

H Jane Dyson

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278
papers

35,255
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92
h-index

185
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402
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38,132
ext. citations

7.7
avg, IF

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L-index

#	Paper	IF	Citations
278	Intrinsically unstructured proteins and their functions. <i>Nature Reviews Molecular Cell Biology</i> , 2005 , 6, 197-208	48.7	2985
277	Intrinsically unstructured proteins: re-assessing the protein structure-function paradigm. <i>Journal of Molecular Biology</i> , 1999 , 293, 321-31	6.5	2343
276	¹ H, ¹³ C and ¹⁵ N chemical shift referencing in biomolecular NMR. <i>Journal of Biomolecular NMR</i> , 1995 , 6, 135-40	3	2054
275	Intrinsically disordered proteins in cellular signalling and regulation. <i>Nature Reviews Molecular Cell Biology</i> , 2015 , 16, 18-29	48.7	1249
274	Coupling of folding and binding for unstructured proteins. <i>Current Opinion in Structural Biology</i> , 2002 , 12, 54-60	8.1	1121
273	Mechanism of coupled folding and binding of an intrinsically disordered protein. <i>Nature</i> , 2007 , 447, 1021-5	5.4	852
272	Linking folding and binding. <i>Current Opinion in Structural Biology</i> , 2009 , 19, 31-8	8.1	813
271	The dynamic energy landscape of dihydrofolate reductase catalysis. <i>Science</i> , 2006 , 313, 1638-42	33.3	778
270	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, 741-52	56.2	648
269	Folding of immunogenic peptide fragments of proteins in water solution. I. Sequence requirements for the formation of a reverse turn. <i>Journal of Molecular Biology</i> , 1988 , 201, 161-200	6.5	634
268	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, 13452-7	11.5	626
267	Unfolded proteins and protein folding studied by NMR. <i>Chemical Reviews</i> , 2004 , 104, 3607-22	68.1	541
266	Sequence-dependent correction of random coil NMR chemical shifts. <i>Journal of the American Chemical Society</i> , 2001 , 123, 2970-8	16.4	509
265	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, 2042-7	11.5	485
264	Conformation of peptide fragments of proteins in aqueous solution: implications for initiation of protein folding. <i>Biochemistry</i> , 1988 , 27, 7167-75	3.2	464
263	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	3	453
262	Folding of immunogenic peptide fragments of proteins in water solution. II. The nascent helix. <i>Journal of Molecular Biology</i> , 1988 , 201, 201-17	6.5	449

261	Structure, dynamics, and catalytic function of dihydrofolate reductase. <i>Annual Review of Biophysics and Biomolecular Structure</i> , 2004 , 33, 119-40		384
260	Mutual synergistic folding in recruitment of CBP/p300 by p160 nuclear receptor coactivators. <i>Nature</i> , 2002 , 415, 549-53	50.4	373
259	An NMR perspective on enzyme dynamics. <i>Chemical Reviews</i> , 2006 , 106, 3055-79	68.1	369
258	Folding of peptide fragments comprising the complete sequence of proteins. Models for initiation of protein folding. I. Myohemerythrin. <i>Journal of Molecular Biology</i> , 1992 , 226, 795-817	6.5	351
257	A dynamic knockout reveals that conformational fluctuations influence the chemical step of enzyme catalysis. <i>Science</i> , 2011 , 332, 234-8	33.3	350
256	Structural basis for Hif-1 alpha /CBP recognition in the cellular hypoxic response. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2002 , 99, 5271-6	11.5	333
255	Structural and dynamic characterization of partially folded states of apomyoglobin and implications for protein folding. <i>Nature Structural Biology</i> , 1998 , 5, 148-55		318
254	Recognition of the mRNA AU-rich element by the zinc finger domain of TIS11d. <i>Nature Structural and Molecular Biology</i> , 2004 , 11, 257-64	17.6	280
253	High-resolution solution structures of oxidized and reduced Escherichia coli thioredoxin. <i>Structure</i> , 1994 , 2, 853-68	5.2	267
252	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, 43-8	3	258
251	Backbone dynamics in dihydrofolate reductase complexes: role of loop flexibility in the catalytic mechanism. <i>Biochemistry</i> , 2001 , 40, 9846-59	3.2	226
250	The immunodominant site of a synthetic immunogen has a conformational preference in water for a type-II reverse turn. <i>Nature</i> , 1985 , 318, 480-3	50.4	221
249	Folding of peptide fragments comprising the complete sequence of proteins. Models for initiation of protein folding. II. Plastocyanin. <i>Journal of Molecular Biology</i> , 1992 , 226, 819-35	6.5	215
248	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, 3561-71	3.2	203
247	Peptide conformation and protein folding. <i>Current Opinion in Structural Biology</i> , 1993 , 3, 60-65	8.1	202
246	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, 6337-47	3.2	198
245	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, 13057-61	11.5	195
244	Insights into the structure and dynamics of unfolded proteins from nuclear magnetic resonance. <i>Advances in Protein Chemistry</i> , 2002 , 62, 311-40		183

