

H Jane Dyson

List of Publications by Year in descending order

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291
papers

40,446
citations

2669

95
h-index

2617

194
g-index

402
all docs

402
docs citations

402
times ranked

27477
citing authors

#	ARTICLE	IF	CITATIONS
1	Multivalency enables unidirectional switch-like competition between intrinsically disordered proteins. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022, 119, .	3.3	22
2	Interactions of a Long Noncoding RNA with Domains of NF- κ B and I κ B β : Implications for the Inhibition of Non-Signal-Related Phosphorylation. <i>Biochemistry</i> , 2022, 61, 367-376.	1.2	4
3	A transthyretin monomer intermediate undergoes local unfolding and transient interaction with oligomers in a kinetically concerted aggregation pathway. <i>Journal of Biological Chemistry</i> , 2022, 298, 102162.	1.6	5
4	Modeling of Hidden Structures Using Sparse Chemical Shift Data from NMR Relaxation Dispersion. <i>Biophysical Journal</i> , 2021, 120, 296-305.	0.2	4
5	Backbone and side-chain chemical shift assignments of p50 subunit of NF- κ B transcription factor. <i>Biomolecular NMR Assignments</i> , 2021, 15, 29-33.	0.4	2
6	Using NMR to identify binding regions for N and C-terminal Hsp90 inhibitors using Hsp90 domains. <i>RSC Medicinal Chemistry</i> , 2021, 12, 410-415.	1.7	4
7	Thermodynamic Stability and Aggregation Kinetics of EF Helix and EF Loop Variants of Transthyretin. <i>Biochemistry</i> , 2021, 60, 756-764.	1.2	14
8	Early Strides in NMR Dynamics Measurements. <i>Biochemistry</i> , 2021, 60, 3452-3454.	1.2	0
9	Diversity at BJ: The editors, the reviewers, the authors. <i>Biophysical Journal</i> , 2021, 120, E1-E2.	0.2	0
10	Role of Active Site Loop Dynamics in Mediating Ligand Release from <i>E. coli</i> Dihydrofolate Reductase. <i>Biochemistry</i> , 2021, 60, 2663-2671.	1.2	4
11	The molecular basis of allostery in a facilitated dissociation process. <i>Structure</i> , 2021, 29, 1327-1338.e5.	1.6	6
12	NMR illuminates intrinsic disorder. <i>Current Opinion in Structural Biology</i> , 2021, 70, 44-52.	2.6	60
13	More pandemic reflections. <i>Biophysical Journal</i> , 2021, 120, E1-E2.	0.2	0
14	A phosphorylation-dependent switch in the disordered p53 transactivation domain regulates DNA binding. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	3.3	33
15	Characterization of the High-Affinity Fuzzy Complex between the Disordered Domain of the E7 Oncoprotein from High-Risk HPV and the TAZ2 Domain of CBP. <i>Biochemistry</i> , 2021, 60, 3887-3898.	1.2	9
16	Reflections on the Pandemic. <i>Biophysical Journal</i> , 2020, 119, E1.	0.2	0
17	RNA Binding by the KTS Splice Variants of Wilms's Tumor Suppressor Protein WT1. <i>Biochemistry</i> , 2020, 59, 3889-3901.	1.2	4
18	A Conformational Switch in the Zinc Finger Protein Kaiso Mediates Differential Readout of Specific and Methylated DNA Sequences. <i>Biochemistry</i> , 2020, 59, 1909-1926.	1.2	7

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19	Management of Hsp90-Dependent Protein Folding by Small Molecules Targeting the Aha1 Co-Chaperone. <i>Cell Chemical Biology</i> , 2020, 27, 292-305.e6.	2.5	13
20	Editorsâ€™ Note. <i>Biophysical Journal</i> , 2020, 119, E1.	0.2	0
21	Comparison of backbone dynamics of the p50 dimerization domain of NF κ B in the homodimeric transcription factor NF κ B1 and in its heterodimeric complex with RelA (p65). <i>Protein Science</i> , 2019, 28, 2064-2072.	3.1	7
22	Perspective: the essential role of NMR in the discovery and characterization of intrinsically disordered proteins. <i>Journal of Biomolecular NMR</i> , 2019, 73, 651-659.	1.6	48
23	Aggregation of zincâ€free p53 is inhibited by Hsp90 but not other chaperones. <i>Protein Science</i> , 2019, 28, 2020-2023.	3.1	7
24	Economics and Politics of Publishing in Our Mission-Driven Society. <i>Biophysical Journal</i> , 2019, 116, E1-E2.	0.2	0
25	A Journal for All Biophysicists. <i>Biophysical Journal</i> , 2019, 116, E1.	0.2	0
26	Role of Backbone Dynamics in Modulating the Interactions of Disordered Ligands with the TAZ1 Domain of the CREB-Binding Protein. <i>Biochemistry</i> , 2019, 58, 1354-1362.	1.2	33
27	A Dynamic Switch in Inactive p38 β Leads to an Excited State on the Pathway to an Active Kinase. <i>Biochemistry</i> , 2019, 58, 5160-5172.	1.2	7
28	Expanding the Paradigm: Intrinsically Disordered Proteins and Allosteric Regulation. <i>Journal of Molecular Biology</i> , 2018, 430, 2309-2320.	2.0	105
29	Slow Dynamics of Tryptophanâ€Water Networks in Proteins. <i>Journal of the American Chemical Society</i> , 2018, 140, 675-682.	6.6	26
30	Characterization of an Hsp90-Independent Interaction between Co-Chaperone p23 and Transcription Factor p53. <i>Biochemistry</i> , 2018, 57, 935-944.	1.2	13
31	Is the BJ Review Process Gender-Biased?. <i>Biophysical Journal</i> , 2018, 114, E1.	0.2	0
32	CHâ€O Hydrogen Bonds Mediate Highly Specific Recognition of Methylated CpG Sites by the Zinc Finger Protein Kaiso. <i>Biochemistry</i> , 2018, 57, 2109-2120.	1.2	19
33	Long-range regulation of p53 DNA binding by its intrinsically disordered N-terminal transactivation domain. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018, 115, E11302-E11310.	3.3	93
34	Mispacking of the Phe87 Side Chain Reduces the Kinetic Stability of Human Transthyretin. <i>Biochemistry</i> , 2018, 57, 6919-6922.	1.2	8
35	Structural Basis for Graded Inhibition of CREB:DNA Interactions by Multisite Phosphorylation. <i>Biochemistry</i> , 2018, 57, 6964-6972.	1.2	7
36	Structural basis for cooperative regulation of KIX-mediated transcription pathways by the HTLV-1 HBZ activation domain. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018, 115, 10040-10045.	3.3	18

