

Nico Tjandra

List of Publications by Year in descending order

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89
papers

10,162
citations

126907

33
h-index

54911

84
g-index

92
all docs

92
docs citations

92
times ranked

8847
citing authors

#	ARTICLE	IF	CITATIONS
1	The Xplor-NIH NMR molecular structure determination package. <i>Journal of Magnetic Resonance</i> , 2003, 160, 65-73.	2.1	2,165
2	Structure of Bax. <i>Cell</i> , 2000, 103, 645-654.	28.9	1,008
3	Rotational diffusion anisotropy of human ubiquitin from ¹⁵ N NMR relaxation. <i>Journal of the American Chemical Society</i> , 1995, 117, 12562-12566.	13.7	678
4	Solution structure of calcium-free calmodulin. <i>Nature Structural and Molecular Biology</i> , 1995, 2, 768-776.	8.2	677
5	BAX activation is initiated at a novel interaction site. <i>Nature</i> , 2008, 455, 1076-1081.	27.8	617
6	Bcl-xL Retrotranslocates Bax from the Mitochondria into the Cytosol. <i>Cell</i> , 2011, 145, 104-116.	28.9	512
7	Use of dipolar ¹ H- ¹⁵ N and ¹ H- ¹³ C couplings in the structure determination of magnetically oriented macromolecules in solution. <i>Nature Structural Biology</i> , 1997, 4, 732-738.	9.7	456
8	Dipolar Couplings in Macromolecular Structure Determination. <i>Methods in Enzymology</i> , 2001, 339, 127-174.	1.0	388
9	Magnetic Field Dependence of Nitrogen- ¹⁵ Proton Splittings in ¹⁵ N-Enriched Human Ubiquitin Resulting from Relaxation Interference and Residual Dipolar Coupling. <i>Journal of the American Chemical Society</i> , 1996, 118, 6264-6272.	13.7	318
10	Protein Backbone Dynamics and ¹⁵ N Chemical Shift Anisotropy from Quantitative Measurement of Relaxation Interference Effects. <i>Journal of the American Chemical Society</i> , 1996, 118, 6986-6991.	13.7	317
11	Anisotropic rotational diffusion of perdeuterated HIV protease from ¹⁵ N NMR relaxation measurements at two magnetic fields. <i>Journal of Biomolecular NMR</i> , 1996, 8, 273-284.	2.8	236
12	Residual Dipolar Couplings in NMR Structure Analysis. <i>Annual Review of Biophysics and Biomolecular Structure</i> , 2004, 33, 387-413.	18.3	193
13	High-resolution heteronuclear NMR of human ubiquitin in an aqueous liquid crystalline medium. <i>Journal of Biomolecular NMR</i> , 1997, 10, 289-292.	2.8	176
14	Defining long range order in NMR structure determination from the dependence of heteronuclear relaxation times on rotational diffusion anisotropy. <i>Nature Structural Biology</i> , 1997, 4, 443-449.	9.7	174
15	Direct Measurement of ¹⁵ N Chemical Shift Anisotropy in Solution. <i>Journal of the American Chemical Society</i> , 1998, 120, 10947-10952.	13.7	154
16	Refined solution structure and backbone dynamics of HIV-1 Nef. <i>Protein Science</i> , 1997, 6, 1248-1263.	7.6	146
17	NMR dipolar couplings for the structure determination of biopolymers in solution. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2002, 40, 175-197.	7.5	145
18	Solution NMR Measurement of Amide Proton Chemical Shift Anisotropy in ¹⁵ N-Enriched Proteins. Correlation with Hydrogen Bond Length. <i>Journal of the American Chemical Society</i> , 1997, 119, 8076-8082.	13.7	134

