the 4D Dynamics of Heterochromatic Repair Sites in Drosophila Cells. <i>Methods in Enzymology</i> , 2018, 601, 359-389.	1.0	24
8	The three-dimensional genome organization of <i>Drosophila melanogaster</i> through data integration. <i>Genome Biology</i> , 2017, 18, 145.	8.8	86
9	Mining 3D genome structure populations identifies major factors governing the stability of regulatory communities. <i>Nature Communications</i> , 2016, 7, 11549.	12.8	36
10	Electrostatic effects on the folding stability of FKBP12. <i>Protein Engineering, Design and Selection</i> , 2016, 29, 301-308.	2.1	6
11	Global reorganization of budding yeast chromosome conformation in different physiological conditions. <i>Journal of Cell Biology</i> , 2016, 212, 321-334.	5.2	36
12	Population-based 3D genome structure analysis reveals driving forces in spatial genome organization. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, E1663-72.	7.1	182
13	TopDom: an efficient and deterministic method for identifying topological domains in genomes. <i>Nucleic Acids Research</i> , 2016, 44, e70-e70.	14.5	246
14	Comparative 3D Genome Structure Analysis of the Fission and the Budding Yeast. <i>PLoS ONE</i> , 2015, 10, e0119672.	2.5	20
15	Physical tethering and volume exclusion determine higher-order genome organization in budding yeast. <i>Genome Research</i> , 2012, 22, 1295-1305.	5.5	190
16	Genome architectures revealed by tethered chromosome conformation capture and population-based modeling. <i>Nature Biotechnology</i> , 2012, 30, 90-98.	17.5	524
17	Exploring the spatial and temporal organization of a cell's proteome. <i>Journal of Structural Biology</i> , 2011, 173, 483-496.	2.8	36
18	The Folding Transition-State Ensemble of a Four-Helix Bundle Protein: Helix Propensity as a Determinant and Macromolecular Crowding as a Probe. <i>Biophysical Journal</i> , 2010, 98, 2273-2280.	0.5	22

#	ARTICLE	IF	CITATIONS
19	Inverse tuning of metal binding affinity and protein stability by altering charged coordination residues in designed calcium binding proteins. <i>PMC Biophysics</i> , 2009, 2, 11.	2.3	12
20	Rational design of a conformationally switchable Ca <sup>2+</sup> and Tb <sup>3+</sup> binding protein without the use of multiple coupled metal binding sites. <i>FEBS Journal</i> , 2008, 275, 5048-5061.	4.7	12
21	Prediction of Protein Solubility from Calculation of Transfer Free Energy. <i>Biophysical Journal</i> , 2008, 95, 2601-2609.	0.5	50
22	On the Dielectric Boundary in Poisson-Boltzmann Calculations. <i>Journal of Chemical Theory and Computation</i> , 2008, 4, 507-514.	5.3	30
23	Accurate Calculations of Binding, Folding, and Transfer Free Energies by a Scaled Generalized Born Method. <i>Journal of Chemical Theory and Computation</i> , 2008, 4, 1733-1744.	5.3	10
24	Modeling Protein-Protein and Protein-Nucleic Acid Interactions: Structure, Thermodynamics, and Kinetics. <i>Annual Reports in Computational Chemistry</i> , 2008, , 67-87.	1.7	3
25	Spontaneous conformational change and toxin binding in $\alpha 7$ acetylcholine receptor: Insight into channel activation and inhibition. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 8280-8285.	7.1	46
26	GBr6NL: A generalized Born method for accurately reproducing solvation energy of the nonlinear Poisson-Boltzmann equation. <i>Journal of Chemical Physics</i> , 2007, 126, 195102.	3.0	41
27	DISPLAR: an accurate method for predicting DNA-binding sites on protein surfaces. <i>Nucleic Acids Research</i> , 2007, 35, 1465-1477.	14.5	142
28	PI2PE: protein interface/interior prediction engine. <i>Nucleic Acids Research</i> , 2007, 35, W357-W362.	14.5	32
29	GBr6: A Parameterization-Free, Accurate, Analytical Generalized Born Method. <i>Journal of Physical Chemistry B</i> , 2007, 111, 3055-3061.	2.6	84
30	The dependence of electrostatic solvation energy on dielectric constants in Poisson-Boltzmann calculations. <i>Journal of Chemical Physics</i> , 2006, 125, 206101.	3.0	16
31	Modeling Protein Solubility in Implicit Solvent. , 0, , 191-207.		0