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37	How Do Intrinsically Disordered Viral Proteins Hijack the Cell?. <i>Biochemistry</i> , 2018, 57, 4045-4046.	1.2	22
38	NMR Measurements Reveal the Structural Basis of Transthyretin Destabilization by Pathogenic Mutations. <i>Biochemistry</i> , 2018, 57, 4421-4430.	1.2	30
39	Kinetic analysis of the multistep aggregation pathway of human transthyretin. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018, 115, E6201-E6208.	3.3	29
40	Hypersensitive termination of the hypoxic response by a disordered protein switch. <i>Nature</i> , 2017, 543, 447-451.	13.7	140
41	Functional importance of stripping in NF- κ B signaling revealed by a stripping-impaired $\hat{\text{I}}^{\text{p}}\hat{\text{B}}\hat{\text{I}}^{\text{s}}$ mutant. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, 1916-1921.	3.3	32
42	Role of the CBP catalytic core in intramolecular SUMOylation and control of histone H3 acetylation. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E5335-E5342.	3.3	56
43	How Does Your Protein Fold? Elucidating the Apomyoglobin Folding Pathway. <i>Accounts of Chemical Research</i> , 2017, 50, 105-111.	7.6	44
44	Fluorotryptophan Incorporation Modulates the Structure and Stability of Transthyretin in a Site-Specific Manner. <i>Biochemistry</i> , 2017, 56, 5570-5581.	1.2	20
45	Defining the Structural Basis for Allosteric Product Release from <i>E. coli</i> Dihydrofolate Reductase Using NMR Relaxation Dispersion. <i>Journal of the American Chemical Society</i> , 2017, 139, 11233-11240.	6.6	27
46	Structural Basis for Interaction of the Tandem Zinc Finger Domains of Human Muscleblind with Cognate RNA from Human Cardiac Troponin T. <i>Biochemistry</i> , 2017, 56, 4154-4168.	1.2	27
47	Greetings from Your New Editor-in-Chief. <i>Biophysical Journal</i> , 2017, 113, E1.	0.2	0
48	Finding Our Way in the Dark Proteome. <i>Journal of the American Chemical Society</i> , 2016, 138, 9730-9742.	6.6	111
49	Structural characterization of the ternary complex that mediates termination of NF- $\hat{\text{I}}^{\text{p}}\hat{\text{B}}\hat{\text{I}}^{\text{s}}$ signaling by $\hat{\text{I}}^{\text{p}}\hat{\text{B}}\hat{\text{I}}^{\text{s}}$. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, 6212-6217.	3.3	8
50	Mapping the interactions of adenoviral E1A proteins with the p160 nuclear receptor coactivator binding domain of CBP. <i>Protein Science</i> , 2016, 25, 2256-2267.	3.1	18
51	$\langle \text{scp} \rangle$ NMR $\langle \text{scp} \rangle$ characterization of a 72 k $\langle \text{scp} \rangle$ D $\langle \text{scp} \rangle$ transcription factor using differential isotopic labeling. <i>Protein Science</i> , 2016, 25, 597-604.	3.1	8
52	The Dependence of Carbohydrate- $\hat{\text{A}}^{\text{r}}$ Aromatic Interaction Strengths on the Structure of the Carbohydrate. <i>Journal of the American Chemical Society</i> , 2016, 138, 7636-7648.	6.6	44
53	Recognition of the disordered p53 transactivation domain by the transcriptional adapter zinc finger domains of CREB-binding protein. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, E1853-62.	3.3	94
54	Role of Intrinsic Protein Disorder in the Function and Interactions of the Transcriptional Coactivators CREB-binding Protein (CBP) and p300. <i>Journal of Biological Chemistry</i> , 2016, 291, 6714-6722.	1.6	251

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55	Making Sense of Intrinsically Disordered Proteins. <i>Biophysical Journal</i> , 2016, 110, 1013-1016.	0.2	81
56	Classic Analysis of Biopolymer Dynamics Is Model Free. <i>Biophysical Journal</i> , 2016, 110, 3-6.	0.2	4
57	Functional advantages of dynamic protein disorder. <i>FEBS Letters</i> , 2015, 589, 2433-2440.	1.3	162
58	Cofactor-Mediated Conformational Dynamics Promote Product Release From <i>Escherichia coli</i> Dihydrofolate Reductase via an Allosteric Pathway. <i>Journal of the American Chemical Society</i> , 2015, 137, 9459-9468.	6.6	45
59	Conformational propensities of intrinsically disordered proteins influence the mechanism of binding and folding. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015, 112, 9614-9619.	3.3	222
60	Biomolecular Systems Interactions, Dynamics, and Allostery: Reflections and New Directions. <i>Biophysical Journal</i> , 2015, 109, E01-E02.	0.2	0
61	Intrinsically disordered proteins in cellular signalling and regulation. <i>Nature Reviews Molecular Cell Biology</i> , 2015, 16, 18-29.	16.1	1,849
62	The High-Risk HPV16 E7 Oncoprotein Mediates Interaction between the Transcriptional Coactivator CBP and the Retinoblastoma Protein pRb. <i>Journal of Molecular Biology</i> , 2014, 426, 4030-4048.	2.0	61
63	Side Chain Conformational Averaging in Human Dihydrofolate Reductase. <i>Biochemistry</i> , 2014, 53, 1134-1145.	1.2	8
64	Probing the Non-Native H Helix Translocation in Apomyoglobin Folding Intermediates. <i>Biochemistry</i> , 2014, 53, 3767-3780.	1.2	16
65	Structural Characterization of Interactions between the Double-Stranded RNA-Binding Zinc Finger Protein JAZ and Nucleic Acids. <i>Biochemistry</i> , 2014, 53, 1495-1510.	1.2	20
66	The CH2 domain of CBP/p300 is a novel zinc finger. <i>FEBS Letters</i> , 2013, 587, 2506-2511.	1.3	12
67	Divergent evolution of protein conformational dynamics in dihydrofolate reductase. <i>Nature Structural and Molecular Biology</i> , 2013, 20, 1243-1249.	3.6	153
68	Localized Structural Fluctuations Promote Amyloidogenic Conformations in Transthyretin. <i>Journal of Molecular Biology</i> , 2013, 425, 977-988.	2.0	65
69	Long-Range Effects and Functional Consequences of Stabilizing Mutations in the Ankyrin Repeat Domain of I β B \pm . <i>Journal of Molecular Biology</i> , 2013, 425, 902-913.	2.0	10
70	A Distal Mutation Perturbs Dynamic Amino Acid Networks in Dihydrofolate Reductase. <i>Biochemistry</i> , 2013, 52, 4605-4619.	1.2	77
71	Structural and Energetic Basis of Carbohydrate-Aromatic Packing Interactions in Proteins. <i>Journal of the American Chemical Society</i> , 2013, 135, 9877-9884.	6.6	85
72	Side-Chain Conformational Heterogeneity of Intermediates in the <i>Escherichia coli</i> Dihydrofolate Reductase Catalytic Cycle. <i>Biochemistry</i> , 2013, 52, 3464-3477.	1.2	16