243	Effects of buried charged groups on cysteine thiol ionization and reactivity in Escherichia coli thioredoxin: structural and functional characterization of mutants of Asp 26 and Lys 57. <i>Biochemistry</i> , 1997 , 36, 2622-36	3.2	176
242	What's in a name? Why these proteins are intrinsically disordered: Why these proteins are intrinsically disordered. <i>Intrinsically Disordered Proteins</i> , 2013 , 1, e24157		171
241	Equilibrium NMR studies of unfolded and partially folded proteins. <i>Nature Structural Biology</i> , 1998 , 5 Suppl, 499-503		169
240	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, 6714-22	5.4	168
239	Local structural plasticity of the prion protein. Analysis of NMR relaxation dynamics. <i>Biochemistry</i> , 2001 , 40, 2743-53	3.2	164
238	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, 9614-9	11.5	161
237	Three-dimensional solution structure of the reduced form of Escherichia coli thioredoxin determined by nuclear magnetic resonance spectroscopy. <i>Biochemistry</i> , 1990 , 29, 4129-36	3.2	161
236	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, 6591-6	11.5	159
235	Structural characterization of unfolded states of apomyoglobin using residual dipolar couplings. <i>Journal of Molecular Biology</i> , 2004 , 340, 1131-42	6.5	157
234	Solution structure of the KIX domain of CBP bound to the transactivation domain of c-Myb. <i>Journal of Molecular Biology</i> , 2004 , 337, 521-34	6.5	154
233	Solution conformational preferences of immunogenic peptides derived from the principal neutralizing determinant of the HIV-1 envelope glycoprotein gp120. <i>Biochemistry</i> , 1991 , 30, 9187-94	3.2	147
232	¹ H NMR studies of the solution conformations of an analogue of the C-peptide of ribonuclease A. <i>Biochemistry</i> , 1989 , 28, 7059-64	3.2	146
231	Structural basis for cooperative transcription factor binding to the CBP coactivator. <i>Journal of Molecular Biology</i> , 2006 , 355, 1005-13	6.5	144
230	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, 262-73	6.5	143
229	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, 5032-7	11.5	143
228	Stabilization of a type VI turn in a family of linear peptides in water solution. <i>Journal of Molecular Biology</i> , 1994 , 243, 736-53	6.5	143
227	Comparison of backbone and tryptophan side-chain dynamics of reduced and oxidized Escherichia coli thioredoxin using ¹⁵ N NMR relaxation measurements. <i>Biochemistry</i> , 1993 , 32, 426-35	3.2	140
226	Cooperativity in transcription factor binding to the coactivator CREB-binding protein (CBP). The mixed lineage leukemia protein (MLL) activation domain binds to an allosteric site on the KIX domain. <i>Journal of Biological Chemistry</i> , 2002 , 277, 43168-74	5.4	139

225	Nuclear magnetic resonance methods for elucidation of structure and dynamics in disordered states. <i>Methods in Enzymology</i> , 2001 , 339, 258-70	1.7	137
224	Insights into protein folding from NMR. <i>Annual Review of Physical Chemistry</i> , 1996 , 47, 369-95	15.7	137
223	Solution structure and acetyl-lysine binding activity of the GCN5 bromodomain. <i>Journal of Molecular Biology</i> , 2000 , 304, 355-70	6.5	135
222	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, 2352-7	11.5	134
221	Expanding the proteome: disordered and alternatively folded proteins. <i>Quarterly Reviews of Biophysics</i> , 2011 , 44, 467-518	7	132
220	Mapping long-range contacts in a highly unfolded protein. <i>Journal of Molecular Biology</i> , 2002 , 322, 655-665	6.5	130
219	Structure of the p53 transactivation domain in complex with the nuclear receptor coactivator binding domain of CREB binding protein. <i>Biochemistry</i> , 2010 , 49, 9964-71	3.2	129
218	DNA-induced alpha-helix capping in conserved linker sequences is a determinant of binding affinity in Cys(2)-His(2) zinc fingers. <i>Journal of Molecular Biology</i> , 2000 , 295, 719-27	6.5	129
217	Conformational preferences in the Ser133-phosphorylated and non-phosphorylated forms of the kinase inducible transactivation domain of CREB. <i>FEBS Letters</i> , 1998 , 430, 317-22	3.8	125
216	Functional advantages of dynamic protein disorder. <i>FEBS Letters</i> , 2015 , 589, 2433-40	3.8	124
215	Molecular hinges in protein folding: the urea-denatured state of apomyoglobin. <i>Biochemistry</i> , 2002 , 41, 12681-6	3.2	123
214	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, 42241-8	5.4	122
213	Chemical shift dispersion and secondary structure prediction in unfolded and partly folded proteins. <i>FEBS Letters</i> , 1997 , 419, 285-9	3.8	121
212	Amylin proprotein processing generates progressively more amyloidogenic peptides that initially sample the helical state. <i>Biochemistry</i> , 2008 , 47, 9900-10	3.2	121
211	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, 1373-8	11.5	118
210	Native and non-native secondary structure and dynamics in the pH 4 intermediate of apomyoglobin. <i>Biochemistry</i> , 2000 , 39, 2894-901	3.2	116
209	Absence of a stable intermediate on the folding pathway of protein A. <i>Protein Science</i> , 1997 , 6, 1449-57	6.3	113
208	Structural basis for recruitment of CBP/p300 coactivators by STAT1 and STAT2 transactivation domains. <i>EMBO Journal</i> , 2009 , 28, 948-58	13	108

