

Peter Gntert

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

240 papers	19,505 citations	62 h-index	136 g-index
250 ext. papers	21,003 ext. citations	7.1 avg, IF	6.81 L-index

#	Paper	IF	Citations
240	Atomic-resolution chemical characterization of (2x)72-kDa tryptophan synthase via four- and five-dimensional H-detected solid-state NMR.. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022 , 119,	11.5	6
239	Optimization and validation of multi-state NMR protein structures using structural correlations.. <i>Journal of Biomolecular NMR</i> , 2022 , 1	3	0
238	A B-factor for NOEs?. <i>Journal of Magnetic Resonance</i> , 2022 , 338, 107189	3	0
237	H, C and N resonance assignments and solution structures of the two RRM domains of Matrin-3. <i>Biomolecular NMR Assignments</i> , 2021 , 1	0.7	
236	Paramagnetic Solid-State NMR to Localize the Metal-Ion Cofactor in an Oligomeric DnaB Helicase. <i>Chemistry - A European Journal</i> , 2021 , 27, 7745-7755	4.8	4
235	Evaluation of Multi-Objective Optimization Algorithms for NMR Chemical Shift Assignment. <i>Molecules</i> , 2021 , 26,	4.8	1
234	H, C and N resonance assignment of the YTH domain of YTHDC2. <i>Biomolecular NMR Assignments</i> , 2021 , 15, 1-7	0.7	0
233	An automated iterative approach for protein structure refinement using pseudocontact shifts. <i>Journal of Biomolecular NMR</i> , 2021 , 75, 319-334	3	2
232	Automated Backbone NMR Resonance Assignment of Large Proteins Using Redundant Linking from a Single Simultaneous Acquisition. <i>Journal of the American Chemical Society</i> , 2020 , 142, 5793-5799	16.4	22
231	Automated assignment of methyl NMR spectra from large proteins. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2020 , 118-119, 54-73	10.4	13
230	Backbone and methyl assignment of bacteriorhodopsin incorporated into nanodiscs. <i>Journal of Biomolecular NMR</i> , 2020 , 74, 45-60	3	7
229	The three-dimensional structure of human β -endorphin amyloid fibrils. <i>Nature Structural and Molecular Biology</i> , 2020 , 27, 1178-1184	17.6	18
228	Protein Allostery at Atomic Resolution. <i>Angewandte Chemie - International Edition</i> , 2020 , 59, 22132-22136	16.4	6
227	Protein Allostery at Atomic Resolution. <i>Angewandte Chemie</i> , 2020 , 132, 22316-22323	3.6	
226	Dynamics of Bacteriorhodopsin in the Dark-Adapted State from Solution Nuclear Magnetic Resonance Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2020 , 59, 20965-20972	16.4	1
225	Dynamics of Bacteriorhodopsin in the Dark-Adapted State from Solution Nuclear Magnetic Resonance Spectroscopy. <i>Angewandte Chemie</i> , 2020 , 132, 21151-21158	3.6	1
224	An atypical LIR motif within UBA5 (ubiquitin like modifier activating enzyme 5) interacts with GABARAP proteins and mediates membrane localization of UBA5. <i>Autophagy</i> , 2020 , 16, 256-270	10.2	16