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73	What's in a name? Why these proteins are intrinsically disordered. <i>Intrinsically Disordered Proteins</i> , 2013, 1, e24157.	1.9	226
74	Identification of Cys255 in HIF1 α as a novel site for development of covalent inhibitors of HIF1 α /ARNT PasB domain protein-protein interaction. <i>Protein Science</i> , 2012, 21, 1885-1896.	3.1	64
75	CheShift-2 resolves a local inconsistency between two X-ray crystal structures. <i>Journal of Biomolecular NMR</i> , 2012, 54, 193-198.	1.6	4
76	Homodimerization of the PAS-B Domains of Hypoxia-Inducible Factors. <i>Journal of Physical Chemistry B</i> , 2012, 116, 6960-6965.	1.2	5
77	Molecular basis for recognition of methylated and specific DNA sequences by the zinc finger protein Kaiso. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2012, 109, 15229-15234.	3.3	101
78	Roles of intrinsic disorder in protein-nucleic acid interactions. <i>Molecular BioSystems</i> , 2012, 8, 97-104.	2.9	76
79	Role of disorder in NF κ B interaction. <i>IUBMB Life</i> , 2012, 64, 499-505.	1.5	41
80	Kaiso uses all three zinc fingers and adjacent sequence motifs for high affinity binding to sequence-specific and methylated CpG DNA targets. <i>FEBS Letters</i> , 2012, 586, 734-739.	1.3	17
81	Structural Basis for Cooperative Transcription Factor Binding to the CBP Coactivator. <i>FASEB Journal</i> , 2012, 26, lb266.	0.2	0
82	A Dynamic Knockout Reveals That Conformational Fluctuations Influence the Chemical Step of Enzyme Catalysis. <i>Science</i> , 2011, 332, 234-238.	6.0	414
83	Expanding the proteome: disordered and alternatively folded proteins. <i>Quarterly Reviews of Biophysics</i> , 2011, 44, 467-518.	2.4	150
84	The RelA Nuclear Localization Signal Folds upon Binding to I κ B α . <i>Journal of Molecular Biology</i> , 2011, 405, 754-764.	2.0	29
85	Consequences of Stabilizing the Natively Disordered F Helix for the Folding Pathway of Apomyoglobin. <i>Journal of Molecular Biology</i> , 2011, 411, 248-263.	2.0	16
86	Dynamic Interaction of Hsp90 with Its Client Protein p53. <i>Journal of Molecular Biology</i> , 2011, 411, 158-173.	2.0	72
87	The client protein p53 adopts a molten globule-like state in the presence of Hsp90. <i>Nature Structural and Molecular Biology</i> , 2011, 18, 537-541.	3.6	121
88	Detection of a ternary complex of NF κ B and I κ B α with DNA provides insights into how I κ B α removes NF κ B from transcription sites. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2011, 108, 1367-1372.	3.3	26
89	Leu628 of the KIX domain of CBP is a key residue for the interaction with the MLL transactivation domain. <i>FEBS Letters</i> , 2010, 584, 4500-4504.	1.3	32
90	3P040 Mapping the Interactions of the Intrinsically Disordered p53 Transactivation Subdomains with the TAZ2 Domain of CBP by NMR (Protein: Structure & Function, The 48th Annual Meeting of the) Tj ETQq0 0 0 rgBTj/Overlock 10 Tf 50 5		

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91	Millisecond timescale fluctuations in dihydrofolate reductase are exquisitely sensitive to the bound ligands. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010, 107, 1373-1378.	3.3	133
92	Structure of the p53 Transactivation Domain in Complex with the Nuclear Receptor Coactivator Binding Domain of CREB Binding Protein. <i>Biochemistry</i> , 2010, 49, 9964-9971.	1.2	162
93	Energetic Frustration of Apomyoglobin Folding: Role of the B Helix. <i>Journal of Molecular Biology</i> , 2010, 396, 1319-1328.	2.0	17
94	Cooperative regulation of p53 by modulation of ternary complex formation with CBP/p300 and HDM2. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009, 106, 6591-6596.	3.3	197
95	Structural basis for subversion of cellular control mechanisms by the adenoviral E1A oncoprotein. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009, 106, 13260-13265.	3.3	119
96	Evaluating β -turn mimics as β -sheet folding nucleators. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009, 106, 11067-11072.	3.3	97
97	Linking folding and binding. <i>Current Opinion in Structural Biology</i> , 2009, 19, 31-38.	2.6	932
98	Diagnostic chemical shift markers for loop conformation and substrate and cofactor binding in dihydrofolate reductase complexes. <i>Protein Science</i> , 2009, 12, 2230-2238.	3.1	38
99	Mapping the Interactions of the p53 Transactivation Domain with the KIX Domain of CBP. <i>Biochemistry</i> , 2009, 48, 2115-2124.	1.2	109
100	Prion Proteins with Pathogenic and Protective Mutations Show Similar Structure and Dynamics. <i>Biochemistry</i> , 2009, 48, 8120-8128.	1.2	53
101	Interaction of the β -terminal PEST Sequence with NF- κ B: Insights into the Inhibition of NF- κ B DNA Binding by β . <i>Journal of Molecular Biology</i> , 2009, 388, 824-838.	2.0	28
102	Functional Dynamics of the Folded Ankyrin Repeats of β Revealed by Nuclear Magnetic Resonance. <i>Biochemistry</i> , 2009, 48, 8023-8031.	1.2	22
103	Structural basis for recruitment of CBP/p300 coactivators by STAT1 and STAT2 transactivation domains. <i>EMBO Journal</i> , 2009, 28, 948-958.	3.5	147
104	Prediction of the Rotational Tumbling Time for Proteins with Disordered Segments. <i>Journal of the American Chemical Society</i> , 2009, 131, 6814-6821.	6.6	48
105	Structure discrimination for the C-terminal domain of Escherichia coli trigger factor in solution. <i>Journal of Biomolecular NMR</i> , 2008, 40, 23-30.	1.6	17
106	Structural characterization of partially folded intermediates of apomyoglobin H64F. <i>Protein Science</i> , 2008, 17, 313-321.	3.1	16
107	Hydrogen-deuterium exchange strategy for delineation of contact sites in protein complexes. <i>FEBS Letters</i> , 2008, 582, 1495-1500.	1.3	16
108	The Kinetic and Equilibrium Molten Globule Intermediates of Apoleghemoglobin Differ in Structure. <i>Journal of Molecular Biology</i> , 2008, 378, 715-725.	2.0	26