207	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, 754-66	6.5	108
206	The client protein p53 adopts a molten globule-like state in the presence of Hsp90. <i>Nature Structural and Molecular Biology</i> , 2011 , 18, 537-41	17.6	105
205	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, 317-31	3	105
204	SANE (Structure Assisted NOE Evaluation): an automated model-based approach for NOE assignment. <i>Journal of Biomolecular NMR</i> , 2001 , 19, 321-9	3	105
203	Divergent evolution of protein conformational dynamics in dihydrofolate reductase. <i>Nature Structural and Molecular Biology</i> , 2013 , 20, 1243-9	17.6	104
202	Conformational changes in the active site loops of dihydrofolate reductase during the catalytic cycle. <i>Biochemistry</i> , 2004 , 43, 16046-55	3.2	103
201	Molecular basis for modulation of biological function by alternate splicing of the WilmsPtumor suppressor protein. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2000 , 97, 11932-5	11.5	103
200	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, 13260-5	11.5	102
199	Solution structure of the cysteine-rich domain of the Escherichia coli chaperone protein DnaJ. <i>Journal of Molecular Biology</i> , 2000 , 300, 805-18	6.5	102
198	Solution structure of the Hdm2 C2H2C4 RING, a domain critical for ubiquitination of p53. <i>Journal of Molecular Biology</i> , 2006 , 363, 433-50	6.5	101
197	Solution structure of the TAZ2 (CH3) domain of the transcriptional adaptor protein CBP. <i>Journal of Molecular Biology</i> , 2000 , 303, 243-53	6.5	101
196	Peptide models of protein folding initiation sites. 3. The G-H helical hairpin of myoglobin. <i>Biochemistry</i> , 1993 , 32, 6356-64	3.2	101
195	Hypersensitive termination of the hypoxic response by a disordered protein switch. <i>Nature</i> , 2017 , 543, 447-451	50.4	99
194	NMR characterization of the metallo-beta-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, 14507-14	3.2	99
193	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	1.7	99
192	Identification of native and non-native structure in kinetic folding intermediates of apomyoglobin. <i>Journal of Molecular Biology</i> , 2006 , 355, 139-56	6.5	98
191	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, 6278-83	11.5	95
190	Characterization of a folding intermediate of apoplastocyanin trapped by proline isomerization. <i>Biochemistry</i> , 1993 , 32, 12299-310	3.2	95

189	Assignment of the proton NMR spectrum of reduced and oxidized thioredoxin: sequence-specific assignments, secondary structure, and global fold. <i>Biochemistry</i> , 1989 , 28, 7074-87	3.2	95
188	Proton sharing between cysteine thiols in Escherichia coli thioredoxin: implications for the mechanism of protein disulfide reduction. <i>Biochemistry</i> , 1995 , 34, 10101-5	3.2	93
187	Finding Our Way in the Dark Proteome. <i>Journal of the American Chemical Society</i> , 2016 , 138, 9730-42	16.4	93
186	Mapping the interactions of the p53 transactivation domain with the KIX domain of CBP. <i>Biochemistry</i> , 2009 , 48, 2115-24	3.2	90
185	Antigenic peptides. <i>FASEB Journal</i> , 1995 , 9, 37-42	0.9	89
184	Conservation of folding pathways in evolutionarily distant globin sequences. <i>Nature Structural Biology</i> , 2000 , 7, 679-86		88
183	The apomyoglobin folding pathway revisited: structural heterogeneity in the kinetic burst phase intermediate. <i>Journal of Molecular Biology</i> , 2002 , 322, 483-9	6.5	86
182	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, 6348-55	3.2	86
181	NMR solution structure of Cu(I) rusticyanin from Thiobacillus ferrooxidans: structural basis for the extreme acid stability and redox potential. <i>Journal of Molecular Biology</i> , 1996 , 263, 752-67	6.5	85
180	Evaluating beta-turn mimics as beta-sheet folding nucleators. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009 , 106, 11067-72	11.5	84
179	Expanding the Paradigm: Intrinsically Disordered Proteins and Allosteric Regulation. <i>Journal of Molecular Biology</i> , 2018 , 430, 2309-2320	6.5	82
178	Structure of the Wilms tumor suppressor protein zinc finger domain bound to DNA. <i>Journal of Molecular Biology</i> , 2007 , 372, 1227-45	6.5	82
177	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, 15229-34	11.5	80
176	Quench-flow experiments combined with mass spectrometry show apomyoglobin folds through and obligatory intermediate. <i>Protein Science</i> , 1999 , 8, 45-9	6.3	80
175	Elucidation of the protein folding landscape by NMR. <i>Methods in Enzymology</i> , 2005 , 394, 299-321	1.7	80
174	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, 13859-64	11.5	79
173	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, 819-32	6.5	79
172	Conformational preferences of synthetic peptides derived from the immunodominant site of the circumsporozoite protein of Plasmodium falciparum by 1H NMR. <i>Biochemistry</i> , 1990 , 29, 7828-37	3.2	79

