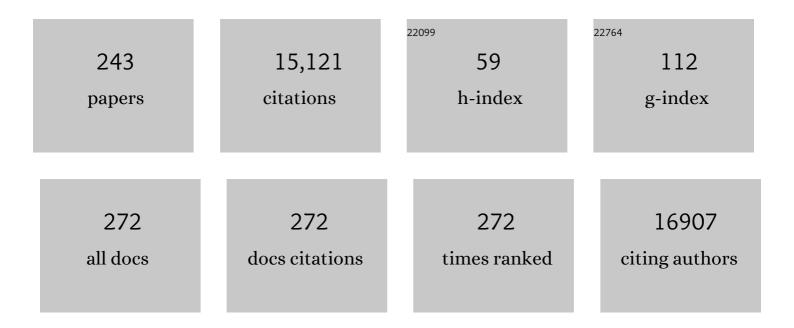
Rolf Boelens

List of Publications by Year in descending order

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POLE ROFLENS

#	Article	IF	CITATIONS
1	HADDOCK:Â A Proteinâ^Protein Docking Approach Based on Biochemical or Biophysical Information. Journal of the American Chemical Society, 2003, 125, 1731-1737.	6.6	2,642
2	Dynamic Readers for 5-(Hydroxy)Methylcytosine and Its Oxidized Derivatives. Cell, 2013, 152, 1146-1159.	13.5	888
3	New insights into the structure and composition of technical lignins: a comparative characterisation study. Green Chemistry, 2016, 18, 2651-2665.	4.6	648
4	Structure and Flexibility Adaptation in Nonspecific and Specific Protein-DNA Complexes. Science, 2004, 305, 386-389.	6.0	506
5	Hsp90-Tau Complex Reveals Molecular Basis for Specificity in Chaperone Action. Cell, 2014, 156, 963-974.	13.5	269
6	Altered flexibility in the substrate-binding site of related native and engineered high-alkaline Bacillus subtilisins 1 1Edited by P. E. Wright. Journal of Molecular Biology, 1999, 292, 111-123.	2.0	256
7	The DNA-binding domain of HIV-1 integrase has an SH3-like fold. Nature Structural and Molecular Biology, 1995, 2, 807-810.	3.6	242
8	Identification of a ubiquitin-protein ligase subunit within the CCR4-NOT transcription repressor complex. EMBO Journal, 2002, 21, 355-364.	3.5	186
9	Structure of Arc represser in solution: evidence for a family of β-sheet DMA-binding proteins. Nature, 1990, 346, 586-589.	13.7	180
10	WeNMR: Structural Biology on the Grid. Journal of Grid Computing, 2012, 10, 743-767.	2.5	170
11	Information-driven protein-DNA docking using HADDOCK: it is a matter of flexibility. Nucleic Acids Research, 2006, 34, 3317-3325.	6.5	169
12	Solution structure of the POU-specific DNA-binding domain of Oct-1. Nature, 1993, 362, 852-855.	13.7	161
13	Structural and dynamic changes of photoactive yellow protein during its photocycle in solution. Nature Structural Biology, 1998, 5, 568-570.	9.7	155
14	The nucleotide-binding site of bacterial translation initiation factor 2 (IF2) as a metabolic sensor. Proceedings of the National Academy of Sciences of the United States of America, 2006, 103, 13962-13967.	3.3	155
15	Identification of a diagnostic structural motif reveals a new reaction intermediate and condensation pathway in kraft lignin formation. Chemical Science, 2018, 9, 6348-6360.	3.7	143
16	Nucleosomal DNA binding drives the recognition of H3K36-methylated nucleosomes by the PSIP1-PWWP domain. Epigenetics and Chromatin, 2013, 6, 12.	1.8	141
17	Redox-Dependent Control of FOXO/DAF-16 by Transportin-1. Molecular Cell, 2013, 49, 730-742.	4.5	138
18	The solution structure of the amino-terminal HHCC domain of HIV-2 integrase: a three-helix bundle stabilized by zinc. Current Biology, 1997, 7, 739-746.	1.8	134

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19	Solution Structure and Backbone Dynamics of the Photoactive Yellow Proteinâ€,‡. Biochemistry, 1998, 37, 12689-12699.	1.2	129
20	An Off-resonance Rotating Frame Relaxation Experiment for the Investigation of Macromolecular Dynamics Using Adiabatic Rotations. Journal of Magnetic Resonance, 1998, 131, 351-357.	1.2	129
21	Gradient-enhanced HMQC and HSQC spectroscopy. Applications to 15N-labeled Mnt repressor. Journal of the American Chemical Society, 1991, 113, 9688-9690.	6.6	128
