

Vikram Khipple Mulligan

List of Publications by Citations

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

35 papers	2,431 citations	19 h-index	36 g-index
36 ext. papers	3,373 ext. citations	14 avg, IF	4.84 L-index

#	Paper	IF	Citations
35	The Rosetta All-Atom Energy Function for Macromolecular Modeling and Design. <i>Journal of Chemical Theory and Computation</i> , 2017 , 13, 3031-3048	6.4	486
34	Global analysis of protein folding using massively parallel design, synthesis, and testing. <i>Science</i> , 2017 , 357, 168-175	33.3	241
33	Accurate de novo design of hyperstable constrained peptides. <i>Nature</i> , 2016 , 538, 329-335	50.4	231
32	Simultaneous Optimization of Biomolecular Energy Functions on Features from Small Molecules and Macromolecules. <i>Journal of Chemical Theory and Computation</i> , 2016 , 12, 6201-6212	6.4	199
31	Macromolecular modeling and design in Rosetta: recent methods and frameworks. <i>Nature Methods</i> , 2020 , 17, 665-680	21.6	165
30	De novo design of bioactive protein switches. <i>Nature</i> , 2019 , 572, 205-210	50.4	113
29	Prion disease susceptibility is affected by beta-structure folding propensity and local side-chain interactions in PrP. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010 , 107, 19808-13	11.5	102
28	Comprehensive computational design of ordered peptide macrocycles. <i>Science</i> , 2017 , 358, 1461-1466	33.3	96
27	Accurate computational design of multipass transmembrane proteins. <i>Science</i> , 2018 , 359, 1042-1046	33.3	93
26	Programmable design of orthogonal protein heterodimers. <i>Nature</i> , 2019 , 565, 106-111	50.4	87
25	Computational design of an unnatural amino acid dependent metalloprotein with atomic level accuracy. <i>Journal of the American Chemical Society</i> , 2013 , 135, 13393-9	16.4	84
24	Correction: Drosophila melanogaster Cad99C, the orthologue of human Usher cadherin PCDH15, regulates the length of microvilli. <i>Journal of Cell Biology</i> , 2005 , 171, 1085-1085	7.3	78
23	Protein misfolding in the late-onset neurodegenerative diseases: common themes and the unique case of amyotrophic lateral sclerosis. <i>Proteins: Structure, Function and Bioinformatics</i> , 2013 , 81, 1285-303	4.2	62
22	Drosophila melanogaster Cad99C, the orthologue of human Usher cadherin PCDH15, regulates the length of microvilli. <i>Journal of Cell Biology</i> , 2005 , 171, 549-58	7.3	59
21	CCM3/PDCD10 heterodimerizes with germinal center kinase III (GCKIII) proteins using a mechanism analogous to CCM3 homodimerization. <i>Journal of Biological Chemistry</i> , 2011 , 286, 25056-64	5.4	50
20	De novo design of covalently constrained mesosize protein scaffolds with unique tertiary structures. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017 , 114, 10852-10857	11.5	44
19	Denaturational stress induces formation of zinc-deficient monomers of Cu,Zn superoxide dismutase: implications for pathogenesis in amyotrophic lateral sclerosis. <i>Journal of Molecular Biology</i> , 2008 , 383, 424-36	6.5	40

18	Early steps in oxidation-induced SOD1 misfolding: implications for non-amyloid protein aggregation in familial ALS. <i>Journal of Molecular Biology</i> , 2012 , 421, 631-52	6.5	38
17	ALS-causing SOD1 mutations promote production of copper-deficient misfolded species. <i>Journal of Molecular Biology</i> , 2011 , 409, 839-52	6.5	33
16	Computationally designed peptide macrocycle inhibitors of New Delhi metallo- β -lactamase 1. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021 , 118,	11.5	17
15	The emerging role of computational design in peptide macrocycle drug discovery. <i>Expert Opinion on Drug Discovery</i> , 2020 , 15, 833-852	6.2	15
14	Better together: Elements of successful scientific software development in a distributed collaborative community. <i>PLoS Computational Biology</i> , 2020 , 16, e1007507	5	15
13	A systematic study of minima in alanine dipeptide. <i>Journal of Computational Chemistry</i> , 2019 , 40, 297-309.	9.5	15
12	Designing Peptides on a Quantum Computer		13
11	Anchor extension: a structure-guided approach to design cyclic peptides targeting enzyme active sites. <i>Nature Communications</i> , 2021 , 12, 3384	17.4	12
10	Conversion of Abeta42 into a folded soluble native-like protein using a semi-random library of amphipathic helices. <i>Journal of Molecular Biology</i> , 2010 , 396, 1284-94	6.5	10
9	Computational design of mixed chirality peptide macrocycles with internal symmetry. <i>Protein Science</i> , 2020 , 29, 2433-2445	6.3	9
8	Analyzing complicated protein folding kinetics rapidly by analytical Laplace inversion using a Tikhonov regularization variant. <i>Analytical Biochemistry</i> , 2012 , 421, 181-90	3.1	7
7	MHCEpitopeEnergy, a Flexible Rosetta-Based Biotherapeutic Deimmunization Platform. <i>Journal of Chemical Information and Modeling</i> , 2021 , 61, 2368-2382	6.1	6
6	The Rosetta all-atom energy function for macromolecular modeling and design		3
5	Current directions in combining simulation-based macromolecular modeling approaches with deep learning. <i>Expert Opinion on Drug Discovery</i> , 2021 , 16, 1025-1044	6.2	3
4	Ensuring scientific reproducibility in bio-macromolecular modeling via extensive, automated benchmarks		2
3	XENet: Using a new graph convolution to accelerate the timeline for protein design on quantum computers. <i>PLoS Computational Biology</i> , 2021 , 17, e1009037	5	2
2	A computational method for the design of nested proteins by loop-directed domain insertion. <i>Proteins: Structure, Function and Bioinformatics</i> , 2018 , 86, 354-369	4.2	1
1	Ensuring scientific reproducibility in bio-macromolecular modeling via extensive, automated benchmarks. <i>Nature Communications</i> , 2021 , 12, 6947	17.4	0

