

# Gaohua Liu

## List of Publications by Year in descending order

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Version: 2024-02-01

15  
papers

540  
citations

933447

10  
h-index

1058476

14  
g-index

17  
all docs

17  
docs citations

17  
times ranked

983  
citing authors

#	ARTICLE	IF	CITATIONS
1	Principles for designing proteins with cavities formed by curved $\hat{I}^2$ sheets. <i>Science</i> , 2017, 355, 201-206.	12.6	117
2	Control over overall shape and size in de novo designed proteins. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, E5478-85.	7.1	113
3	De novo protein design by citizen scientists. <i>Nature</i> , 2019, 570, 390-394.	27.8	105
4	Structural/Functional Properties of Human NFU1, an Intermediate [4Fe-4S] Carrier in Human Mitochondrial Iron-Sulfur Cluster Biogenesis. <i>Structure</i> , 2016, 24, 2080-2091.	3.3	45
5	Assessment of prediction methods for protein structures determined by <a href="#">NMR</a> in <a href="#">CASP14</a> : Impact of <a href="#">AlphaFold2</a> . <i>Proteins: Structure, Function and Bioinformatics</i> , 2021, 89, 1959-1976.	2.6	30
6	The second round of Critical Assessment of Automated Structure Determination of Proteins by NMR: CASD-NMR-2013. <i>Journal of Biomolecular NMR</i> , 2015, 62, 413-424.	2.8	27
7	A community resource of experimental data for <a href="#">NMR</a> / <a href="#">X-ray</a> crystal structure pairs. <i>Protein Science</i> , 2016, 25, 30-45.	7.6	24
8	Protein structure prediction assisted with sparse NMR data in CASP13. <i>Proteins: Structure, Function and Bioinformatics</i> , 2019, 87, 1315-1332.	2.6	21
9	Solution NMR structure of the ARID domain of human AT-rich interactive domain-containing protein 3A: A human cancer protein interaction network target. <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 2170-2175.	2.6	16
10	A common binding motif in the ET domain of BRD3 forms polymorphic structural interfaces with host and viral proteins. <i>Structure</i> , 2021, 29, 886-898.e6.	3.3	16
11	REDCRAFT: A computational platform using residual dipolar coupling NMR data for determining structures of perdeuterated proteins in solution. <i>PLoS Computational Biology</i> , 2021, 17, e1008060.	3.2	8
12	NMR structure of <a href="#">F-actin</a> binding domain of Arg/Abl2 from <i>Homo sapiens</i> . <i>Proteins: Structure, Function and Bioinformatics</i> , 2010, 78, 1326-1330.	2.6	7
13	Introduction of a polar core into the de novo designed protein <a href="#">T</a> <sub>op7</sub> . <i>Protein Science</i> , 2016, 25, 1299-1307.	7.6	7
14	Backbone and Ile- $\hat{I}^1$ , Leu, Val methyl $^1H$ , $^{15}N$ , and $^{13}C$ , chemical shift assignments for <i>Rhizopus chinensis</i> lipase. <i>Biomolecular NMR Assignments</i> , 2018, 12, 63-68.	0.8	3
15	Aromatic claw: A new fold with high aromatic content that evades structural prediction. <i>Protein Science</i> , 2017, 26, 208-217.	7.6	0