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19	Analysis of Slow Interdomain Motion of Macromolecules Using NMR Relaxation Data. <i>Journal of the American Chemical Society</i> , 2001, 123, 3953-3959.	13.7	130
20	The Use of Residual Dipolar Coupling in Studying Proteins by NMR. <i>Topics in Current Chemistry</i> , 2011, 326, 47-67.	4.0	100
21	An Approach to Direct Determination of Protein Dynamics from ¹⁵ N NMR Relaxation at Multiple Fields, Independent of Variable ¹⁵ N Chemical Shift Anisotropy and Chemical Exchange Contributions. <i>Journal of the American Chemical Society</i> , 1999, 121, 8577-8582.	13.7	84
22	The use of dipolar couplings for determining the solution structure of rat apo- α 100B(β 2 β 2). <i>Protein Science</i> , 1999, 8, 800-809.	7.6	77
23	Temperature Dependence of Domain Motions of Calmodulin Probed by NMR Relaxation at Multiple Fields. <i>Journal of the American Chemical Society</i> , 2003, 125, 11379-11384.	13.7	68
24	Structure of Transmembrane Domain of Lysosome-associated Membrane Protein Type 2a (LAMP-2A) Reveals Key Features for Substrate Specificity in Chaperone-mediated Autophagy. <i>Journal of Biological Chemistry</i> , 2014, 289, 35111-35123.	3.4	63
25	Conformational Rearrangements in the Pro-apoptotic Protein, Bax, as It Inserts into Mitochondria. <i>Journal of Biological Chemistry</i> , 2014, 289, 32871-32882.	3.4	61
26	Determining the Magnitude of the Fully Asymmetric Diffusion Tensor from Heteronuclear Relaxation Data in the Absence of Structural Information. <i>Journal of the American Chemical Society</i> , 1998, 120, 4889-4890.	13.7	59
27	Structural Insights of tBid, the Caspase-8-activated Bid, and Its BH3 Domain. <i>Journal of Biological Chemistry</i> , 2013, 288, 35840-35851.	3.4	59
28	Structural mechanism of Bax inhibition by cytomegalovirus protein vMIA. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2012, 109, 20901-20906.	7.1	53
29	Bcl-2 proteins bid and bax form a network to permeabilize the mitochondria at the onset of apoptosis. <i>Cell Death and Disease</i> , 2016, 7, e2424-e2424.	6.3	49
30	Carbonyl CSA Restraints from Solution NMR for Protein Structure Refinement. <i>Journal of the American Chemical Society</i> , 2001, 123, 11065-11066.	13.7	44
31	Are proteins even floppier than we thought?. <i>Nature Structural Biology</i> , 1997, 4, 254-256.	9.7	43
32	¹⁵ N chemical shift anisotropy in protein structure refinement and comparison with NH residual dipolar couplings. <i>Journal of Magnetic Resonance</i> , 2003, 164, 171-176.	2.1	41
33	Analysis of the isomer ratios of polymethylated-DOTA complexes and the implications on protein structural studies. <i>Dalton Transactions</i> , 2016, 45, 4673-4687.	3.3	38
34	Tsg101 chaperone function revealed by HIV-1 assembly inhibitors. <i>Nature Communications</i> , 2017, 8, 1391.	12.8	37
35	Structural basis for polyglutamate chain initiation and elongation by TTL family enzymes. <i>Nature Structural and Molecular Biology</i> , 2020, 27, 802-813.	8.2	35
36	Determination of the Solution-Bound Conformation of an Amino Acid Binding Protein by NMR Paramagnetic Relaxation Enhancement: Use of a Single Flexible Paramagnetic Probe with Improved Estimation of Its Sampling Space. <i>Journal of the American Chemical Society</i> , 2009, 131, 9532-9537.	13.7	32