22	A comprehensive framework of E2–RING E3 interactions of the human ubiquitin–proteasome system. Molecular Systems Biology, 2009, 5, 295.	3.2	126
23	Data-driven docking for the study of biomolecular complexes. FEBS Journal, 2005, 272, 293-312.	2.2	125
24	Structural Insight into the Recognition of the H3K4me3 Mark by the TFIID Subunit TAF3. Structure, 2008, 16, 1245-1256.	1.6	123
25	Plasticity in protein-DNA recognition: lac repressor interacts with its natural operator O1 through alternative conformations of its DNA-binding domain. EMBO Journal, 2002, 21, 2866-2876.	3.5	117
26	The Structure of the Human ERCC1/XPF Interaction Domains Reveals a Complementary Role for the Two Proteins in Nucleotide Excision Repair. Structure, 2005, 13, 1849-1858.	1.6	116
27	Structural Model of the UbcH5B/CNOT4 Complex Revealed by Combining NMR, Mutagenesis, and Docking Approaches. Structure, 2004, 12, 633-644.	1.6	113
28	Solution Structure of a Chemosensory Protein from the Desert Locust Schistocerca gregaria,. Biochemistry, 2006, 45, 10606-10613.	1.2	111
29	The Solution Structure of the AppA BLUF Domain: Insight into the Mechanism of Light-Induced Signaling. ChemBioChem, 2006, 7, 187-193.	1.3	111
30	NMR analysis of protein interactions. Current Opinion in Chemical Biology, 2005, 9, 501-508.	2.8	109
31	N-terminal domain of human Hsp90 triggers binding to the cochaperone p23. Proceedings of the National Academy of Sciences of the United States of America, 2011, 108, 580-585.	3.3	109
32	On the Role of Aromatic Side Chains in the Photoactivation of BLUF Domains. Biochemistry, 2007, 46, 7405-7415.	1.2	106
33	E3 ligase Rad18 promotes monoubiquitination rather than ubiquitin chain formation by E2 enzyme Rad6. Proceedings of the National Academy of Sciences of the United States of America, 2011, 108, 5590-5595.	3.3	104
34	Solution Structure of the HU Protein fromBacillus stearothermophilus. Journal of Molecular Biology, 1995, 254, 692-703.	2.0	103
35	Hydration dynamics of the collagen triple helix by NMR11Edited by P. E. Wright. Journal of Molecular Biology, 2000, 300, 1041-1048.	2.0	103
36	Proteins Feel More Than They See: Fine-Tuning of Binding Affinity by Properties of the Non-Interacting Surface. Journal of Molecular Biology, 2014, 426, 2632-2652.	2.0	103

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37	Quantitative Measurement of Relaxation Interference Effects between1HNCSA and1Hâ^'15N Dipolar Interaction:Â Correlation with Secondary Structure. Journal of the American Chemical Society, 1997, 119, 8985-8990.	6.6	94
38	Mapping the fMet-tRNAfMet binding site of initiation factor IF2. EMBO Journal, 2000, 19, 5233-5240.	3.5	94
39	The solution structure of Lac repressor headpiece 62 complexed to a symmetrical lac operator. Structure, 1999, 7, 1483-S3.	1.6	84
40	Refined solution structure of the c-terminal DNA-binding domain of human immunovirus-1 integrase. Proteins: Structure, Function and Bioinformatics, 1999, 36, 556-564.	1.5	83
41	Conformational changes in the oligonucleotide duples d(GCGTTGCG). d (CGCAACGC) induced by formation of a cis-syn thymine dimer. A two-dimensional NMR study. FEBS Journal, 1987, 162, 37-43.	0.2	82
42	Probing the Nature of the Blue-Shifted Intermediate of Photoactive Yellow Protein in Solution by NMR:  Hydrogenâ^'Deuterium Exchange Data and pH Studies. Biochemistry, 2000, 39, 14392-14399.	1.2	81