171	Structural and energetic basis of carbohydrate-aromatic packing interactions in proteins. <i>Journal of the American Chemical Society</i> , 2013 , 135, 9877-84	16.4	77
170	Improved low pH bicelle system for orienting macromolecules over a wide temperature range. <i>Journal of Biomolecular NMR</i> , 1999 , 13, 387-91	3	77
169	Folding propensities of peptide fragments of myoglobin. <i>Protein Science</i> , 1997 , 6, 706-16	6.3	76
168	Interaction of the TAZ1 domain of the CREB-binding protein with the activation domain of CITED2: regulation by competition between intrinsically unstructured ligands for non-identical binding sites. <i>Journal of Biological Chemistry</i> , 2004 , 279, 3042-9	5.4	76
167	Calculations of electrostatic interactions and pKas in the active site of Escherichia coli thioredoxin. <i>Biochemistry</i> , 1998 , 37, 10298-306	3.2	76
166	Activation of the redox-regulated chaperone Hsp33 by domain unfolding. <i>Journal of Biological Chemistry</i> , 2004 , 279, 20529-38	5.4	75
165	Structural and dynamic characterization of an unfolded state of poplar apo-plastocyanin formed under nondenaturing conditions. <i>Protein Science</i> , 2001 , 10, 1056-66	6.3	75
164	Glycosylation of Threonine of the Repeating Unit of RNA Polymerase II with β -Linked N-Acetylglucosamine Leads to a Turnlike Structure. <i>Journal of the American Chemical Society</i> , 1998 , 120, 11567-11575	16.4	73
163	Structure-based design of a constrained peptide mimic of the HIV-1 V3 loop neutralization site. <i>Journal of Molecular Biology</i> , 1997 , 266, 31-9	6.5	72
162	ZZ domain of CBP: an unusual zinc finger fold in a protein interaction module. <i>Journal of Molecular Biology</i> , 2004 , 343, 1081-93	6.5	72
161	Effect of H helix destabilizing mutations on the kinetic and equilibrium folding of apomyoglobin. <i>Journal of Molecular Biology</i> , 1999 , 285, 269-82	6.5	72
160	NMR relaxation study of the complex formed between CBP and the activation domain of the nuclear hormone receptor coactivator ACTR. <i>Biochemistry</i> , 2008 , 47, 1299-308	3.2	71
159	Effect of cofactor binding and loop conformation on side chain methyl dynamics in dihydrofolate reductase. <i>Biochemistry</i> , 2004 , 43, 374-83	3.2	71
158	Direct measurement of the aspartic acid 26 pKa for reduced Escherichia coli thioredoxin by ^{13}C NMR. <i>Biochemistry</i> , 1996 , 35, 1-6	3.2	69
157	A distal mutation perturbs dynamic amino acid networks in dihydrofolate reductase. <i>Biochemistry</i> , 2013 , 52, 4605-19	3.2	68
156	Solution structure of Escherichia coli glutaredoxin-2 shows similarity to mammalian glutathione-S-transferases. <i>Journal of Molecular Biology</i> , 2001 , 310, 907-18	6.5	67
155	Proton-transfer effects in the active-site region of Escherichia coli thioredoxin using two-dimensional ^1H NMR. <i>Biochemistry</i> , 1991 , 30, 4262-8	3.2	67
154	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, 859-65	6.5	66

153	NMR solution structure of the inserted domain of human leukocyte function associated antigen-1. <i>Journal of Molecular Biology</i> , 2000 , 295, 1251-64	6.5	65
152	Genomic-scale comparison of sequence- and structure-based methods of function prediction: does structure provide additional insight?. <i>Protein Science</i> , 2001 , 10, 1005-14	6.3	64
151	Changes in the apomyoglobin folding pathway caused by mutation of the distal histidine residue. <i>Biochemistry</i> , 2000 , 39, 11227-37	3.2	62
150	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, 533-9	6.5	62
149	CBP/p300 TAZ1 domain forms a structured scaffold for ligand binding. <i>Biochemistry</i> , 2005 , 44, 490-7	3.2	61
148	Roles of intrinsic disorder in protein-nucleic acid interactions. <i>Molecular BioSystems</i> , 2012 , 8, 97-104		60
147	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, 275-91	6.5	60
146	Packing, specificity, and mutability at the binding interface between the p160 coactivator and CREB-binding protein. <i>Protein Science</i> , 2004 , 13, 203-10	6.3	59
145	Comparison of the hydrogen-exchange behavior of reduced and oxidized Escherichia coli thioredoxin. <i>Biochemistry</i> , 1995 , 34, 611-9	3.2	59
144	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, 6640-8	3.2	59
143	Dynamic Interaction of Hsp90 with Its Client Protein p53. <i>Journal of Molecular Biology</i> , 2011 , 411, 158-736.5		58
142	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, 8725-34	3.2	57
141	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	11.5	56
140	Localized structural fluctuations promote amyloidogenic conformations in transthyretin. <i>Journal of Molecular Biology</i> , 2013 , 425, 977-88	6.5	56
139	Transfer of flexibility between ankyrin repeats in IkappaB* upon formation of the NF-kappaB complex. <i>Journal of Molecular Biology</i> , 2008 , 380, 917-31	6.5	56
138	Making Sense of Intrinsically Disordered Proteins. <i>Biophysical Journal</i> , 2016 , 110, 1013-6	2.9	55
137	High-resolution solution structure of the retinoid X receptor DNA-binding domain. <i>Journal of Molecular Biology</i> , 1998 , 281, 271-84	6.5	54
136	Localization of sites of interaction between p23 and Hsp90 in solution. <i>Journal of Biological Chemistry</i> , 2006 , 281, 14457-64	5.4	53

