

# Chun Tang

## List of Publications by Year in descending order

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76  
papers

5,021  
citations

159585

30  
h-index

95266

68  
g-index

91  
all docs

91  
docs citations

91  
times ranked

6014  
citing authors

#	ARTICLE	IF	CITATIONS
1	Preferential Interactions of a Crowder Protein with the Specific Binding Site of a Native Protein Complex. <i>Journal of Physical Chemistry Letters</i> , 2022, 13, 792-800.	4.6	8
2	Nucleobase Clustering Contributes to the Formation and Hollowing of Repeat-Expansion RNA Condensate. <i>Journal of the American Chemical Society</i> , 2022, 144, 4716-4720.	13.7	14
3	Characterization of protein unfolding by fast cross-linking mass spectrometry using di-ortho-phthalaldehyde cross-linkers. <i>Nature Communications</i> , 2022, 13, 1468.	12.8	20
4	Preferential Regulation of Transient Protein-Protein Interaction by the Macromolecular Crowders. <i>Journal of Physical Chemistry B</i> , 2022, 126, 4840-4848.	2.6	1
5	FRET-based dynamic structural biology: Challenges, perspectives and an appeal for open-science practices. <i>ELife</i> , 2021, 10, .	6.0	152
6	Kinetic Constraints in the Specific Interaction between Phosphorylated Ubiquitin and Proteasomal Shuttle Factors. <i>Biomolecules</i> , 2021, 11, 1008.	4.0	2
7	Pseudopotentials for coarse-grained cross-link-assisted modeling of protein structures. <i>Journal of Computational Chemistry</i> , 2021, 42, 2054-2067.	3.3	4
8	Discovery of a Novel Androgen Receptor Antagonist Manifesting Evidence to Disrupt the Dimerization of the Ligand-Binding Domain via Attenuating the Hydrogen-Bonding Network Between the Two Monomers. <i>Journal of Medicinal Chemistry</i> , 2021, 64, 17221-17238.	6.4	7
9	Recent Developments in Data-Assisted Modeling of Flexible Proteins. <i>Frontiers in Molecular Biosciences</i> , 2021, 8, 765562.	3.5	8
10	How Phosphorylation by PINK1 Remodels the Ubiquitin System: A Perspective from Structure and Dynamics. <i>Biochemistry</i> , 2020, 59, 26-33.	2.5	9
11	Tightening the Crosslinking Distance Restraints for Better Resolution of Protein Structure and Dynamics. <i>Structure</i> , 2020, 28, 1160-1167.e3.	3.3	15
12	Hierarchical Conformational Dynamics Confers Thermal Adaptability to preQ1 RNA Riboswitches. <i>Journal of Molecular Biology</i> , 2020, 432, 4523-4543.	4.2	1
13	The Conformational Preference of Chemical Cross-linkers Determines the Cross-linking Probability of Reactive Protein Residues. <i>Journal of Physical Chemistry B</i> , 2020, 124, 4446-4453.	2.6	7
14	FLIM-FRET-Based Structural Characterization of a Class-A GPCR Dimer in the Cell Membrane. <i>Journal of Molecular Biology</i> , 2020, 432, 4596-4611.	4.2	9
15	Integrating Non-NMR Distance Restraints to Augment NMR Depiction of Protein Structure and Dynamics. <i>Journal of Molecular Biology</i> , 2020, 432, 2913-2929.	4.2	13
16	Ubiquitin is double-phosphorylated by PINK1 for enhanced pH-sensitivity of conformational switch. <i>Protein and Cell</i> , 2019, 10, 908-913.	11.0	6
17	Nicotinamide phosphoribosyltransferase secreted from microglia <i>via</i> exosome during ischemic injury. <i>Journal of Neurochemistry</i> , 2019, 150, 723-737.	3.9	28
18	On the necessity of an integrative approach to understand protein structural dynamics. <i>Journal of Zhejiang University: Science B</i> , 2019, 20, 496-502.	2.8	5