43	The Structure of the C4C4RING Finger of Human NOT4 Reveals Features Distinct from Those of C3HC4 RING Fingers. Journal of Biological Chemistry, 2001, 276, 10185-10190.	1.6	80
44	High-resolution Structure of the Phosphorylated Form of the Histidine-containing Phosphocarrier Protein HPr fromEscherichia coliDetermined by Restrained Molecular Dynamics from NMR-NOE Data. Journal of Molecular Biology, 1995, 246, 180-193.	2.0	78
45	Formation of the hinge helix in the lac represser is induced upon binding to the lac operator. Nature Structural Biology, 1996, 3, 916-919.	9.7	77
46	Structural and Functional Analysis of the Kid Toxin Protein from E. coli Plasmid R1. Structure, 2002, 10, 1425-1433.	1.6	77
47	Rapid and simple approach for the NMR resonance assignment of the carbohydrate chains of an intact glycoprotein Application of gradient-enhanced natural abundance 1 H-13 C HSQC and HSQC-TOCSY to the α-subunit of human chorionic gonadotropin. FEBS Letters, 1994, 348, 1-6.	1.3	75
48	PhoE Signal Peptide Inserts into Micelles as a Dynamic Helix-Break-Helix Structure, Which Is Modulated by the Environment. A Two-Dimensional 1H NMR Study. Biochemistry, 1995, 34, 11617-11624.	1.2	75
49	NMR Studies of the Free alpha Subunit of Human Chorionic Gonadotropin. Structural Influences of N-Glycosylation and the beta Subunit on the Conformation of the alpha Subunit. FEBS Journal, 1996, 241, 229-242.	0.2	75
50	Light-Induced Flipping of a Conserved Glutamine Sidechain and Its Orientation in the AppA BLUF Domain. Journal of the American Chemical Society, 2006, 128, 15066-15067.	6.6	75
51	"Ensemble―iterative relaxation matrix approach: A new NMR refinement protocol applied to the solution structure of crambin. Proteins: Structure, Function and Bioinformatics, 1993, 15, 385-400.	1.5	74
52	Toward an Integrated Model of Proteinâ^'DNA Recognition as Inferred from NMR Studies on the Lac Repressor System. Chemical Reviews, 2004, 104, 3567-3586.	23.0	74
53	NMR structures of phospholipase A2 reveal conformational changes during interfacial activation. Nature Structural and Molecular Biology, 1995, 2, 402-406.	3.6	72
54	The Solution Structure of a Transient Photoreceptor Intermediate: Δ25 Photoactive Yellow Protein. Structure, 2005, 13, 953-962.	1.6	71

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55	Nuclear Magnetic Resonance Solution Structure of the Arc Repressor Using Relaxation Matrix Calculations. Journal of Molecular Biology, 1994, 236, 328-341.	2.0	69
56	Solution Structure of Dimeric Mnt Repressor (1-76). Biochemistry, 1994, 33, 15036-15045.	1.2	67
57	Conformational changes in phospholipase A2 upon binding to micellar interfaces in the absence and presence of competitive inhibitors. A proton and nitrogen-15 NMR study. Biochemistry, 1992, 31, 10024-10030.	1.2	66
58	Solution Structure of the Sequence-specific HMG Box of the Lymphocyte Transcriptional Activator Sox-4. Journal of Biological Chemistry, 1995, 270, 30516-30524.	1.6	65
59	Structural Properties of the Promiscuous VP16 Activation Domainâ€. Biochemistry, 2005, 44, 827-839.	1.2	63
60	Improved HSQC experiments for the observation of exchange broadened signals. Journal of Biomolecular NMR, 1996, 8, 223-8.	1.6	62
61	Crystal structure and collagen-binding site of immune inhibitory receptor LAIR-1: unexpected implications for collagen binding by platelet receptor GPVI. Blood, 2010, 115, 1364-1373.	0.6	62