135	Role of a solvent-exposed tryptophan in the recognition and binding of antibiotic substrates for a metallo-beta-lactamase. <i>Protein Science</i> , 2003 , 12, 1368-75	6.3	53
134	Dynamics of the metallo-beta-lactamase from <i>Bacteroides fragilis</i> in the presence and absence of a tight-binding inhibitor. <i>Biochemistry</i> , 2000 , 39, 13356-64	3.2	53
133	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, 4765-70	11.5	52
132	Efficient inhibition of <i>Escherichia coli</i> RNA polymerase by the bacteriophage T4 AsiA protein requires that AsiA binds first to free sigma70. <i>Journal of Molecular Biology</i> , 2000 , 304, 731-9	6.5	52
131	Polypeptide backbone resonance assignments and secondary structure of <i>Bacillus subtilis</i> enzyme IIIglc determined by two-dimensional and three-dimensional heteronuclear NMR spectroscopy. <i>Biochemistry</i> , 1991 , 30, 6896-907	3.2	52
130	Conformational relaxation following hydride transfer plays a limiting role in dihydrofolate reductase catalysis. <i>Biochemistry</i> , 2008 , 47, 9227-33	3.2	50
129	Alternative splicing of Wilms tumor suppressor protein modulates DNA binding activity through isoform-specific DNA-induced conformational changes. <i>Biochemistry</i> , 2000 , 39, 5341-8	3.2	50
128	Prion proteins with pathogenic and protective mutations show similar structure and dynamics. <i>Biochemistry</i> , 2009 , 48, 8120-8	3.2	49
127	Tailoring relaxation dispersion experiments for fast-associating protein complexes. <i>Journal of the American Chemical Society</i> , 2007 , 129, 13406-7	16.4	49
126	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	6.5	48
125	Identification of Cys255 in HIF-1 as a novel site for development of covalent inhibitors of HIF-1/ARNT PasB domain protein-protein interaction. <i>Protein Science</i> , 2012 , 21, 1885-96	6.3	47
124	Populating the equilibrium molten globule state of apomyoglobin under conditions suitable for structural characterization by NMR. <i>FEBS Letters</i> , 1997 , 417, 92-6	3.8	47
123	Sequence determinants of a protein folding pathway. <i>Journal of Molecular Biology</i> , 2005 , 351, 383-92	6.5	46
122	The intrinsically disordered RNR inhibitor Sml1 is a dynamic dimer. <i>Biochemistry</i> , 2008 , 47, 13428-37	3.2	45
121	Conformational and dynamic characterization of the molten globule state of an apomyoglobin mutant with an altered folding pathway. <i>Biochemistry</i> , 2001 , 40, 14459-67	3.2	43
120	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	6.5	42
119	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	3	42
118	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	11.5	42

117	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, 5234-44	3.2	41
116	Prediction of the rotational tumbling time for proteins with disordered segments. <i>Journal of the American Chemical Society</i> , 2009 , 131, 6814-21	16.4	39
115	Detection of a catalytic antibody species acylated at the active site by electrospray mass spectrometry. <i>Biochemistry</i> , 1995 , 34, 720-3	3.2	39
114	The Dependence of Carbohydrate-Aromatic Interaction Strengths on the Structure of the Carbohydrate. <i>Journal of the American Chemical Society</i> , 2016 , 138, 7636-48	16.4	38
113	Diagnostic chemical shift markers for loop conformation and substrate and cofactor binding in dihydrofolate reductase complexes. <i>Protein Science</i> , 2003 , 12, 2230-8	6.3	38
112	Cofactor-Mediated Conformational Dynamics Promote Product Release From Escherichia coli Dihydrofolate Reductase via an Allosteric Pathway. <i>Journal of the American Chemical Society</i> , 2015 , 137, 9459-68	16.4	37
111	The CBP/p300 TAZ1 domain in its native state is not a binding partner of MDM2. <i>Biochemical Journal</i> , 2004 , 381, 685-91	3.8	37
110	Inherent flexibility in a potent inhibitor of blood coagulation, recombinant nematode anticoagulant protein c2. <i>FEBS Journal</i> , 1999 , 265, 539-48		36
109	Role of disorder in IB-NFB interaction. <i>IUBMB Life</i> , 2012 , 64, 499-505	4.7	35
108	Disulfide bonding arrangements in active forms of the somatomedin B domain of human vitronectin. <i>Biochemistry</i> , 2004 , 43, 6519-34	3.2	35
107	The zinc-dependent redox switch domain of the chaperone Hsp33 has a novel fold. <i>Journal of Molecular Biology</i> , 2004 , 341, 893-9	6.5	35
106	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	3	35
105	Assignment of the 15N NMR spectra of reduced and oxidized Escherichia coli thioredoxin. <i>FEBS Letters</i> , 1991 , 284, 178-83	3.8	35
104	Structural differences between oxidized and reduced thioredoxin monitored by two-dimensional 1H NMR spectroscopy. <i>FEBS Letters</i> , 1988 , 228, 254-8	3.8	35
103	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	11.5	33
102	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	16.4	33
101	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	3	29
100	How Does Your Protein Fold? Elucidating the Apomyoglobin Folding Pathway. <i>Accounts of Chemical Research</i> , 2017 , 50, 105-111	24.3	28

