

Po-Ssu Huang

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6199043/publications.pdf>

Version: 2024-02-01

46
papers

6,900
citations

201674

27
h-index

223800

46
g-index

54
all docs

54
docs citations

54
times ranked

8565
citing authors

#	ARTICLE	IF	CITATIONS
1	The coming of age of de novo protein design. <i>Nature</i> , 2016, 537, 320-327.	27.8	1,069
2	Rational HIV Immunogen Design to Target Specific Germline B Cell Receptors. <i>Science</i> , 2013, 340, 711-716.	12.6	680
3	A Potent and Broad Neutralizing Antibody Recognizes and Penetrates the HIV Glycan Shield. <i>Science</i> , 2011, 334, 1097-1103.	12.6	644
4	Macromolecular modeling and design in Rosetta: recent methods and frameworks. <i>Nature Methods</i> , 2020, 17, 665-680.	19.0	513
5	Protein structure determination using metagenome sequence data. <i>Science</i> , 2017, 355, 294-298.	12.6	456
6	Design of a hyperstable 60-subunit protein icosahedron. <i>Nature</i> , 2016, 535, 136-139.	27.8	373
7	Accurate de novo design of hyperstable constrained peptides. <i>Nature</i> , 2016, 538, 329-335.	27.8	327
8	RosettaRemodel: A Generalized Framework for Flexible Backbone Protein Design. <i>PLoS ONE</i> , 2011, 6, e24109.	2.5	310
9	De novo design of a fluorescence-activating β -barrel. <i>Nature</i> , 2018, 561, 485-491.	27.8	269
10	High thermodynamic stability of parametrically designed helical bundles. <i>Science</i> , 2014, 346, 481-485.	12.6	264
11	Exploring the repeat protein universe through computational protein design. <i>Nature</i> , 2015, 528, 580-584.	27.8	227
12	De novo design of a four-fold symmetric TIM-barrel protein with atomic-level accuracy. <i>Nature Chemical Biology</i> , 2016, 12, 29-34.	8.0	214
13	Computation-Guided Backbone Grafting of a Discontinuous Motif onto a Protein Scaffold. <i>Science</i> , 2011, 334, 373-376.	12.6	212
14	Computational design of transmembrane pores. <i>Nature</i> , 2020, 585, 129-134.	27.8	120
15	Protein sequence design with a learned potential. <i>Nature Communications</i> , 2022, 13, 746.	12.8	85
16	Control of repeat-protein curvature by computational protein design. <i>Nature Structural and Molecular Biology</i> , 2015, 22, 167-174.	8.2	84
17	A de novo designed protein-protein interface. <i>Protein Science</i> , 2007, 16, 2770-2774.	7.6	83
18	A General Computational Approach for Repeat Protein Design. <i>Journal of Molecular Biology</i> , 2015, 427, 563-575.	4.2	72