62	Time- and ensemble-averaged direct NOE restraints. Journal of Biomolecular NMR, 1994, 4, 143-9.	1.6	60
63	Microsecond time scale dynamics in the RXR DNA-binding domain from a combination of spin-echo and off-resonance rotating frame relaxation measurements. Journal of Biomolecular NMR, 1999, 13, 275-288.	1.6	59
64	Function and Interactions of ERCC1-XPF in DNA Damage Response. Molecules, 2018, 23, 3205.	1.7	59
65	Critical Scaffolding Regions of the Tumor Suppressor Axin1 Are Natively Unfolded. Journal of Molecular Biology, 2011, 405, 773-786.	2.0	58
66	A novel µâ€conopeptide, CnIIIC, exerts potent and preferential inhibition of Na _V 1.2/1.4 channels and blocks neuronal nicotinic acetylcholine receptors. British Journal of Pharmacology, 2012, 166, 1654-1668.	2.7	55
67	Spatial arrangement of the three α helices in the solution conformation of E. coli lac represser DNA-binding domain. FEBS Letters, 1984, 174, 243-247.	1.3	54
68	Time-saving methods for heteronuclear multidimensional NMR of (13C, 15N) doubly labeled proteins. Journal of Biomolecular NMR, 1994, 4, 201-13.	1.6	54
69	The solution structure of serine protease PB92 from Bacillus alcalophilus presents a rigid fold with a flexible substrate-binding site. Structure, 1997, 5, 521-532.	1.6	53
70	Mobilities of the Inner Three Core Residues and the Man(α1→6) Branch of the Glycan at Asn78 of the α-Subunit of Human Chorionic Gonadotropin Are Restricted by the Protein. Biochemistry, 1998, 37, 1933-1940.	1.2	51
71	Hinge-helix formation and DNA bending in various lac repressor–operator complexes. EMBO Journal, 1999, 18, 6472-6480.	3.5	51
72	Interactions of Kid–Kis toxin–antitoxin complexes with the parD operator-promoter region of plasmid R1 are piloted by the Kis antitoxin and tuned by the stoichiometry of Kid–Kis oligomers. Nucleic Acids Research, 2007, 35, 1737-1749.	6.5	51

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73	MONTY: a Monte Carlo approach to protein-DNA recognition. Journal of Molecular Biology, 1994, 235, 318-324.	2.0	49
74	1H, 13C, and 15N resonance assignments and secondary structure analysis of the HU protein from Bacillus stearothermophilus using two- and three-dimensional double- and triple-resonance heteronuclear magnetic resonance spectroscopy. Biochemistry, 1994, 33, 14858-14870.	1.2	49
75	The tetramerization domain of the Mnt repressor consists of two right-handed coiled coils. Nature Structural Biology, 1999, 6, 755-759.	9.7	49
76	Sulforaphane inhibits pancreatic cancer through disrupting Hsp90–p50Cdc37 complex and direct interactions with amino acids residues of Hsp90. Journal of Nutritional Biochemistry, 2012, 23, 1617-1626.	1.9	49
77	Biophysical characterization of mutants of <i>Bacillus subtilis</i> lipase evolved for thermostability: Factors contributing to increased activity retention. Protein Science, 2012, 21, 487-497.	3.1	49
78	Determination of protein structures from nuclear magnetic resonance data using a restrained molecular dynamics approach: The lac repressor DNA binding domain. Biochimie, 1985, 67, 707-715.	1.3	48
79	Model for RNA Binding and the Catalytic Site of the RNase Kid of the Bacterial parD Toxin–Antitoxin System. Journal of Molecular Biology, 2006, 357, 115-126.	2.0	48