99	The RelA nuclear localization signal folds upon binding to I κ B. <i>Journal of Molecular Biology</i> , 2011 , 405, 754-64	6.5	28
98	Folding of a beta-sheet protein monitored by real-time NMR spectroscopy. <i>Journal of Molecular Biology</i> , 2003 , 328, 1161-71	6.5	28
97	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, 4500-4	3.8	26
96	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	16.4	26
95	Interaction of the I κ B C-terminal PEST sequence with NF- κ B: insights into the inhibition of NF- κ B DNA binding by I κ B. <i>Journal of Molecular Biology</i> , 2009 , 388, 824-38	6.5	25
94	Association between the first two immunoglobulin-like domains of the neural cell adhesion molecule N-CAM. <i>FEBS Letters</i> , 1999 , 451, 162-8	3.8	25
93	Replacement of Trp28 in Escherichia coli thioredoxin by site-directed mutagenesis affects thermodynamic stability but not function. <i>Journal of Biological Chemistry</i> , 1996 , 271, 3091-6	5.4	24
92	The kinetic and equilibrium molten globule intermediates of apoleghemoglobin differ in structure. <i>Journal of Molecular Biology</i> , 2008 , 378, 715-25	6.5	23
91	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, 161-72	6.5	23
90	Sequence requirements for stabilization of a peptide reverse turn in water solution--proline is not essential for stability. <i>FEBS Journal</i> , 1998 , 255, 462-71		22
89	Functional importance of stripping in NF κ B signaling revealed by a stripping-impaired I κ B mutant. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017 , 114, 1916-1921	11.5	21
88	Functional dynamics of the folded ankyrin repeats of I kappa B alpha revealed by nuclear magnetic resonance. <i>Biochemistry</i> , 2009 , 48, 8023-31	3.2	21
87	PCR-based gene synthesis and protein NMR spectroscopy. <i>Structure</i> , 1997 , 5, 1407-12	5.2	21
86	NMR illuminates intrinsic disorder. <i>Current Opinion in Structural Biology</i> , 2021 , 70, 44-52	8.1	21
85	Slow Dynamics of Tryptophan-Water Networks in Proteins. <i>Journal of the American Chemical Society</i> , 2018 , 140, 675-682	16.4	20
84	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	3	20
83	Detection of a ternary complex of NF- κ B and I κ B with DNA provides insights into how I κ B removes NF- κ B from transcription sites. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2011 , 108, 1367-72	11.5	20
82	Embryonic neural inducing factor churchill is not a DNA-binding zinc finger protein: solution structure reveals a solvent-exposed beta-sheet and zinc binuclear cluster. <i>Journal of Molecular Biology</i> , 2007 , 371, 1274-89	6.5	20

81	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	3.2	20
80	Defining the Structural Basis for Allosteric Product Release from E. coli Dihydrofolate Reductase Using NMR Relaxation Dispersion. <i>Journal of the American Chemical Society</i> , 2017 , 139, 11233-11240	16.4	19
79	Structure of the nuclear factor ALY: insights into post-transcriptional regulatory and mRNA nuclear export processes. <i>Biochemistry</i> , 2003 , 42, 7348-57	3.2	19
78	Antigen-antibody interactions: an NMR approach. <i>Biochemical Pharmacology</i> , 1990 , 40, 83-8	6	19
77	Amide proton hydrogen exchange rates for sperm whale myoglobin obtained from 15N-1H NMR spectra. <i>Protein Science</i> , 2000 , 9, 186-93	6.3	18
76	NMR solution structure of the peptide fragment 1-30, derived from unprocessed mouse Doppel protein, in DHPC micelles. <i>Biochemistry</i> , 2006 , 45, 159-66	3.2	18
75	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, 107-9	4.7	18
74	Characterization by 1H 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, 11-7	3.8	18
73	NMR Measurements Reveal the Structural Basis of Transthyretin Destabilization by Pathogenic Mutations. <i>Biochemistry</i> , 2018 , 57, 4421-4430	3.2	16
72	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	11.5	16
71	Kaiso uses all three zinc fingers and adjacent sequence motifs for high affinity binding to sequence-specific and methyl-CpG DNA targets. <i>FEBS Letters</i> , 2012 , 586, 734-9	3.8	16
70	Energetic frustration of apomyoglobin folding: role of the B helix. <i>Journal of Molecular Biology</i> , 2010 , 396, 1319-28	6.5	16
69	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, 718-30	6.5	16
68	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	3.2	15
67	Side-chain conformational heterogeneity of intermediates in the Escherichia coli dihydrofolate reductase catalytic cycle. <i>Biochemistry</i> , 2013 , 52, 3464-77	3.2	15
66	A NOESY-HSQC simulation program, SPIRIT. <i>Journal of Biomolecular NMR</i> , 1998 , 11, 17-29	3	15
65	Structural characterization of partially folded intermediates of apomyoglobin H64F. <i>Protein Science</i> , 2008 , 17, 313-21	6.3	15
64	Structural characterization of interactions between the double-stranded RNA-binding zinc finger protein JAZ and nucleic acids. <i>Biochemistry</i> , 2014 , 53, 1495-510	3.2	14