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109	Transfer of Flexibility between Ankyrin Repeats in Î±B1± upon Formation of the NF-Î±B Complex. <i>Journal of Molecular Biology</i> , 2008, 380, 917-931.	2.0	61
110	Hierarchical folding mechanism of apomyoglobin revealed by ultra-fast H/D exchange coupled with 2D NMR. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 13859-13864.	3.3	89
111	Amylin Proprotein Processing Generates Progressively More Amyloidogenic Peptides that Initially Sample the Helical State. <i>Biochemistry</i> , 2008, 47, 9900-9910.	1.2	132
112	NMR Relaxation Study of the Complex Formed Between CBP and the Activation Domain of the Nuclear Hormone Receptor Coactivator ACTR^{â€‹}. <i>Biochemistry</i> , 2008, 47, 1299-1308.	1.2	86
113	The Intrinsically Disordered RNR Inhibitor Sml1 Is a Dynamic Dimer. <i>Biochemistry</i> , 2008, 47, 13428-13437.	1.2	53
114	Conformational Relaxation following Hydride Transfer Plays a Limiting Role in Dihydrofolate Reductase Catalysisâ€‹. <i>Biochemistry</i> , 2008, 47, 9227-9233.	1.2	53
115	Modeling transient collapsed states of an unfolded protein to provide insights into early folding events. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 6278-6283.	3.3	105
116	Embryonic Neural Inducing Factor Churchill Is not a DNA-binding Zinc Finger Protein: Solution Structure Reveals a Solvent-exposed Î²-Sheet and Zinc Binuclear Cluster. <i>Journal of Molecular Biology</i> , 2007, 371, 1274-1289.	2.0	21
117	Structure of the Wilms Tumor Suppressor Protein Zinc Finger Domain Bound to DNA. <i>Journal of Molecular Biology</i> , 2007, 372, 1227-1245.	2.0	91
118	Tailoring Relaxation Dispersion Experiments for Fast-Associating Protein Complexes. <i>Journal of the American Chemical Society</i> , 2007, 129, 13406-13407.	6.6	52
119	NMR detection of adventitious xylose binding to the quorum-sensing protein SdiA of <i>Escherichia coli</i> . <i>Bioorganic and Medicinal Chemistry Letters</i> , 2007, 17, 6202-6205.	1.0	7
120	Mechanism of coupled folding and binding of an intrinsically disordered protein. <i>Nature</i> , 2007, 447, 1021-1025.	13.7	984
121	Dynamics of IÎ±Ba Probed by NMR. <i>FASEB Journal</i> , 2007, 21, A655.	0.2	0
122	The Dynamic Energy Landscape of Dihydrofolate Reductase Catalysis. <i>Science</i> , 2006, 313, 1638-1642.	6.0	877
123	An NMR Perspective on Enzyme Dynamics. <i>Chemical Reviews</i> , 2006, 106, 3055-3079.	23.0	424
124	NMR Solution Structure of the Peptide Fragment 1âˆ’30, Derived from Unprocessed Mouse Doppel Protein, in DHPC Micellesâ€‹. <i>Biochemistry</i> , 2006, 45, 159-166.	1.2	19
125	The Reduced, Denatured Somatomedin B Domain of Vitronectin Refolds into a Stable, Biologically Active Molecule. <i>Biochemistry</i> , 2006, 45, 3297-3306.	1.2	11
126	Structural Basis for Cooperative Transcription Factor Binding to the CBP Coactivator. <i>Journal of Molecular Biology</i> , 2006, 355, 1005-1013.	2.0	166

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127	Structure of the Escherichia coli Quorum Sensing Protein SdiA: Activation of the Folding Switch by Acyl Homoserine Lactones. <i>Journal of Molecular Biology</i> , 2006, 355, 262-273.	2.0	162
128	Identification of Native and Non-native Structure in Kinetic Folding Intermediates of Apomyoglobin. <i>Journal of Molecular Biology</i> , 2006, 355, 139-156.	2.0	112
129	Induced Fit and "Lock and Key" Recognition of 5S RNA by Zinc Fingers of Transcription Factor IIIA. <i>Journal of Molecular Biology</i> , 2006, 357, 275-291.	2.0	72
130	Solution Structure of the Hdm2 C2H2C4 RING, a Domain Critical for Ubiquitination of p53. <i>Journal of Molecular Biology</i> , 2006, 363, 433-450.	2.0	120
131	According to current textbooks, a well-defined three-dimensional structure is a prerequisite for the function of a protein. Is this correct?. <i>IUBMB Life</i> , 2006, 58, 107-109.	1.5	20
132	The role of hydrophobic interactions in initiation and propagation of protein folding. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2006, 103, 13057-13061.	3.3	266
133	Localization of Sites of Interaction between p23 and Hsp90 in Solution. <i>Journal of Biological Chemistry</i> , 2006, 281, 14457-14464.	1.6	58
134	Structure and Function of the CBP/p300 TAZ Domains. , 2005, , 114-120.		4
135	Intrinsically unstructured proteins and their functions. <i>Nature Reviews Molecular Cell Biology</i> , 2005, 6, 197-208.	16.1	3,403
136	Generation of native-like protein structures from limited NMR data, modern force fields and advanced conformational sampling. <i>Journal of Biomolecular NMR</i> , 2005, 31, 59-64.	1.6	36
137	Letter to the Editor: Backbone and side chain 1H, 13C and 15N assignments for Escherichia coli SdiA1-171, the autoinducer-binding domain of a quorum sensing protein. <i>Journal of Biomolecular NMR</i> , 2005, 31, 373-374.	1.6	7
138	Defining the role of active-site loop fluctuations in dihydrofolate reductase catalysis. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 5032-5037.	3.3	152
139	Enhanced picture of protein-folding intermediates using organic solvents in H/D exchange and quench-flow experiments. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 4765-4770.	3.3	62
140	CBP/p300 TAZ1 Domain Forms a Structured Scaffold for Ligand Binding,. <i>Biochemistry</i> , 2005, 44, 490-497.	1.2	76
141	Solution Structure of the N-terminal Zinc Fingers of the Xenopus laevis double-stranded RNA-binding Protein ZFa. <i>Journal of Molecular Biology</i> , 2005, 351, 718-730.	2.0	18
142	Sequence Determinants of a Protein Folding Pathway. <i>Journal of Molecular Biology</i> , 2005, 351, 383-392.	2.0	54
143	Elucidation of the Protein Folding Landscape by NMR. <i>Methods in Enzymology</i> , 2005, 394, 299-321.	0.4	90
144	Interaction of the TAZ1 Domain of the CREB-Binding Protein with the Activation Domain of CITED2. <i>Journal of Biological Chemistry</i> , 2004, 279, 3042-3049.	1.6	97