80	Specificity and Affinity of Lac Repressor for the Auxiliary Operators O2 and O3 Are Explained by the Structures of Their Protein–DNA Complexes. Journal of Molecular Biology, 2009, 390, 478-489.	2.0	46
81	Solution structure of the α-subunit of human chorionic gonadotropin. FEBS Journal, 1999, 260, 490-498.	0.2	45
82	Mutations in the glucocorticoid receptor DNA-binding domain mimic an allosteric effect of DNA 1 1Edited by P. E. Wright. Journal of Molecular Biology, 2000, 301, 947-958.	2.0	45
83	Effects of the N-Linked Glycans on the 3D Structure of the Free α-Subunit of Human Chorionic Gonadotropin. Biochemistry, 2000, 39, 6012-6021.	1.2	43
84	Cys-Ph-TAHA: a lanthanide binding tag for RDC and PCS enhanced protein NMR. Journal of Biomolecular NMR, 2011, 51, 329-337.	1.6	43
85	Structure, stability, and IgE binding of the peach allergen <scp>P</scp> eamaclein (Pru p 7). Biopolymers, 2014, 102, 416-425.	1.2	43
86	An EPR study of the photodissociation reactions of oxidised cytochrome c oxidase-nitric oxide complexes. Biochimica Et Biophysica Acta - Bioenergetics, 1983, 724, 176-183.	0.5	42
87	Symmetry and Asymmetry of the RING–RING Dimer of Rad18. Journal of Molecular Biology, 2011, 410, 424-435.	2.0	41
88	Computer-assisted assignment of 2D1H NMR spectra of proteins: Basic algorithms and application to phoratoxin B. Journal of Biomolecular NMR, 1991, 1, 23-47.	1.6	40
89	Identification of the ribosome binding sites of translation initiation factor IF3 by multidimensional heteronuclear NMR spectroscopy. Rna, 1999, 5, 82-92.	1.6	40
90	Combining NMR Relaxation with Chemical Shift Perturbation Data to Drive Protein–protein Docking. Journal of Biomolecular NMR, 2006, 34, 237-244.	1.6	39

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91	Analysis of the XPA and ssDNA-binding surfaces on the central domain of human ERCC1 reveals evidence for subfunctionalization. Nucleic Acids Research, 2007, 35, 5789-5798.	6.5	39
92	Solution Structure of the Immunodominant Region of Protein G of Bovine Respiratory Syncytial Virusâ€,â€j. Biochemistry, 1996, 35, 14684-14688.	1.2	38
93	Measurement of (15)N- (1)H coupling constants in uniformly (15)N-labeled proteins: Application to the photoactive yellow protein. Journal of Biomolecular NMR, 1997, 10, 301-306.	1.6	38
94	E2–c-Cbl Recognition Is Necessary but not Sufficient for Ubiquitination Activity. Journal of Molecular Biology, 2009, 385, 507-519.	2.0	37
95	Homonuclear three-dimensional proton NMR spectroscopy of pike parvalbumin. Comparison of short- and medium-range NOEs from 2D and 3D NMR. Journal of the American Chemical Society, 1990, 112, 5024-5030.	6.6	36
96	LexA repressor and iron uptake regulator from Escherichia coli: new members of the CAP-like DNA binding domain superfamily. Protein Engineering, Design and Selection, 1994, 7, 1449-1453.	1.0	36
97	DNA repair factor APLF acts as a H2A-H2B histone chaperone through binding its DNA interaction surface. Nucleic Acids Research, 2018, 46, 7138-7152.	6.5	36
98	An Altered-specificity Ubiquitin-conjugating Enzyme/Ubiquitin–Protein Ligase Pair. Journal of Molecular Biology, 2004, 337, 157-165.	2.0	35
99	Observation of intersubunit NOEs in a dimeric P22 Mnt repressor mutant by a time-shared [15N, 13C] double half-filter technique. Journal of Biomolecular NMR, 1993, 3, 709.	1.6	34
100	A model for the LexA repressor DNA complex. Proteins: Structure, Function and Bioinformatics, 1995, 21, 226-236.	1.5	34