63	Consequences of stabilizing the natively disordered F helix for the folding pathway of apomyoglobin. <i>Journal of Molecular Biology</i> , 2011 , 411, 248-63	6.5	14
62	NMR characterization of a single-cysteine mutant of Escherichia coli thioredoxin and a covalent thioredoxin-peptide complex. <i>FEBS Journal</i> , 1998 , 257, 299-308		14
61	Structure discrimination for the C-terminal domain of Escherichia coli trigger factor in solution. <i>Journal of Biomolecular NMR</i> , 2008 , 40, 23-30	3	14
60	Identification of the regions involved in DNA binding by the mouse PEBP2alpha protein. <i>FEBS Letters</i> , 2000 , 470, 125-30	3.8	14
59	Fluorotryptophan Incorporation Modulates the Structure and Stability of Transthyretin in a Site-Specific Manner. <i>Biochemistry</i> , 2017 , 56, 5570-5581	3.2	13
58	Changes in structure and dynamics of the Fv fragment of a catalytic antibody upon binding of inhibitor. <i>Protein Science</i> , 2003 , 12, 1386-94	6.3	13
57	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, 1458-63	3.2	13
56	How Do Intrinsically Disordered Viral Proteins Hijack the Cell?. <i>Biochemistry</i> , 2018 , 57, 4045-4046	3.2	12
55	Hydrogen-deuterium exchange strategy for delineation of contact sites in protein complexes. <i>FEBS Letters</i> , 2008 , 582, 1495-500	3.8	12
54	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	6.3	12
53	The reduced, denatured somatomedin B domain of vitronectin refolds into a stable, biologically active molecule. <i>Biochemistry</i> , 2006 , 45, 3297-306	3.2	11
52	CH ₂ Hydrogen Bonds Mediate Highly Specific Recognition of Methylated CpG Sites by the Zinc Finger Protein Kaiso. <i>Biochemistry</i> , 2018 , 57, 2109-2120	3.2	10
51	Probing the non-native H helix translocation in apomyoglobin folding intermediates. <i>Biochemistry</i> , 2014 , 53, 3767-80	3.2	10
50	Long-range effects and functional consequences of stabilizing mutations in the ankyrin repeat domain of I β . <i>Journal of Molecular Biology</i> , 2013 , 425, 902-13	6.5	10
49	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	11.5	10
48	The CH ₂ domain of CBP/p300 is a novel zinc finger. <i>FEBS Letters</i> , 2013 , 587, 2506-11	3.8	9
47	Gene synthesis, high-level expression and assignment of backbone ¹⁵ N and ¹³ C resonances of soybean leghemoglobin. <i>FEBS Letters</i> , 1996 , 399, 283-9	3.8	9
46	Characterization of an Hsp90-Independent Interaction between Co-Chaperone p23 and Transcription Factor p53. <i>Biochemistry</i> , 2018 , 57, 935-944	3.2	8