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145	Activation of the Redox-regulated Chaperone Hsp33 by Domain Unfolding. <i>Journal of Biological Chemistry</i> , 2004, 279, 20529-20538.	1.6	100
146	Recognition of the mRNA AU-rich element by the zinc finger domain of TIS11d. <i>Nature Structural and Molecular Biology</i> , 2004, 11, 257-264.	3.6	320
147	Unfolded Proteins and Protein Folding Studied by NMR. <i>ChemInform</i> , 2004, 35, no.	0.1	1
148	Disulfide Bonding Arrangements in Active Forms of the Somatomedin B Domain of Human Vitronectin. <i>Biochemistry</i> , 2004, 43, 6519-6534.	1.2	37
149	The LEF-1 High-Mobility Group Domain Undergoes a Disorder-to-Order Transition upon Formation of a Complex with Cognate DNA. <i>Biochemistry</i> , 2004, 43, 8725-8734.	1.2	62
150	Effect of Cofactor Binding and Loop Conformation on Side Chain Methyl Dynamics in Dihydrofolate Reductase. <i>Biochemistry</i> , 2004, 43, 374-383.	1.2	73
151	Conformational Changes in the Active Site Loops of Dihydrofolate Reductase during the Catalytic Cycle. <i>Biochemistry</i> , 2004, 43, 16046-16055.	1.2	119
152	Unfolded Proteins and Protein Folding Studied by NMR. <i>Chemical Reviews</i> , 2004, 104, 3607-3622.	23.0	596
153	Packing, specificity, and mutability at the binding interface between the p160 coactivator and CREB-binding protein. <i>Protein Science</i> , 2004, 13, 203-210.	3.1	66
154	Introduction: Biological Nuclear Magnetic Resonance. <i>Chemical Reviews</i> , 2004, 104, 3517-3518.	23.0	2
155	Solution Structure of the KIX Domain of CBP Bound to the Transactivation Domain of c-Myb. <i>Journal of Molecular Biology</i> , 2004, 337, 521-534.	2.0	181
156	Structural Characterization of Unfolded States of Apomyoglobin using Residual Dipolar Couplings. <i>Journal of Molecular Biology</i> , 2004, 340, 1131-1142.	2.0	165
157	The Zinc-dependent Redox Switch Domain of the Chaperone Hsp33 has a Novel Fold. <i>Journal of Molecular Biology</i> , 2004, 341, 893-899.	2.0	52
158	ZZ Domain of CBP: an Unusual Zinc Finger Fold in a Protein Interaction Module. <i>Journal of Molecular Biology</i> , 2004, 343, 1081-1093.	2.0	81
159	Structure, Dynamics, and Catalytic Function of Dihydrofolate Reductase. <i>Annual Review of Biophysics and Biomolecular Structure</i> , 2004, 33, 119-140.	18.3	444
160	The CBP/p300 TAZ1 domain in its native state is not a binding partner of MDM2. <i>Biochemical Journal</i> , 2004, 381, 685-691.	1.7	41
161	Changes in structure and dynamics of the Fv fragment of a catalytic antibody upon binding of inhibitor. <i>Protein Science</i> , 2003, 12, 1386-1394.	3.1	14
162	Role of a solvent-exposed tryptophan in the recognition and binding of antibiotic substrates for a metallo- β -lactamase. <i>Protein Science</i> , 2003, 12, 1368-1375.	3.1	56

#	ARTICLE	IF	CITATIONS
163	Structure of the Nuclear Factor ALY: Insights into Post-Transcriptional Regulatory and mRNA Nuclear Export Processes. <i>Biochemistry</i> , 2003, 42, 7348-7357.	1.2	20
164	Role of the B Helix in Early Folding Events in Apomyoglobin: Evidence from Site-directed Mutagenesis for Native-like Long Range Interactions. <i>Journal of Molecular Biology</i> , 2003, 334, 293-307.	2.0	51
165	Monomeric Complex of Human Orphan Estrogen Related Receptor-2 with DNA: A Pseudo-dimer Interface Mediates Extended Half-site Recognition. <i>Journal of Molecular Biology</i> , 2003, 327, 819-832.	2.0	97
166	Folding of a β^2 -sheet Protein Monitored by Real-time NMR Spectroscopy. <i>Journal of Molecular Biology</i> , 2003, 328, 1161-1171.	2.0	29
167	Roles of Phosphorylation and Helix Propensity in the Binding of the KIX Domain of CREB-binding Protein by Constitutive (c-Myb) and Inducible (CREB) Activators. <i>Journal of Biological Chemistry</i> , 2002, 277, 42241-42248.	1.6	134
168	Structural basis for Hif-1 α /CBP recognition in the cellular hypoxic response. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002, 99, 5271-5276.	3.3	376
169	Molecular Hinges in Protein Folding: The Urea-Denatured State of Apomyoglobin. <i>Biochemistry</i> , 2002, 41, 12681-12686.	1.2	130
170	Insights into the structure and dynamics of unfolded proteins from nuclear magnetic resonance. <i>Advances in Protein Chemistry</i> , 2002, 62, 311-340.	4.4	208
171	The Apomyoglobin Folding Pathway Revisited: Structural Heterogeneity in the Kinetic Burst Phase Intermediate. <i>Journal of Molecular Biology</i> , 2002, 322, 483-489.	2.0	89
172	Cooperativity in Transcription Factor Binding to the Coactivator CREB-binding Protein (CBP). <i>Journal of Biological Chemistry</i> , 2002, 277, 43168-43174.	1.6	166
173	Mapping Long-range Contacts in a Highly Unfolded Protein. <i>Journal of Molecular Biology</i> , 2002, 322, 655-662.	2.0	140
174	Coupling of folding and binding for unstructured proteins. <i>Current Opinion in Structural Biology</i> , 2002, 12, 54-60.	2.6	1,223
175	Mutual synergistic folding in recruitment of CBP/p300 by p160 nuclear receptor coactivators. <i>Nature</i> , 2002, 415, 549-553.	13.7	423
176	Comparison of protein solution structures refined by molecular dynamics simulation in vacuum, with a generalized Born model, and with explicit water. <i>Journal of Biomolecular NMR</i> , 2002, 22, 317-331.	1.6	112
177	Assignment of a 15 kDa protein complex formed between the p160 coactivator ACTR and CREB binding protein. <i>Journal of Biomolecular NMR</i> , 2002, 22, 377-378.	1.6	6
178	Structural and dynamic characterization of an unfolded state of poplar apo-plastocyanin formed under nondenaturing conditions. <i>Protein Science</i> , 2001, 10, 1056-1066.	3.1	79
179	NMR Structural and Dynamic Characterization of the Acid-Unfolded State of Apomyoglobin Provides Insights into the Early Events in Protein Folding. <i>Biochemistry</i> , 2001, 40, 3561-3571.	1.2	212
180	Backbone Dynamics in Dihydrofolate Reductase Complexes: Role of Loop Flexibility in the Catalytic Mechanism. <i>Biochemistry</i> , 2001, 40, 9846-9859.	1.2	246