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37	Parameterization of solvent-protein interaction and its use on NMR protein structure determination. <i>Journal of Magnetic Resonance</i> , 2012, 221, 76-84.	2.1	31
38	Selective Targeting of Virus Replication by Proton Pump Inhibitors. <i>Scientific Reports</i> , 2020, 10, 4003.	3.3	31
39	Increasing the Chemical Shift Dispersion of Unstructured Proteins with a Covalent Lanthanide Shift Reagent. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 14847-14851.	13.8	29
40	Conformational Ensemble of Disordered Proteins Probed by Solvent Paramagnetic Relaxation Enhancement (sPRE). <i>Angewandte Chemie - International Edition</i> , 2018, 57, 13519-13522.	13.8	28
41	Humanin induces conformational changes in the apoptosis regulator BAX and sequesters it into fibers, preventing mitochondrial outer-membrane permeabilization. <i>Journal of Biological Chemistry</i> , 2019, 294, 19055-19065.	3.4	27
42	Estimation of Interdomain Flexibility of N-Terminus of Factor H Using Residual Dipolar Couplings. <i>Biochemistry</i> , 2011, 50, 8138-8149.	2.5	26
43	Extended Model Free Approach To Analyze Correlation Functions of Multidomain Proteins in the Presence of Motional Coupling. <i>Journal of the American Chemical Society</i> , 2008, 130, 12745-12751.	13.7	24
44	Sirt1 carboxyl-domain is an ATP-repressible domain that is transferrable to other proteins. <i>Nature Communications</i> , 2017, 8, 15560.	12.8	24
45	Characterizing the magnetic susceptibility tensor of lanthanide-containing polymethylated-DOTA complexes. <i>Journal of Biomolecular NMR</i> , 2016, 66, 125-139.	2.8	23
46	Analysis of NMR Relaxation Data of Biomolecules with Slow Domain Motions Using Wobble-in-a-Cone Approximation. <i>Journal of the American Chemical Society</i> , 2001, 123, 11484-11485.	13.7	22
47	Characterization of the membrane-inserted C-terminus of cytoprotective BCL-XL. <i>Protein Expression and Purification</i> , 2016, 122, 56-63.	1.3	22
48	Residue-Specific ^{13}C CSA Tensor Principal Components for Ubiquitin: Correlation between Tensor Components and Hydrogen Bonding. <i>Journal of the American Chemical Society</i> , 2007, 129, 1321-1326.	13.7	20
49	Application of Solution NMR Spectroscopy to Study Protein Dynamics. <i>Entropy</i> , 2012, 14, 581-598.	2.2	20
50	Simple multidimensional NMR experiments to obtain different types of one-bond dipolar couplings simultaneously. <i>Journal of Biomolecular NMR</i> , 2001, 19, 63-67.	2.8	19
51	Flexible IgE epitope-containing domains of Phl p 5 cause high allergenic activity. <i>Journal of Allergy and Clinical Immunology</i> , 2017, 140, 1187-1191.	2.9	19
52	Structure of the NPR:EIN N^{tr} Complex: Mechanism for Specificity in Paralogous Phosphotransferase Systems. <i>Structure</i> , 2016, 24, 2127-2137.	3.3	16
53	Long-Range RNA Structural Information via a Paramagnetically Tagged Reporter Protein. <i>Journal of the American Chemical Society</i> , 2019, 141, 1430-1434.	13.7	16
54	Humanin selectively prevents the activation of pro-apoptotic protein BID by sequestering it into fibers. <i>Journal of Biological Chemistry</i> , 2020, 295, 18226-18238.	3.4	16

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55	Determination of the residue-specific ¹⁵ N CSA tensor principal components using multiple alignment media. <i>Journal of Biomolecular NMR</i> , 2006, 35, 249-259.	2.8	14
56	The fluorescent aptamer Squash extensively repurposes the adenine riboswitch fold. <i>Nature Chemical Biology</i> , 2022, 18, 191-198.	8.0	12
57	NMR Analysis of Apo Glutamine Binding Protein Exposes Challenges in the Study of Interdomain Dynamics. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 16899-16902.	13.8	10
58	Single color FRET based measurements of conformational changes of proteins resulting from translocation inside cells. <i>Methods</i> , 2014, 66, 180-187.	3.8	9
59	Prazoles Targeting Tsg101 Inhibit Release of Epstein-Barr Virus following Reactivation from Latency. <i>Journal of Virology</i> , 2021, 95, e0246620.	3.4	9
60	Rotational Dynamics of Calcium-Free Calmodulin Studied by ¹⁵ N NMR Relaxation Measurements. <i>FEBS Journal</i> , 1995, 230, 1014-1024.	0.2	8
61	A practical implementation of cross-spectrum in protein backbone resonance assignment. <i>Journal of Magnetic Resonance</i> , 2010, 203, 208-212.	2.1	8
62	Potential Regulatory Role of Competitive Encounter Complexes in Paralogous Phosphotransferase Systems. <i>Journal of Molecular Biology</i> , 2019, 431, 2331-2342.	4.2	8
63	Backbone ¹⁵ N relaxation analysis of the N-terminal domain of the HTLV-I capsid protein and comparison with the capsid protein of HIV-1. <i>Protein Science</i> , 2003, 12, 973-981.	7.6	7
64	Conformational Heterogeneity in the Activation Mechanism of Bax. <i>Structure</i> , 2017, 25, 1310-1316.e3.	3.3	7
65	Nuclear Magnetic Resonance Solution Structure of the Recombinant Fragment Containing Three Fibrin-Binding Cysteine-Rich Domains of the Very Low Density Lipoprotein Receptor. <i>Biochemistry</i> , 2018, 57, 4395-4403.	2.5	7
66	Comparison of Solution Properties of Polymethylated DOTA-like Lanthanide Complexes with Opposite Chirality of the Pendant Arms. <i>Inorganic Chemistry</i> , 2019, 58, 15788-15800.	4.0	7
67	Bax expression is optimal at low oxygen tension and constant agitation. <i>Protein Expression and Purification</i> , 2020, 165, 105501.	1.3	7
68	Novel Tsg101 Binding Partners Regulate Viral L Domain Trafficking. <i>Viruses</i> , 2021, 13, 1147.	3.3	7
69	Top-down approach in protein RDC data analysis: de novo estimation of the alignment tensor. <i>Journal of Biomolecular NMR</i> , 2007, 38, 303-313.	2.8	6
70	The Structure of Melanoregulin Reveals a Role for Cholesterol Recognition in the Protein's Ability to Promote Dynein Function. <i>Structure</i> , 2018, 26, 1373-1383.e4.	3.3	6
71	Model of a Kinetically Driven Crosstalk between Paralogous Protein Encounter Complexes. <i>Biophysical Journal</i> , 2019, 117, 1655-1665.	0.5	6
72	RNA Binding Suppresses Tsg101 Recognition of Ub-Modified Gag and Facilitates Recruitment to the Plasma Membrane. <i>Viruses</i> , 2020, 12, 447.	3.3	6