45	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, 157-65		8
44	Management of Hsp90-Dependent Protein Folding by Small Molecules Targeting the Aha1 Co-Chaperone. <i>Cell Chemical Biology</i> , 2020 , 27, 292-305.e6	8.2	8
43	NMR detection of adventitious xylose binding to the quorum-sensing protein SdiA of Escherichia coli. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2007 , 17, 6202-5	2.9	7
42	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,	11.5	7
41	NMR characterization of a 72 kDa transcription factor using differential isotopic labeling. <i>Protein Science</i> , 2016 , 25, 597-604	6.3	6
40	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, 377-8	3	6
39	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, 373-4	3	6
38	Assignment of 1H, 13C and 15N resonances of the I-domain of human leukocyte function associated antigen-1. <i>Journal of Biomolecular NMR</i> , 2000 , 16, 271-2	3	6
37	Complete 13C assignments for recombinant Cu(I) rusticyanin. Prediction of secondary structure from patterns of chemical shifts. <i>FEBS Letters</i> , 1995 , 365, 35-41	3.8	6
36	Structural characterization of the ternary complex that mediates termination of NF- κ B signaling by I κ B. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016 , 113, 6212-7	11.5	6
35	Homodimerization of the PAS-B domains of hypoxia-inducible factors. <i>Journal of Physical Chemistry B</i> , 2012 , 116, 6960-5	3.4	5
34	1H, 13C and 15N NMR backbone assignments of 25.5 kDa metallo-beta-lactamase from Bacteroides fragilis. <i>Journal of Biomolecular NMR</i> , 1998 , 12, 201-2	3	5
33	The identification of metal-binding ligand residues in metalloproteins using nuclear magnetic resonance spectroscopy. <i>Protein Science</i> , 1998 , 7, 2476-9	6.3	5
32	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	6.3	4
31	Side chain conformational averaging in human dihydrofolate reductase. <i>Biochemistry</i> , 2014 , 53, 1134-45	3.2	4
30	Characterization of monomeric and dimeric B domain of Staphylococcal protein A. <i>Chemical Biology and Drug Design</i> , 1999 , 54, 344-52		4
29	Backbone resonance assignments for the Fv fragment of the catalytic antibody NPN43C9 with bound p-nitrophenol. <i>Journal of Biomolecular NMR</i> , 1999 , 15, 83-4	3	4
28	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,	11.5	4

27	The Folding Pathway of Apomyoglobin. <i>NATO ASI Series Series B: Physics</i> , 1994 , 7-18		4
26	Thermodynamic Stability and Aggregation Kinetics of EF Helix and EF Loop Variants of Transthyretin. <i>Biochemistry</i> , 2021 , 60, 756-764	3.2	4
25	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	3.2	3
24	Aggregation of zinc-free p53 is inhibited by Hsp90 but not other chaperones. <i>Protein Science</i> , 2019 , 28, 2020-2023	6.3	3
23	CheShift-2 resolves a local inconsistency between two X-ray crystal structures. <i>Journal of Biomolecular NMR</i> , 2012 , 54, 193-8	3	3
22	Backbone H(N), N, Calpha, CPand Cbeta assignments of the 19 kDa DHFR/NADPH complex at 9 degrees C and pH 7.6. <i>Journal of Biomolecular NMR</i> , 2000 , 16, 349-50	3	3
21	A Dynamic Switch in Inactive p38 Leads to an Excited State on the Pathway to an Active Kinase. <i>Biochemistry</i> , 2019 , 58, 5160-5172	3.2	3
20	Mispacking of the Phe87 Side Chain Reduces the Kinetic Stability of Human Transthyretin. <i>Biochemistry</i> , 2018 , 57, 6919-6922	3.2	3
19	Structure and Function of the CBP/p300 TAZ Domains 2005 , 114-120		2
18	Assignment of 1H, 13C, and 15N resonances of reduced Escherichia coli glutaredoxin 2. <i>Journal of Biomolecular NMR</i> , 1999 , 14, 197-8	3	2
17	Modeling of Hidden Structures Using Sparse Chemical Shift Data from NMR Relaxation Dispersion. <i>Biophysical Journal</i> , 2021 , 120, 296-305	2.9	2
16	Structural Basis for Graded Inhibition of CREB:DNA Interactions by Multisite Phosphorylation. <i>Biochemistry</i> , 2018 , 57, 6964-6972	3.2	2
15	Thioredoxin and Glutaredoxin 2007 ,		1
14	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
13	RNA Binding by the KTS Splice Variants of WilmsPTumor Suppressor Protein WT1. <i>Biochemistry</i> , 2020 , 59, 3889-3901	3.2	1
12	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	3.5	1
11	Role of Active Site Loop Dynamics in Mediating Ligand Release from Dihydrofolate Reductase. <i>Biochemistry</i> , 2021 , 60, 2663-2671	3.2	1
10	The molecular basis of allostery in a facilitated dissociation process. <i>Structure</i> , 2021 , 29, 1327-1338.e5	5.2	1

9	Classic Analysis of Biopolymer Dynamics Is Model Free. <i>Biophysical Journal</i> , 2016 , 110, 3-6	2.9	o
8	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	0.7	o
7	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	3.2	o
6	Chapter 5:NMR Studies of Disordered but Functional Proteins. <i>RSC Biomolecular Sciences</i> ,111-129		
5	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 Biophysical Society of Japan). <i>Seibutsu Butsuri</i> , 2010 , 50, S152		o
4	Nuclear magnetic resonance of thioredoxin and glutaredoxin. <i>Methods in Enzymology</i> , 1995 , 252, 293-306.	6.7	
3	Dynamics of I κ B α Probed by NMR. <i>FASEB Journal</i> , 2007 , 21, A655	0.9	
2	Structural Basis for Cooperative Transcription Factor Binding to the CBP Coactivator. <i>FASEB Journal</i> , 2012 , 26, lb266	0.9	
1	Early Strides in NMR Dynamics Measurements. <i>Biochemistry</i> , 2021 , 60, 3452-3454	3.2	