101	Sequence-specific Recognition of DNA by the C-terminal Domain of Nucleoid-associated Protein H-NS. Journal of Biological Chemistry, 2009, 284, 30453-30462.	1.6	34
102	Identification, structural and pharmacological characterization of Ï,,-CnVA, a conopeptide that selectively interacts with somatostatin sst3 receptor. Biochemical Pharmacology, 2013, 85, 1663-1671.	2.0	34
103	Millisecond to Microsecond Time Scale Dynamics of the Retinoid X and Retinoic Acid Receptor DNA-Binding Domains and Dimeric Complex Formationâ€. Biochemistry, 1999, 38, 1951-1956.	1.2	33
104	Lack of Negative Charge in the E46Q Mutant of Photoactive Yellow Protein Prevents Partial Unfolding of the Blue-Shifted Intermediateâ€. Biochemistry, 2003, 42, 14501-14506.	1.2	33
105	<i>parD</i> toxin–antitoxin system of plasmid R1 – basic contributions, biotechnological applications and relationships with closelyâ€related toxin–antitoxin systems. FEBS Journal, 2010, 277, 3097-3117.	2.2	32
106	A Naturalâ€Product Switch for a Dynamic Protein Interface. Angewandte Chemie - International Edition, 2014, 53, 6443-6448.	7.2	32
107	Solution Structure of the Ubiquitin-conjugating Enzyme UbcH5B. Journal of Molecular Biology, 2004, 344, 513-526.	2.0	31
108	Structure and Function of Bacterial Kid-Kis and Related Toxin-Antitoxin Systems. Protein and Peptide Letters, 2007, 14, 113-124.	0.4	31

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109	Axin cancer mutants form nanoaggregates to rewire the Wnt signaling network. Nature Structural and Molecular Biology, 2016, 23, 324-332.	3.6	31
110	A residue-specific view of the association and dissociation pathway in protein–DNA recognition. Nature Structural Biology, 2002, 9, 193-7.	9.7	30
111	Two-dimensional NMR studies on des-pentapeptide-insulin. Proton resonance assignments and secondary structure analysis. FEBS Journal, 1990, 191, 147-153.	0.2	29
112	Porcine pancreatic phospholipase A2: sequence-specific proton and nitrogen-15 NMR assignments and secondary structure. Biochemistry, 1991, 30, 3135-3146.	1.2	29
113	1H, 13C and 15N NMR backbone assignments of the 269-residue serine protease PB92 from Bacillus alcalophilus. Journal of Biomolecular NMR, 1994, 4, 123-8.	1.6	29
114	Monte Carlo docking of protein-DNA complexes: incorporation of DNA flexibility and experimental data. Protein Engineering, Design and Selection, 1994, 7, 761-768.	1.0	29
115	Interactions between the toxin Kid of the bacterial parD system and the antitoxins Kis and MazE. Proteins: Structure, Function and Bioinformatics, 2007, 67, 219-231.	1.5	29
116	Structural Dynamics in the Activation of Epac. Journal of Biological Chemistry, 2008, 283, 6501-6508.	1.6	29
117	Assignment of the 1H-NMR spectrum of a lac repressor headpiece-operator complex in H2O and identification of NOEs. Consequences for protein-DNA interaction. FEBS Journal, 1990, 194, 629-637.	0.2	28
118	Solution Structure of the C-terminal Domain of TFIIH P44 Subunit Reveals a Novel Type of C4C4 Ring Domain Involved in Protein-Protein Interactions. Journal of Biological Chemistry, 2005, 280, 20785-20792.	1.6	28
119	High-level expression of biologically active glycoprotein hormones in Pichia pastoris strains—selection of strain GS115, and not X-33, for the production of biologically active N-glycosylated 15N-labeled phCG. Glycoconjugate Journal, 2008, 25, 245-257.	1.4	28