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73	Conformational Ensemble of Disordered Proteins Probed by Solvent Paramagnetic Relaxation Enhancement (sPRE). <i>Angewandte Chemie</i> , 2018, 130, 13707-13710.	2.0	5
74	Tsg101/ESCRT-I recruitment regulated by the dual binding modes of K63-linked diubiquitin. <i>Structure</i> , 2022, 30, 289-299.e6.	3.3	5
75	Refinement of protein structure against non-redundant carbonyl ¹³ C NMR relaxation. <i>Journal of Biomolecular NMR</i> , 2007, 38, 243-253.	2.8	4
76	Acquiring snapshots of the orientation of transmembrane protein domains using a hybrid FRET pair. <i>FEBS Letters</i> , 2015, 589, 885-889.	2.8	4
77	Solvent saturation transfer to proteins (SSTP) for structural and functional characterization of proteins. <i>Journal of Biomolecular NMR</i> , 2018, 70, 11-20.	2.8	3
78	Squeezing lipids: NMR characterization of lipoprotein particles under pressure. <i>Chemistry and Physics of Lipids</i> , 2020, 228, 104874.	3.2	3
79	Decoding the components of dynamics in three-domain proteins. <i>Journal of Computational Chemistry</i> , 2014, 35, 518-525.	3.3	2
80	Inducible fold-switching as a mechanism to fibrillate proapoptotic BCL-2 proteins. <i>Biopolymers</i> , 2021, 112, e23424.	2.4	2
81	Structural Basis for the Interaction of Fibrin with the Very Low-Density Lipoprotein Receptor Revealed by NMR and Site-Directed Mutagenesis. <i>Biochemistry</i> , 2021, 60, 2537-2548.	2.5	2
82	Exploiting image registration for automated resonance assignment in NMR. <i>Journal of Biomolecular NMR</i> , 2015, 62, 143-156.	2.8	1
83	Verbesserung der Dispersion der chemischen Verschiebungen von unstrukturierten Proteinen durch einen kovalent gebundenen Lanthanoidkomplex. <i>Angewandte Chemie</i> , 2016, 128, 15069-15073.	2.0	1
84	Incorporation of residual chemical shift anisotropy into the treatment of ¹⁵ N pseudocontact shifts for structural refinement. <i>Journal of Magnetic Resonance</i> , 2022, 340, 107213.	2.1	1
85	The Kindlin Outside Connection. <i>Structure</i> , 2019, 27, 1615-1616.	3.3	0
86	NMR Analysis of Apo Glutamine-Binding Protein Exposes Challenges in the Study of Interdomain Dynamics. <i>Angewandte Chemie</i> , 2019, 131, 17055-17058.	2.0	0
87	Simultaneous measurement of ¹ H/ ¹⁵ N-R ₂ ρs for rapid acquisition of backbone and sidechain paramagnetic relaxation enhancements (PREs) in proteins. <i>Journal of Biomolecular NMR</i> , 2021, 75, 109-118.	2.8	0
88	Residual Dipolar Coupling for Conformational and Dynamic Studies. , 2017, , 1-16.		0
89	Residual Dipolar Coupling for Conformational and Dynamic Studies. , 2018, , 419-434.		0