#	ARTICLE	IF	CITATIONS
181	Nuclear Magnetic Resonance Methods for Elucidation of Structure and Dynamics in Disordered States. <i>Methods in Enzymology</i> , 2001, 339, 258-270.	0.4	148
182	Sequence-Dependent Correction of Random Coil NMR Chemical Shifts. <i>Journal of the American Chemical Society</i> , 2001, 123, 2970-2978.	6.6	562
183	Solution structure of Escherichia coli glutaredoxin-2 shows similarity to mammalian glutathione-S-transferases. <i>Journal of Molecular Biology</i> , 2001, 310, 907-918.	2.0	71
184	Solution structure of the third immunoglobulin domain of the neural cell adhesion molecule N-CAM: can solution studies define the mechanism of homophilic binding?. <i>Journal of Molecular Biology</i> , 2001, 311, 161-172.	2.0	25
185	Conformational and Dynamic Characterization of the Molten Globule State of an Apomyoglobin Mutant with an Altered Folding Pathway. <i>Biochemistry</i> , 2001, 40, 14459-14467.	1.2	45
186	Local Structural Plasticity of the Prion Protein. Analysis of NMR Relaxation Dynamics. <i>Biochemistry</i> , 2001, 40, 2743-2753.	1.2	171
187	SANE (Structure Assisted NOE Evaluation): an automated model-based approach for NOE assignment. <i>Journal of Biomolecular NMR</i> , 2001, 19, 321-329.	1.6	113
188	Potential bias in NMR relaxation data introduced by peak intensity analysis and curve fitting methods. <i>Journal of Biomolecular NMR</i> , 2001, 21, 1-9.	1.6	44
189	Genomic-scale comparison of sequence- and structure-based methods of function prediction: Does structure provide additional insight?. <i>Protein Science</i> , 2001, 10, 1005-1014.	3.1	68
190	Two different neurodegenerative diseases caused by proteins with similar structures. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2001, 98, 2352-2357.	3.3	147
191	Conservation of folding pathways in evolutionarily distant globin sequences. <i>Nature Structural Biology</i> , 2000, 7, 679-686.	9.7	95
192	Backbone H(N), N, C α , C β and C γ assignments of the 19 kDa DHFR/NADPH complex at 9 degrees C and pH 7.6. <i>Journal of Biomolecular NMR</i> , 2000, 16, 349-350.	1.6	3
193	Assignment of ^1H , ^{13}C and ^{15}N resonances of the I-domain of human leukocyte function associated antigen-1. <i>Journal of Biomolecular NMR</i> , 2000, 16, 271-272.	1.6	7
194	Random coil chemical shifts in acidic 8 M urea: implementation of random coil shift data in NMRView. <i>Journal of Biomolecular NMR</i> , 2000, 18, 43-48.	1.6	272
195	Molecular basis for modulation of biological function by alternate splicing of the Wilms' tumor suppressor protein. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2000, 97, 11932-11935.	3.3	115
196	DNA-induced α -helix capping in conserved linker sequences is a determinant of binding affinity in Cys2-His2 zinc fingers. <i>Journal of Molecular Biology</i> , 2000, 295, 719-727.	2.0	137
197	NMR solution structure of the inserted domain of human leukocyte function associated antigen-1. <i>Journal of Molecular Biology</i> , 2000, 295, 1251-1264.	2.0	74
198	Solution Structure of the Cysteine-rich Domain of the Escherichia coli Chaperone Protein DnaJ. <i>Journal of Molecular Biology</i> , 2000, 300, 805-818.	2.0	121

#	ARTICLE	IF	CITATIONS
199	Efficient Inhibition of Escherichia Coli RNA Polymerase by the Bacteriophage T4 AsiA Protein Requires that AsiA Binds First to Free λ 70. <i>Journal of Molecular Biology</i> , 2000, 304, 731-739.	2.0	55
200	Solution structure of the TAZ2 (CH3) domain of the transcriptional adaptor protein CBP. <i>Journal of Molecular Biology</i> , 2000, 303, 243-253.	2.0	121
201	Solution structure and acetyl-lysine binding activity of the GCN5 bromodomain. <i>Journal of Molecular Biology</i> , 2000, 304, 355-370.	2.0	141
202	Identification of the regions involved in DNA binding by the mouse PEBP2 β protein. <i>FEBS Letters</i> , 2000, 470, 125-130.	1.3	15
203	Dynamics of the Metallo- β -Lactamase from <i>Bacteroides fragilis</i> in the Presence and Absence of a Tight-Binding Inhibitor. <i>Biochemistry</i> , 2000, 39, 13356-13364.	1.2	59
204	Alternative Splicing of Wilms' Tumor Suppressor Protein Modulates DNA Binding Activity through Isoform-Specific DNA-Induced Conformational Changes. <i>Biochemistry</i> , 2000, 39, 5341-5348.	1.2	58
205	Changes in the Apomyoglobin Folding Pathway Caused by Mutation of the Distal Histidine Residue. <i>Biochemistry</i> , 2000, 39, 11227-11237.	1.2	68
206	Native and Non-native Secondary Structure and Dynamics in the pH 4 Intermediate of Apomyoglobin. <i>Biochemistry</i> , 2000, 39, 2894-2901.	1.2	121
207	Amide proton hydrogen exchange rates for sperm whale myoglobin obtained from ^{15}N NMR spectra. <i>Protein Science</i> , 2000, 9, 186-193.	3.1	23
208	Copper binding to the prion protein: Structural implications of four identical cooperative binding sites. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1999, 96, 2042-2047.	3.3	520
209	Characterization of monomeric and dimeric B domain of Staphylococcal protein A. <i>Chemical Biology and Drug Design</i> , 1999, 54, 344-352.	1.2	5
210	Inherent flexibility in a potent inhibitor of blood coagulation, recombinant nematode anticoagulant protein c2. <i>FEBS Journal</i> , 1999, 265, 539-548.	0.2	42
211	Backbone resonance assignments for the Fv fragment of the catalytic antibody NPN43C9 with bound p-nitrophenol. <i>Journal of Biomolecular NMR</i> , 1999, 15, 83-84.	1.6	4
212	Improved low pH bicelle system for orienting macromolecules over a wide temperature range. <i>Journal of Biomolecular NMR</i> , 1999, 13, 387-391.	1.6	84
213	Assignment of ^1H , ^{13}C , and ^{15}N resonances of reduced <i>Escherichia coli</i> glutaredoxin 2. <i>Journal of Biomolecular NMR</i> , 1999, 14, 197-198.	1.6	3
214	Association between the first two immunoglobulin-like domains of the neural cell adhesion molecule N-CAM. <i>FEBS Letters</i> , 1999, 451, 162-168.	1.3	30
215	NMR Characterization of the Metallo- β -lactamase from <i>Bacteroides fragilis</i> and Its Interaction with a Tight-Binding Inhibitor: Role of an Active-Site Loop. <i>Biochemistry</i> , 1999, 38, 14507-14514.	1.2	104
216	Effect of H helix destabilizing mutations on the kinetic and equilibrium folding of apomyoglobin 1 Edited by F. Cohen. <i>Journal of Molecular Biology</i> , 1999, 285, 269-282.	2.0	79