120	The Structure of the XPF-ssDNA Complex Underscores the Distinct Roles of the XPF and ERCC1 Helix- Hairpin-Helix Domains in ss/ds DNA Recognition. Structure, 2012, 20, 667-675.	1.6	28
121	<i>E. coli</i> MG1655 modulates its phospholipid composition through the cell cycle. FEBS Letters, 2015, 589, 2726-2730.	1.3	28
122	The cytochrome c oxidase-azide-nitric oxide complex as a model for the oxygen-binding site. Biochimica Et Biophysica Acta - Bioenergetics, 1984, 765, 196-209.	0.5	27
123	Rapid acquisition of three-dimensional triple-resonance experiments using pulsed field gradient techniques. Journal of Biomolecular NMR, 1992, 2, 395-400.	1.6	27
124	Gradient-enhanced 3D NOESY-HMQC spectroscopy. Journal of Biomolecular NMR, 1992, 2, 301-305.	1.6	27
125	Structure and dynamics of the DNA binding protein HU from Bacillus stearothermophilus by NMR spectroscopy. , 1996, 40, 553-559.		27
126	Refined solution structure of the dimeric N-terminal HHCC domain of HIV-2 integrase. Journal of Biomolecular NMR, 2000, 18, 119-128.	1.6	27

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127	Modeling Proteinâ^'Protein Complexes Involved in the CytochromecOxidase Copper-Delivery Pathway. Journal of Proteome Research, 2007, 6, 1530-1539.	1.8	27
128	The solution structure of a monomeric insulin. A two-dimensional 1H-NMR study of des-(B26-B30)-insulin in combination with distance geometry and restrained molecular dynamics. FEBS Journal, 1991, 202, 447-458.	0.2	25
129	Use of very long-distance NOEs in a fully deuterated protein: an approach for rapid protein fold determination. Journal of Magnetic Resonance, 2003, 163, 228-235.	1.2	25
130	3D DOSY-TROSY to determine the translational diffusion coefficient of large protein complexes. Protein Engineering, Design and Selection, 2011, 24, 99-103.	1.0	25
131	Glycosylation of Conotoxins. Marine Drugs, 2013, 11, 623-642.	2.2	25
132	Conformational Plasticity of the POTRA 5 Domain in the Outer Membrane Protein Assembly Factor BamA. Structure, 2015, 23, 1317-1324.	1.6	25
133	Applications of two-dimensional 1H nuclear magnetic resonance methods in photochemically induced dynamic nuclear polarisation spectroscopy. Faraday Discussions of the Chemical Society, 1984, 78, 245.	2.2	24
134	Title is missing!. Journal of Biomolecular NMR, 1998, 11, 265-277.	1.6	24
135	Identification of the Single-stranded DNA Binding Surface of the Transcriptional Coactivator PC4 by NMR. Journal of Biological Chemistry, 1999, 274, 3693-3699.	1.6	24
136	SAMPLEX: Automatic mapping of perturbed and unperturbed regions of proteins and complexes. BMC Bioinformatics, 2010, 11, 51.	1.2	24
137	Structural Dynamics of Bacterial Translation Initiation Factor IF2. Journal of Biological Chemistry, 2012, 287, 10922-10932.	1.6	24
138	Temperature dependence of the magnetic volume susceptibility of human breast fat tissue: an NMR study. Magnetic Resonance Materials in Physics, Biology, and Medicine, 2012, 25, 33-39.	1.1	24
139	Two-dimensional 1H-NMR studies of phospholipase-A2-inhibitor complexes bound to a micellar lipid-water interface. FEBS Journal, 1991, 199, 601-607.	0.2	23
140	Rapid protein fold determination using secondary chemical shifts and cross-hydrogen bond 15N-13C' scalar couplings (3hbJNC'). Journal of Biomolecular NMR, 2001, 21, 221-233.	1.6	23
