Pietro Sormanni

List of Publications by Year in descending order

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56 papers	2,951 citations	218677 26 h-index	50 g-index
61	61	61	4139
all docs	docs citations	times ranked	citing authors

#	Article	IF	CITATIONS
1	The CamSol Method of Rational Design of Protein Mutants with Enhanced Solubility. Journal of Molecular Biology, 2015, 427, 478-490.	4.2	341
2	A natural product inhibits the initiation of $\hat{l}\pm$ -synuclein aggregation and suppresses its toxicity. Proceedings of the National Academy of Sciences of the United States of America, 2017, 114, E1009-E1017.	7.1	231
3	MobiDB 3.0: more annotations for intrinsic disorder, conformational diversity and interactions in proteins. Nucleic Acids Research, 2018, 46, D471-D476.	14.5	190
4	Different soluble aggregates of $A\hat{l}^2$ 42 can give rise to cellular toxicity through different mechanisms. Nature Communications, 2019, 10, 1541.	12.8	140
5	Parapred: antibody paratope prediction using convolutional and recurrent neural networks. Bioinformatics, 2018, 34, 2944-2950.	4.1	124
6	Selective targeting of primary and secondary nucleation pathways in \hat{Al}^242 aggregation using a rational antibody scanning method. Science Advances, 2017, 3, e1700488.	10.3	116
7	Proteome-wide observation of the phenomenon of life on the edge of solubility. Proceedings of the National Academy of Sciences of the United States of America, 2020, 117, 1015-1020.	7.1	115
8	Rational design of antibodies targeting specific epitopes within intrinsically disordered proteins. Proceedings of the National Academy of Sciences of the United States of America, 2015, 112, 9902-9907.	7.1	113
9	Simultaneous quantification of protein order and disorder. Nature Chemical Biology, 2017, 13, 339-342.	8.0	113
10	Rapid and accurate in silico solubility screening of a monoclonal antibody library. Scientific Reports, 2017, 7, 8200.	3.3	97
11	Third generation antibody discovery methods: <i>in silico</i> rational design. Chemical Society Reviews, 2018, 47, 9137-9157.	38.1	94
12	Targeting disordered proteins with small molecules using entropy. Trends in Biochemical Sciences, 2015, 40, 491-496.	7.5	87
13	A protein homeostasis signature in healthy brains recapitulates tissue vulnerability to Alzheimer's disease. Science Advances, 2016, 2, e1600947.	10.3	84
14	The s2D Method: Simultaneous Sequence-Based Prediction of the Statistical Populations of Ordered and Disordered Regions in Proteins. Journal of Molecular Biology, 2015, 427, 982-996.	4.2	77
15	In vitro and in silico assessment of the developability of a designed monoclonal antibody library. MAbs, 2019, 11, 388-400.	5. 2	72
16	Rational design of a conformation-specific antibody for the quantification of $A\hat{I}^2$ oligomers. Proceedings of the National Academy of Sciences of the United States of America, 2020, 117, 13509-13518.	7.1	61
17	Quantifying misfolded protein oligomers as drug targets and biomarkers in Alzheimer and Parkinson diseases. Nature Reviews Chemistry, 2021, 5, 277-294.	30.2	56
18	Conformational dynamics in crystals reveal the molecular bases for D76N beta-2 microglobulin aggregation propensity. Nature Communications, 2018, 9, 1658.	12.8	53

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19	Massively parallel C. elegans tracking provides multi-dimensional fingerprints for phenotypic discovery. Journal of Neuroscience Methods, 2018, 306, 57-67.	2.5	52
20	Developability Assessment of Engineered Monoclonal Antibody Variants with a Complex Self-Association Behavior Using Complementary Analytical and in Silico Tools. Molecular Pharmaceutics, 2018, 15, 5697-5710.	4.6	52
21	A chemical kinetic basis for measuring translation initiation and elongation rates from ribosome profiling data. PLoS Computational Biology, 2019, 15, e1007070.	3.2	50
22	Rational design of mutations that change the aggregation rate of a protein while maintaining its native structure and stability. Scientific Reports, 2016, 6, 25559.	3.3	47
23	Attentive Cross-Modal Paratope Prediction. Journal of Computational Biology, 2019, 26, 536-545.	1.6	45
24	Understanding the frustration arising from the competition between function, misfolding, and aggregation in a globular protein. Proceedings of the National Academy of Sciences of the United States of America, 2014, 111, 14141-14146.	7.1	43
25	Protein Solubility Predictions Using the CamSol Method in the Study of Protein Homeostasis. Cold Spring Harbor Perspectives in Biology, 2019, 11, a033845.	5.5	42
26	Oxetane Grafts Installed Siteâ€Selectively on Native Disulfides to Enhance Protein Stability and Activity Inâ€Vivo. Angewandte Chemie - International Edition, 2017, 56, 14963-14967.	13.8	39
27	Current advances in biopharmaceutical informatics: guidelines, impact and challenges in the computational developability assessment of antibody therapeutics. MAbs, 2022, 14, 2020082.	5.2	35
28	Inherent Biophysical Properties Modulate the Toxicity of Soluble Amyloidogenic Light Chains. Journal of Molecular Biology, 2020, 432, 845-860.	4.2	26
29	Assessment of Therapeutic Antibody Developability by Combinations of In Vitro and In Silico Methods. Methods in Molecular Biology, 2022, 2313, 57-113.	0.9	26
30	A Rational Design Strategy for the Selective Activity Enhancement of a Molecular Chaperone toward a Target Substrate. Biochemistry, 2015, 54, 5103-5112.	2.5	25
31	Supersaturated proteins are enriched at synapses and underlie cell and tissue vulnerability in Alzheimer's disease. Heliyon, 2019, 5, e02589.	3.2	23
32	Identifying A- and P-site locations on ribosome-protected mRNA fragments using Integer Programming. Scientific Reports, 2019, 9, 6256.	3.3	18
33	Corneal Dystrophy Mutations Drive Pathogenesis by Targeting TGFBIp Stability and Solubility in a Latent Amyloid-forming Domain. Journal of Molecular Biology, 2018, 430, 1116-1140.	4.2	17
34	Delivery of Native Proteins into C. elegans Using a Transduction Protocol Based on Lipid Vesicles. Scientific Reports, 2017, 7, 15045.	3.3	16
35	An open-source automated PEG precipitation assay to measure the relative solubility of proteins with low material requirement. Scientific Reports, 2021, 11, 21932.	3.3	16
36	Fibrillogenic propensity of the GroEL apical domain: A Janusâ€faced minichaperone. FEBS Letters, 2012, 586, 1120-1127.	2.8	15