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19	Structural insights into DNA recognition by AimR of the arbitrium communication system in the SPbeta phage. <i>Cell Discovery</i> , 2019, 5, 29.	6.7	12
20	Structural basis for the recognition of K48-linked Ub chain by proteasomal receptor Rpn13. <i>Cell Discovery</i> , 2019, 5, 19.	6.7	27
21	Refining RNA solution structures with the integrative use of label-free paramagnetic relaxation enhancement NMR. <i>Biophysics Reports</i> , 2019, 5, 244-253.	0.8	4
22	Requirement for p62 acetylation in the aggregation of ubiquitylated proteins under nutrient stress. <i>Nature Communications</i> , 2019, 10, 5792.	12.8	83
23	Solution structure of the RNA recognition domain of METTL3-METTL14 N6-methyladenosine methyltransferase. <i>Protein and Cell</i> , 2019, 10, 272-284.	11.0	99
24	Theory and practice of using solvent paramagnetic relaxation enhancement to characterize protein conformational dynamics. <i>Methods</i> , 2018, 148, 48-56.	3.8	28
25	Carboxylate-Selective Chemical Cross-Linkers for Mass Spectrometric Analysis of Protein Structures. <i>Analytical Chemistry</i> , 2018, 90, 1195-1201.	6.5	42
26	Characterizing Protein Dynamics with Integrative Use of Bulk and Single-Molecule Techniques. <i>Biochemistry</i> , 2018, 57, 305-313.	2.5	21
27	Understanding the graphene quantum dots-ubiquitin interaction by identifying the interaction sites. <i>Carbon</i> , 2017, 121, 285-291.	10.3	17
28	Ubiquitin S65 phosphorylation engenders a pH-sensitive conformational switch. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, 6770-6775.	7.1	40
29	Lanthanoid tagging via an unnatural amino acid for protein structure characterization. <i>Journal of Biomolecular NMR</i> , 2017, 67, 273-282.	2.8	17
30	Lipid-dependent conformational dynamics underlie the functional versatility of T-cell receptor. <i>Cell Research</i> , 2017, 27, 505-525.	12.0	38
31	Modeling Protein Excited-state Structures from $\alpha$ -Over-length $\beta$ -Chemical Cross-links. <i>Journal of Biological Chemistry</i> , 2017, 292, 1187-1196.	3.4	48
32	Protein Structural Ensembles Visualized by Solvent Paramagnetic Relaxation Enhancement. <i>Angewandte Chemie</i> , 2017, 129, 1022-1026.	2.0	4
33	Protein Structural Ensembles Visualized by Solvent Paramagnetic Relaxation Enhancement. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 1002-1006.	13.8	34
34	Protocol for analyzing protein ensemble structures from chemical cross-links using DynaXL. <i>Biophysics Reports</i> , 2017, 3, 100-108.	0.8	10
35	Specific cell surface labeling of GPCRs using split GFP. <i>Scientific Reports</i> , 2016, 6, 20568.	3.3	15
36	Estimation of protein dynamic states with single molecule fluorescence data analysis at microsecond scale., 2016, , .		0

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37	Structural basis of N6-adenosine methylation by the METTL3&#x2013;METTL14 complex. <i>Nature</i> , 2016, 534, 575-578.	27.8	807
38	Structural basis of nonribosomal peptide macrocyclization in fungi. <i>Nature Chemical Biology</i> , 2016, 12, 1001-1003.	8.0	54
39	Hemi-methylated DNA opens a closed conformation of UHRF1 to facilitate its histone recognition. <i>Nature Communications</i> , 2016, 7, 11197.	12.8	100
40	Ensemble structure description of Lys63-linked diubiquitin. <i>Data in Brief</i> , 2016, 7, 81-88.	1.0	6
41	Transient protein&#x2013;protein interactions visualized by solution NMR. <i>Biochimica Et Biophysica Acta - Proteins and Proteomics</i> , 2016, 1864, 115-122.	2.3	49
42	Visualizing the Ensemble Structures of Protein Complexes Using Chemical Cross-Linking Coupled with Mass Spectrometry. <i>Biophysics Reports</i> , 2015, 1, 127-138.	0.8	26
43	NAMPT inhibitor and metabolite protect mouse brain from cryoinjury through distinct mechanisms. <i>Neuroscience</i> , 2015, 291, 230-240.	2.3	18
44	Conjoined Use of EM and NMR in RNA Structure Refinement. <i>PLoS ONE</i> , 2015, 10, e0120445.	2.5	24
45	Lys63-linked ubiquitin chain adopts multiple conformational states for specific target recognition. <i>ELife</i> , 2015, 4, .	6.0	50
46	PolyUbiquitin Chain Linkage Topology Selects the Functions from the Underlying Binding Landscape. <i>PLoS Computational Biology</i> , 2014, 10, e1003691.	3.2	30
47	A decadentate Gd(III)-coordinating paramagnetic cosolvent for protein relaxation enhancement measurement. <i>Journal of Biomolecular NMR</i> , 2014, 58, 149-154.	2.8	17
48	Characterization of the Raptor/4E-BP1 Interaction by Chemical Cross-linking Coupled with Mass Spectrometry Analysis. <i>Journal of Biological Chemistry</i> , 2014, 289, 4723-4734.	3.4	14
49	Subtle Dynamics of <i>holo</i> Glutamine Binding Protein Revealed with a Rigid Paramagnetic Probe. <i>Biochemistry</i> , 2014, 53, 1403-1409.	2.5	19
50	Visualizing an Ultra&#x2013;Weak Protein&#x2013;Protein Interaction in Phosphorylation Signaling. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 11501-11505.	13.8	24
51	NMR Model of Prgl&#x2013;SipD Interaction and Its Implications in the Needle-Tip Assembly of the Salmonella Type III Secretion System. <i>Journal of Molecular Biology</i> , 2014, 426, 2958-2969.	4.2	36
52	Molecular mechanism for Rabex-5 GEF activation by Rabaptin-5. <i>ELife</i> , 2014, 3, .	6.0	47
53	Protein dynamics elucidated by NMR technique. <i>Protein and Cell</i> , 2013, 4, 726-730.	11.0	6
54	NASR: An Effective Approach for Simultaneous Noise and Artifact Suppression in NMR Spectroscopy. <i>Analytical Chemistry</i> , 2013, 85, 2523-2528.	6.5	15