141	X-ray absorption spectroscopic studies of zinc in the N-terminal domain of HIV-2 integrase and model compounds. Journal of Synchrotron Radiation, 2003, 10, 86-95.	1.0	23
142	Solution structure of the C1-subdomain ofBacillus stearothermophilustranslation initiation factor IF2. Protein Science, 2005, 14, 2461-2468.	3.1	23
143	Binding Hotspots of BAZ2B Bromodomain: Histone Interaction Revealed by Solution NMR Driven Docking. Biochemistry, 2014, 53, 6706-6716.	1.2	23
144	Hydrogen exchange studies of the Arc repressor: Evidence for a monomeric folding intermediate. Biopolymers, 1995, 35, 217-226.	1.2	22

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145	Dynamics and Metal Exchange Properties of C4C4 RING Domains from CNOT4 and the p44 Subunit of TFIIH. Journal of Molecular Biology, 2005, 349, 621-637.	2.0	21
146	The Solution Structure of DNA-free Pax-8 Paired Box Domain Accounts for Redox Regulation of Transcriptional Activity in the Pax Protein Family. Journal of Biological Chemistry, 2008, 283, 33321-33328.	1.6	21
147	A comprehensive framework of E2–RING E3 interactions of the human ubiquitin–proteasome system. Molecular Systems Biology, 2009, 5, .	3.2	21
148	Structure of the Oâ€Glycosylated Conopeptide CcTx from <i>Conus consors</i> Venom. Chemistry - A European Journal, 2013, 19, 870-879.	1.7	21
149	Structure refinement of the glucocorticoid receptor-DNA binding domain from NMR data by relaxation matrix calculations. Journal of Molecular Biology, 1995, 247, 689-700.	2.0	20
150	The Solution Structure and Dynamics of an Arc Repressor Mutant Reveal Premelting Conformational Changes Related to DNA Bindingâ€. Biochemistry, 1999, 38, 6035-6042.	1.2	20
151	The Intrinsically Unstructured Domain of PC4 Modulates the Activity of the Structured Core through Inter- and Intramolecular Interactions. Biochemistry, 2006, 45, 5067-5081.	1.2	20
152	Quantitative use of chemical shifts for the modeling of protein complexes. Proteins: Structure, Function and Bioinformatics, 2011, 79, 2662-2670.	1.5	20
153	Solution Structure of Porcine Pancreatic Procolipase as Determined from 1H Homonuclear Two-Dimensional and Three-Dimensional NMR. FEBS Journal, 1995, 227, 663-672.	0.2	20
154	The use of two-dimensional nuclear-magnetic-resonance spectroscopy and two-dimensional difference spectra in the elucidation of the active center of Megasphaera elsdenii flavodoxin. FEBS Journal, 1984, 141, 323-330.	0.2	19
155	The HhH domain of the human DNA repair protein XPF forms stable homodimers. Proteins: Structure, Function and Bioinformatics, 2008, 70, 1551-1563.	1.5	19
156	Structural basis of nucleic acid binding by <i>Nicotiana tabacum</i> glycine-rich RNA-binding protein: implications for its RNA chaperone function. Nucleic Acids Research, 2014, 42, 8705-8718.	6.5	19
157	Insight into the conformational stability of membrane-embedded BamA using a combined solution and solid-state NMR approach. Journal of Biomolecular NMR, 2015, 61, 321-332.	1.6	19
158	Ligand-binding effects on the kringle 4 domain from human plasminogen: a study by laser photo-CIDNP1H-NMR spectroscopy. BBA - Proteins and Proteomics, 1989, 994, 121-137.	2.1	18
159	Editing of Chemical Exchange-Relayed NOEs in NMR Experiments for the Observation of Protein–Water Interactions. Journal of Magnetic Resonance, 1999, 136, 214-218.	1.2	18
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