#	ARTICLE	IF	CITATIONS
217	Structural analyses of CREB-CBP transcriptional activator-coactivator complexes by NMR spectroscopy: implications for mapping the boundaries of structural domains. <i>Journal of Molecular Biology</i> , 1999, 287, 859-865.	2.0	68
218	Intrinsically unstructured proteins: re-assessing the protein structure-function paradigm. <i>Journal of Molecular Biology</i> , 1999, 293, 321-331.	2.0	2,668
219	Quench-flow experiments combined with mass spectrometry show apomyoglobin folds through an obligatory intermediate. <i>Protein Science</i> , 1999, 8, 45-49.	3.1	93
220	A NOESY-HSQC simulation program, SPIRIT. <i>Journal of Biomolecular NMR</i> , 1998, 11, 17-29.	1.6	17
221	¹ H, ¹³ C and ¹⁵ N NMR backbone assignments of 25.5 kDa metallo-beta-lactamase from <i>Bacteroides fragilis</i> . <i>Journal of Biomolecular NMR</i> , 1998, 12, 201-202.	1.6	5
222	Structural and dynamic characterization of partially folded states of apomyoglobin and implications for protein folding. <i>Nature Structural Biology</i> , 1998, 18, 148-155.	9.7	344
223	Equilibrium NMR studies of unfolded and partially folded proteins. <i>Nature Structural Biology</i> , 1998, 5, 499-503.	9.7	187
224	Sequence requirements for stabilization of a peptide reverse turn in water solution. Proline is not essential for stability. <i>FEBS Journal</i> , 1998, 255, 462-471.	0.2	25
225	NMR characterization of a single-cysteine mutant of <i>Escherichia coli</i> thioredoxin and a covalent thioredoxin-peptide complex. <i>FEBS Journal</i> , 1998, 257, 299-308.	0.2	14
226	The identification of metal-binding ligand residues in metalloproteins using nuclear magnetic resonance spectroscopy. <i>Protein Science</i> , 1998, 7, 2476-2479.	3.1	6
227	Conformational preferences in the Ser133 -phosphorylated and non-phosphorylated forms of the kinase inducible transactivation domain of CREB. <i>FEBS Letters</i> , 1998, 430, 317-322.	1.3	137
228	Calculations of Electrostatic Interactions and pKas in the Active Site of <i>Escherichia coli</i> Thioredoxin. <i>Biochemistry</i> , 1998, 37, 10298-10306.	1.2	82
229	Glycosylation of Threonine of the Repeating Unit of RNA Polymerase II with ² -Linked N-Acetylglucosamine Leads to a Turnlike Structure. <i>Journal of the American Chemical Society</i> , 1998, 120, 11567-11575.	6.6	73
230	High-resolution solution structure of the retinoid X receptor DNA-binding domain. <i>Journal of Molecular Biology</i> , 1998, 281, 271-284.	2.0	58
231	DNA-induced conformational changes are the basis for cooperative dimerization by the DNA binding domain of the retinoid X receptor. <i>Journal of Molecular Biology</i> , 1998, 284, 533-539.	2.0	66
232	Structure of the recombinant full-length hamster prion protein PrP(29-231): The N terminus is highly flexible. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1997, 94, 13452-13457.	3.3	665
233	Effects of Buried Charged Groups on Cysteine Thiol Ionization and Reactivity in <i>Escherichia coli</i> Thioredoxin: A Structural and Functional Characterization of Mutants of Asp 26 and Lys 57. <i>Biochemistry</i> , 1997, 36, 2622-2636.	1.2	192
234	Contribution of Increased Length and Intact Capping Sequences to the Conformational Preference for Helix in a 31-Residue Peptide from the C Terminus of Myohemerythrin. <i>Biochemistry</i> , 1997, 36, 5234-5244.	1.2	44

#	ARTICLE	IF	CITATIONS
235	Structure-based design of a constrained peptide mimic of the HIV-1 V3 loop neutralization site 1 Edited by F.E. Cohen. <i>Journal of Molecular Biology</i> , 1997, 266, 31-39.	2.0	77
236	Solution Structure of the KIX Domain of CBP Bound to the Transactivation Domain of CREB: A Model for Activator:Coactivator Interactions. <i>Cell</i> , 1997, 91, 741-752.	13.5	705
237	Populating the equilibrium molten globule state of apomyoglobin under conditions suitable for structural characterization by NMR. <i>FEBS Letters</i> , 1997, 417, 92-96.	1.3	53
238	Chemical shift dispersion and secondary structure prediction in unfolded and partly folded proteins. <i>FEBS Letters</i> , 1997, 419, 285-289.	1.3	135
239	PCR-based gene synthesis and protein NMR spectroscopy. <i>Structure</i> , 1997, 5, 1407-1412.	1.6	22
240	Folding propensities of peptide fragments of myoglobin. <i>Protein Science</i> , 1997, 6, 706-716.	3.1	82
241	Absence of a stable intermediate on the folding pathway of protein A. <i>Protein Science</i> , 1997, 6, 1449-1457.	3.1	117
242	INSIGHTS INTO PROTEIN FOLDING FROM NMR. <i>Annual Review of Physical Chemistry</i> , 1996, 47, 369-395.	4.8	144
243	NMR Solution Structure of Cu(I) Rusticyanin from <i>Thiobacillus ferrooxidans</i> : Structural Basis for the Extreme Acid Stability and Redox Potential. <i>Journal of Molecular Biology</i> , 1996, 263, 752-767.	2.0	104
244	Gene synthesis, high-level expression and assignment of backbone ¹⁵ N and ¹³ C resonances of soybean leghemoglobin. <i>FEBS Letters</i> , 1996, 399, 283-289.	1.3	12
245	Direct Measurement of the Aspartic Acid 26 pKa for Reduced <i>Escherichia coli</i> Thioredoxin by ¹³ C NMR. <i>Biochemistry</i> , 1996, 35, 1-6.	1.2	72
246	Solution conformation of an immunogenic peptide derived from the principal neutralizing determinant of the HIV-2 envelope glycoprotein gp125. <i>Folding & Design</i> , 1996, 1, 157-165.	4.5	8
247	Replacement of Trp28 in <i>Escherichia coli</i> Thioredoxin by Site-directed Mutagenesis Affects Thermodynamic Stability but Not Function. <i>Journal of Biological Chemistry</i> , 1996, 271, 3091-3096.	1.6	24
248	[³⁰] Nuclear magnetic resonance of thioredoxin and glutaredoxin. <i>Methods in Enzymology</i> , 1995, 252, 293-306.	0.4	0
249	¹ H, ¹³ C and ¹⁵ N chemical shift referencing in biomolecular NMR. <i>Journal of Biomolecular NMR</i> , 1995, 6, 135-140.	1.6	2,216
250	Random coil ¹ H chemical shifts obtained as a function of temperature and trifluoroethanol concentration for the peptide series GGXGG. <i>Journal of Biomolecular NMR</i> , 1995, 5, 14-24.	1.6	476
251	Antigenic peptides. <i>FASEB Journal</i> , 1995, 9, 37-42.	0.2	96
252	Detection of a Catalytic Antibody Species Acylated at the Active Site by Electrospray Mass Spectrometry. <i>Biochemistry</i> , 1995, 34, 720-723.	1.2	41