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37	A rationally designed bicyclic peptide remodels Aβ42 aggregation in vitro and reduces its toxicity in a worm model of Alzheimer's disease. Scientific Reports, 2020, 10, 15280.	3.3	15
38	Systematic Activity Maturation of a Single-Domain Antibody with Non-canonical Amino Acids through Chemical Mutagenesis. Cell Chemical Biology, 2021, 28, 70-77.e5.	5.2	15
39	A Rationally Designed Hsp70 Variant Rescues the Aggregation-Associated Toxicity of Human IAPP in Cultured Pancreatic Islet β-Cells. International Journal of Molecular Sciences, 2018, 19, 1443.	4.1	14
40	MonteGrappa: An iterative Monte Carlo program to optimize biomolecular potentials in simplified models. Computer Physics Communications, 2015, 186, 93-104.	7.5	12
41	Rationally Designed Antibodies as Research Tools to Study the Structure–Toxicity Relationship of Amyloid-l² Oligomers. International Journal of Molecular Sciences, 2020, 21, 4542.	4.1	12
42	Comparative Studies in the A30P and A53T α-Synuclein C. elegans Strains to Investigate the Molecular Origins of Parkinson's Disease. Frontiers in Cell and Developmental Biology, 2021, 9, 552549.	3.7	12
43	Biochemical and biophysical comparison of human and mouse betaâ€⊋ microglobulin reveals the molecular determinants of low amyloid propensity. FEBS Journal, 2020, 287, 546-560.	4.7	11
44	Correlation between the binding affinity and the conformational entropy of nanobody SARS-CoV-2 spike protein complexes. Proceedings of the National Academy of Sciences of the United States of America, 2022, 119 , .	7.1	11
45	Subdomain Architecture and Stability of a Giant Repeat Protein. Journal of Physical Chemistry B, 2013, 117, 13029-13037.	2.6	10
46	Oxetane Grafts Installed Siteâ€Selectively on Native Disulfides to Enhance Protein Stability and Activity Inâ€Vivo. Angewandte Chemie, 2017, 129, 15159-15163.	2.0	10
47	Neuroserpin and transthyretin are extracellular chaperones that preferentially inhibit amyloid formation. Science Advances, 2021, 7, eabf7606.	10.3	10
48	Pairs of amino acids at the P- and A-sites of the ribosome predictably and causally modulate translation-elongation rates. Journal of Molecular Biology, 2020, 432, 166696.	4.2	9
49	Ï€-Clamp-Mediated Homo- and Heterodimerization of Single-Domain Antibodies via Site-Specific Homobifunctional Conjugation. Journal of the American Chemical Society, 2022, 144, 13026-13031.	13.7	9
50	Conformational Entropy as a Potential Liability of Computationally Designed Antibodies. Biomolecules, 2022, 12, 718.	4.0	8
51	Iterative derivation of effective potentials to sample the conformational space of proteins at atomistic scale. Journal of Chemical Physics, 2014, 140, 195101.	3.0	6
52	A method of predicting the in vitro fibril formation propensity of $A\hat{I}^240$ mutants based on their inclusion body levels in E. coli. Scientific Reports, 2019, 9, 3680.	3.3	6
53	Rationally Designed Bicyclic Peptides Prevent the Conversion of A \hat{l}^2 42 Assemblies Into Fibrillar Structures. Frontiers in Neuroscience, 2021, 15, 623097.	2.8	6
54	An aggregation inhibitor specific to oligomeric intermediates of $A\hat{l}^2$ 42 derived from phage display libraries of stable, small proteins. Proceedings of the National Academy of Sciences of the United States of America, 2022, 119, e2121966119.	7.1	5

#	Article	lF	CITATIONS
55	Modulating the cardiotoxic behaviour of immunoglobulin light chain dimers through point mutations. Amyloid: the International Journal of Experimental and Clinical Investigation: the Official Journal of the International Society of Amyloidosis, 2019, 26, 105-106.	3.0	4
56	Computational maturation of a single-domain antibody against $\hat{Al^2}42$ aggregation. Chemical Science, 2021, 12, 13940-13948.	7.4	4