#	ARTICLE	IF	CITATIONS
19	Theoretical basis for stabilizing messenger RNA through secondary structure design. <i>Nucleic Acids Research</i> , 2021, 49, 10604-10617.	14.5	67
20	Structure and Functional Binding Epitope of V-domain Ig Suppressor of T Cell Activation. <i>Cell Reports</i> , 2019, 28, 2509-2516.e5.	6.4	61
21	Structure-based protein design with deep learning. <i>Current Opinion in Chemical Biology</i> , 2021, 65, 136-144.	6.1	53
22	Ig-VAE: Generative modeling of protein structure by direct 3D coordinate generation. <i>PLoS Computational Biology</i> , 2022, 18, e1010271.	3.2	51
23	Computational design of closely related proteins that adopt two well-defined but structurally divergent folds. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020, 117, 7208-7215.	7.1	48
24	Immune Focusing and Enhanced Neutralization Induced by HIV-1 gp140 Chemical Cross-Linking. <i>Journal of Virology</i> , 2013, 87, 10163-10172.	3.4	43
25	Designing repeat proteins: a modular approach to protein design. <i>Current Opinion in Structural Biology</i> , 2017, 45, 116-123.	5.7	40
26	Computational design and experimental verification of a symmetric protein homodimer. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 10714-10719.	7.1	34
27	Modulation of Integrin Activation by an Entropic Spring in the Î²-Knee. <i>Journal of Biological Chemistry</i> , 2010, 285, 32954-32966.	3.4	32
28	Tight and specific lanthanide binding in a de novo TIM barrel with a large internal cavity designed by symmetric domain fusion. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020, 117, 30362-30369.	7.1	31
29	Adaptation of a fast Fourier transform-based docking algorithm for protein design. <i>Journal of Computational Chemistry</i> , 2005, 26, 1222-1232.	3.3	27
30	HIV-1 VRC01 Germline-Targeting Immunogens Select Distinct Epitope-Specific B Cell Receptors. <i>Immunity</i> , 2020, 53, 840-851.e6.	14.3	27
31	A computationally engineered RAS rheostat reveals RAS-ERK signaling dynamics. <i>Nature Chemical Biology</i> , 2017, 13, 119-126.	8.0	21
32	Computational De Novo Design of a Self-Assembling Peptide with Predefined Structure. <i>Journal of Molecular Biology</i> , 2015, 427, 550-562.	4.2	20
33	The molecular basis of chaperone-mediated interleukin 23 assembly control. <i>Nature Communications</i> , 2019, 10, 4121.	12.8	18
34	Engineering a potent receptor superagonist or antagonist from a novel IL-6 family cytokine ligand. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020, 117, 14110-14118.	7.1	17
35	Optical control of fast and processive engineered myosins in vitro and in living cells. <i>Nature Chemical Biology</i> , 2021, 17, 540-548.	8.0	17
36	A Chimeric HIV-1 Envelope Glycoprotein Trimer with an Embedded Granulocyte-Macrophage Colony-stimulating Factor (GM-CSF) Domain Induces Enhanced Antibody and T Cell Responses. <i>Journal of Biological Chemistry</i> , 2011, 286, 22250-22261.	3.4	15

#	ARTICLE	IF	CITATIONS
37	Domain 1 of Mucosal Addressin Cell Adhesion Molecule Has an I1-set Fold and a Flexible Integrin-binding Loop. <i>Journal of Biological Chemistry</i> , 2013, 288, 6284-6294.	3.4	13
38	A Designed Protein Interface That Blocks Fibril Formation. <i>Journal of the American Chemical Society</i> , 2004, 126, 13914-13915.	13.7	12
39	Interleukin-2 superkines by computational design. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022, 119, e21117401119.	7.1	12
40	Using Molecular Dynamics Simulations as an Aid in the Prediction of Domain Swapping of Computationally Designed Protein Variants. <i>Journal of Molecular Biology</i> , 2015, 427, 2697-2706.	4.2	11
41	Multi-scale structural analysis of proteins by deep semantic segmentation. <i>Bioinformatics</i> , 2020, 36, 1740-1749.	4.1	10
42	Identification of N-Terminally Diversified GLP-1R Agonists Using Saturation Mutagenesis and Chemical Design. <i>ACS Chemical Biology</i> , 2021, 16, 58-66.	3.4	5
43	High-resolution structure prediction of a circular permutation loop. <i>Protein Science</i> , 2011, 20, 1929-1934.	7.6	3
44	Chimeric mutants of staphylococcal hemolysin, which act as both one-component and two-component hemolysin, created by grafting the stem domain. <i>FEBS Journal</i> , 2022, 289, 3505-3520.	4.7	1
45	Refocussing Antibody Responses by Chemical Modification of Vaccine Antigens. <i>AIDS Research and Human Retroviruses</i> , 2014, 30, A66-A67.	1.1	0
46	Harnessing Human Neural Networks for Protein Design. <i>Biochemistry</i> , 2019, 58, 5100-5101.	2.5	0