#	ARTICLE	IF	CITATIONS
253	Comparison of the Hydrogen-Exchange Behavior of Reduced and Oxidized Escherichia Coli Thioredoxin. <i>Biochemistry</i> , 1995, 34, 611-619.	1.2	61
254	Proton Sharing between Cysteine Thiols in Escherichia coli Thioredoxin: Implications for the Mechanism of Protein Disulfide Reduction. <i>Biochemistry</i> , 1995, 34, 10101-10105.	1.2	96
255	Complete ¹³ C assignments for recombinant Cu(I) rusticyanin prediction of secondary structure from patterns of chemical shifts. <i>FEBS Letters</i> , 1995, 365, 35-41.	1.3	6
256	Gene Synthesis, High-Level Expression, and Mutagenesis of Thiobacillus ferrooxidans Rusticyanin: His 85 Is a Ligand to the Blue Copper Center. <i>Biochemistry</i> , 1995, 34, 6640-6648.	1.2	66
257	Effect of disulfide bridge formation on the NMR spectrum of a protein: Studies on oxidized and reduced Escherichia coli thioredoxin. <i>Journal of Biomolecular NMR</i> , 1994, 4, 411-32.	1.6	30
258	High-resolution solution structures of oxidized and reduced Escherichia coli thioredoxin. <i>Structure</i> , 1994, 2, 853-868.	1.6	281
259	Characterization by ¹ H NMR of a C32S,C35S double mutant of Escherichia coli thioredoxin confirms its resemblance to the reduced wild-type protein. <i>FEBS Letters</i> , 1994, 339, 11-17.	1.3	22
260	Stabilization of a type VI turn in a family of linear peptides in water solution. <i>Journal of Molecular Biology</i> , 1994, 243, 736-753.	2.0	152
261	Three-dimensional structure of a type VI turn in a linear peptide in water solution Evidence for stacking of aromatic rings as a major stabilizing factor. <i>Journal of Molecular Biology</i> , 1994, 243, 754-766.	2.0	118
262	[¹³ C] Use of chemical shifts and coupling constants in nuclear magnetic resonance structural studies on peptides and proteins. <i>Methods in Enzymology</i> , 1994, 239, 392-416.	0.4	111
263	Binding of hapten to a single-chain catalytic antibody demonstrated by electrospray mass spectrometry. <i>Journal of the American Chemical Society</i> , 1994, 116, 7937-7938.	6.6	29
264	Differential Side Chain Hydration in a Linear Peptide Containing a Type VI Turn. <i>Journal of the American Chemical Society</i> , 1994, 116, 12051-12052.	6.6	35
265	The Folding Pathway of Apomyoglobin. <i>NATO ASI Series Series B: Physics</i> , 1994, , 7-18.	0.2	4
266	Peptide models of protein folding initiation sites. 3. The G-H helical hairpin of myoglobin. <i>Biochemistry</i> , 1993, 32, 6356-6364.	1.2	109
267	Peptide models of protein folding initiation sites. 1. Secondary structure formation by peptides corresponding to the G- and H-helices of myoglobin. <i>Biochemistry</i> , 1993, 32, 6337-6347.	1.2	209
268	Peptide models of protein folding initiation sites. 2. The G-H turn region of myoglobin acts as a helix stop signal. <i>Biochemistry</i> , 1993, 32, 6348-6355.	1.2	92
269	Peptide conformation and protein folding. <i>Current Opinion in Structural Biology</i> , 1993, 3, 60-65.	2.6	232
270	Characterization of a folding intermediate of apoplastocyanin trapped by proline isomerization. <i>Biochemistry</i> , 1993, 32, 12299-12310.	1.2	103

#	ARTICLE	IF	CITATIONS
271	Comparison of backbone and tryptophan side-chain dynamics of reduced and oxidized Escherichia coli thioredoxin using nitrogen-15 NMR relaxation measurements. <i>Biochemistry</i> , 1993, 32, 426-435.	1.2	148
272	Immunogenic peptides corresponding to the dominant antigenic region alanine-597 to cysteine-619 in the transmembrane protein of simian immunodeficiency virus have a propensity to fold in aqueous solution. <i>Biochemistry</i> , 1992, 31, 1458-1463.	1.2	15
273	Folding of peptide fragments comprising the complete sequence of proteins. <i>Journal of Molecular Biology</i> , 1992, 226, 795-817.	2.0	385
274	Folding of peptide fragments comprising the complete sequence of proteins. <i>Journal of Molecular Biology</i> , 1992, 226, 819-835.	2.0	226
275	Solution conformational preferences of immunogenic peptides derived from the principal neutralizing determinant of the HIV-1 envelope glycoprotein gp120. <i>Biochemistry</i> , 1991, 30, 9187-9194.	1.2	155
276	Proton-transfer effects in the active-site region of Escherichia coli thioredoxin using two-dimensional proton NMR. <i>Biochemistry</i> , 1991, 30, 4262-4268.	1.2	76
277	Polypeptide backbone resonance assignments and secondary structure of Bacillus subtilis enzyme Illglc determined by two-dimensional and three-dimensional heteronuclear NMR spectroscopy. <i>Biochemistry</i> , 1991, 30, 6896-6907.	1.2	56
278	Assignment of the ¹⁵ N NMR spectra of reduced and oxidized Escherichia coli thioredoxin. <i>FEBS Letters</i> , 1991, 284, 178-183.	1.3	39
279	Three-dimensional solution structure of the reduced form of Escherichia coli thioredoxin determined by nuclear magnetic resonance spectroscopy. <i>Biochemistry</i> , 1990, 29, 4129-4136.	1.2	177
280	Conformational preferences of synthetic peptides derived from the immunodominant site of the circumsporozoite protein of Plasmodium falciparum by proton NMR. <i>Biochemistry</i> , 1990, 29, 7828-7837.	1.2	94
281	Antigen-antibody interactions: An NMR approach. <i>Biochemical Pharmacology</i> , 1990, 40, 83-88.	2.0	21
282	Proton NMR studies of the solution conformations of an analog of the C-peptide of ribonuclease A. <i>Biochemistry</i> , 1989, 28, 7059-7064.	1.2	162
283	Assignment of the proton NMR spectrum of reduced and oxidized thioredoxin: sequence-specific assignments, secondary structure, and global fold. <i>Biochemistry</i> , 1989, 28, 7074-7087.	1.2	102
284	Structural differences between oxidized and reduced thioredoxin monitored by two-dimensional ¹ H NMR spectroscopy. <i>FEBS Letters</i> , 1988, 228, 254-258.	1.3	38
285	Folding of immunogenic peptide fragments of proteins in water solution. <i>Journal of Molecular Biology</i> , 1988, 201, 161-200.	2.0	685
286	Folding of immunogenic peptide fragments of proteins in water solution. <i>Journal of Molecular Biology</i> , 1988, 201, 201-217.	2.0	477
287	Conformation of peptide fragments of proteins in aqueous solution: implications for initiation of protein folding. <i>Biochemistry</i> , 1988, 27, 7167-7175.	1.2	505
288	Selection by Site-Directed Antibodies of Small Regions of Peptides which are Ordered in Water. <i>Novartis Foundation Symposium</i> , 1986, 119, 58-75.	1.2	1

#	ARTICLE	IF	CITATIONS
289	The immunodominant site of a synthetic immunogen has a conformational preference in water for a type-II reverse turn. <i>Nature</i> , 1985, 318, 480-483.	13.7	246
290	Chapter 5. NMR Studies of Disordered but Functional Proteins. <i>RSC Biomolecular Sciences</i> , 0, , 111-129.	0.4	0
291	Structural and dynamic studies of DNA recognition by NF- κ B p50 RHR homodimer using methyl NMR spectroscopy. <i>Nucleic Acids Research</i> , 0, , .	6.5	1