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55	Cerebral Ischemia Is Exacerbated by Extracellular Nicotinamide Phosphoribosyltransferase via a Non-Enzymatic Mechanism. PLoS ONE, 2013, 8, e85403.	2.5	24
56	Exploration of Multi-State Conformational Dynamics and Underlying Global Functional Landscape of Maltose Binding Protein. PLoS Computational Biology, 2012, 8, e1002471.	3.2	50
57	Noncovalent Dimerization of Ubiquitin. Angewandte Chemie - International Edition, 2012, 51, 469-472.	13.8	80
58	Characterization of the Interaction between the Salmonella Type III Secretion System Tip Protein SipD and the Needle Protein PrgI by Paramagnetic Relaxation Enhancement. Journal of Biological Chemistry, 2011, 286, 4922-4930.	3.4	30
59	Characterizing Dynamic Protein-Protein Interactions Using Differentially Scaled Paramagnetic Relaxation Enhancement. Journal of the American Chemical Society, 2009, 131, 17291-17297.	13.7	30
60	Visualizing transient events in amino-terminal autoprocessing of HIV-1 protease. Nature, 2008, 455, 693-696.	27.8	123
61	Visualization of Transient Ultra-Weak Protein Self-Association in Solution Using Paramagnetic Relaxation Enhancement. Journal of the American Chemical Society, 2008, 130, 4048-4056.	13.7	80
62	Replica exchange simulations of transient encounter complexes in protein-protein association. Proceedings of the National Academy of Sciences of the United States of America, 2008, 105, 12855-12860.	7.1	107
63	Structure of the Antiviral Assembly Inhibitor CAP-1 Complex with the HIV-1 CA Protein. Journal of Molecular Biology, 2007, 373, 355-366.	4.2	144
64	Role of Electrostatic Interactions in Transient Encounter Complexes in Protein-Protein Association Investigated by Paramagnetic Relaxation Enhancement. Journal of the American Chemical Society, 2007, 129, 12954-12955.	13.7	73
65	Two Distinct Buckling Modes in Carbon Nanotube Bending. Nano Letters, 2007, 7, 143-148.	9.1	62
66	Practical aspects of 1H transverse paramagnetic relaxation enhancement measurements on macromolecules. Journal of Magnetic Resonance, 2007, 184, 185-195.	2.1	239
67	Open-to-closed transition in apo maltose-binding protein observed by paramagnetic NMR. Nature, 2007, 449, 1078-1082.	27.8	390
68	Elucidating transient macromolecular interactions using paramagnetic relaxation enhancement. Current Opinion in Structural Biology, 2007, 17, 603-616.	5.7	201
69	Visualization of transient encounter complexes in protein-protein association. Nature, 2006, 444, 383-386.	27.8	397
70	A simple and reliable approach to docking protein-protein complexes from very sparse NOE-derived intermolecular distance restraints. Journal of Biomolecular NMR, 2006, 36, 37-44.	2.8	32
71	Accurate Determination of Leucine and Valine Side-chain Conformations using U-[15N/13C/2H]/[1H-(methine/methyl)-Leu/Val] Isotope Labeling, NOE Pattern Recognition, and Methine C-H/C-H Residual Dipolar Couplings: Application to the 34-kDa Enzyme II-Achitobiose. Journal of Biomolecular NMR, 2005, 33, 105-121.	2.8	12
72	Solution Structure of Enzyme II-Achitobiose from the N, N'-Diacetylchitobiose Branch of the Escherichia coli Phosphotransferase System. Journal of Biological Chemistry, 2005, 280, 11770-11780.	3.4	9

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73	Entropic switch regulates myristate exposure in the HIV-1 matrix protein. Proceedings of the National Academy of Sciences of the United States of America, 2004, 101, 517-522.	7.1	293
74	Antiviral Inhibition of the HIV-1 Capsid Protein. Journal of Molecular Biology, 2003, 327, 1013-1020.	4.2	204
75	Structure of the N-terminal 283-residue fragment of the immature HIV-1 Gag polyprotein. Nature Structural Biology, 2002, 9, 537-43.	9.7	151
76	Two Methods to Synthesize C60Nitroxide Derivatives. Fullerenes, Nanotubes, and Carbon Nanostructures, 1999, 7, 297-303.	0.6	0