223	The Solution Structure and Dynamics of Cd-Metallothionein from Reveal Optimization for Binding Cd over Zn. <i>Biochemistry</i> , 2019 , 58, 4570-4581	3.2	10
222	Integrated NMR and cryo-EM atomic-resolution structure determination of a half-megadalton enzyme complex. <i>Nature Communications</i> , 2019 , 10, 2697	17.4	46
221	Protein Structure Determination in Living Cells. <i>International Journal of Molecular Sciences</i> , 2019 , 20,	6.3	12
220	Global response of diacylglycerol kinase towards substrate binding observed by 2D and 3D MAS NMR. <i>Scientific Reports</i> , 2019 , 9, 3995	4.9	8
219	High-Resolution Protein 3D Structure Determination in Living Eukaryotic Cells. <i>Angewandte Chemie - International Edition</i> , 2019 , 58, 7284-7288	16.4	34
218	High-Resolution Protein 3D Structure Determination in Living Eukaryotic Cells. <i>Angewandte Chemie</i> , 2019 , 131, 7362-7366	3.6	5
217	Including Protons in Solid-State NMR Resonance Assignment and Secondary Structure Analysis: The Example of RNA Polymerase II Subunits Rpo4/7. <i>Frontiers in Molecular Biosciences</i> , 2019 , 6, 100	5.6	12
216	Automatic structure-based NMR methyl resonance assignment in large proteins. <i>Nature Communications</i> , 2019 , 10, 4922	17.4	18
215	CYANA 2019 , 1-2		
214	Protein aggregation of the p63 transcription factor underlies severe skin fragility in AEC syndrome. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018 , 115, E906-E915	11.5	19
213	Chain Assembly and Disassembly Processes Differently Affect the Conformational Space of Ubiquitin Chains. <i>Structure</i> , 2018 , 26, 249-258.e4	5.2	11
212	Noise peak filtering in multi-dimensional NMR spectra using convolutional neural networks. <i>Bioinformatics</i> , 2018 , 34, 4300-4301	7.2	12
211	Extending the Applicability of Exact Nuclear Overhauser Enhancements to Large Proteins and RNA. <i>ChemBioChem</i> , 2018 , 19, 1695	3.8	9
210	High-resolution small RNA structures from exact nuclear Overhauser enhancement measurements without additional restraints. <i>Communications Biology</i> , 2018 , 1, 61	6.7	12
209	Structural investigation of glycan recognition by the ERAD quality control lectin Yos9. <i>Journal of Biomolecular NMR</i> , 2018 , 72, 1-10	3	4
208	Automated Structure Determination from NMR Spectra 2018 , 401-418		
207	Fully automated assignment of methyl resonances of a 36 kDa protein dimer from sparse NOESY data. <i>Journal of Physics: Conference Series</i> , 2018 , 1036, 012008	0.3	1
206	Structure and dynamics conspire in the evolution of affinity between intrinsically disordered proteins. <i>Science Advances</i> , 2018 , 4, eaau4130	14.3	20

205	Proteome-wide analysis of phospho-regulated PDZ domain interactions. <i>Molecular Systems Biology</i> , 2018 , 14, e8129	12.2	32
204	Regulation of the Activity in the p53 Family Depends on the Organization of the Transactivation Domain. <i>Structure</i> , 2018 , 26, 1091-1100.e4	5.2	10
203	Peak picking multidimensional NMR spectra with the contour geometry based algorithm CYPICK. <i>Journal of Biomolecular NMR</i> , 2017 , 67, 63-76	3	21
202	NMR-based automated protein structure determination. <i>Archives of Biochemistry and Biophysics</i> , 2017 , 628, 24-32	4.1	27
201	Solution structure of discoidal high-density lipoprotein particles with a shortened apolipoprotein A-I. <i>Nature Structural and Molecular Biology</i> , 2017 , 24, 187-193	17.6	85
200	Determination of helix orientations in a flexible DNA by multi-frequency EPR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 29801-29811	3.6	14
199	Solution structure of the first RNA recognition motif domain of human spliceosomal protein SF3b49 and its mode of interaction with a SF3b145 fragment. <i>Protein Science</i> , 2017 , 26, 280-291	6.3	5
198	Structural and functional dissection of the DH and PH domains of oncogenic Bcr-Abl tyrosine kinase. <i>Nature Communications</i> , 2017 , 8, 2101	17.4	21
197	The Exact Nuclear Overhauser Enhancement: Recent Advances. <i>Molecules</i> , 2017 , 22,	4.8	13
196	Automated Structure Determination from NMR Spectra 2017 , 1-18		1
195	Atomic-resolution structure of a disease-relevant A β (1-42) amyloid fibril. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016 , 113, E4976-84	11.5	557
194	NMR Investigation of Structures of G-protein Coupled Receptor Folding Intermediates. <i>Journal of Biological Chemistry</i> , 2016 , 291, 27170-27186	5.4	6
193	The CUE Domain of Cue1 Aligns Growing Ubiquitin Chains with Ubc7 for Rapid Elongation. <i>Molecular Cell</i> , 2016 , 62, 918-928	17.6	25
192	NMR-Based Determination of the 3D Structure of the Ligand-Protein Interaction Site without Protein Resonance Assignment. <i>Journal of the American Chemical Society</i> , 2016 , 138, 4393-400	16.4	38
191	The Exact NOE as an Alternative in Ensemble Structure Determination. <i>Biophysical Journal</i> , 2016 , 110, 113-26	2.9	30
190	Solid-state NMR sequential assignments of the N-terminal domain of HpDnaB helicase. <i>Biomolecular NMR Assignments</i> , 2016 , 10, 13-23	0.7	13
189	Protein NMR Structure Refinement based on Bayesian Inference. <i>Journal of Physics: Conference Series</i> , 2016 , 699, 012005	0.3	8
188	Improved in-cell structure determination of proteins at near-physiological concentration. <i>Scientific Reports</i> , 2016 , 6, 38312	4.9	29

187	Solid-state NMR sequential assignment of an Amyloid- β (1-42) fibril polymorph. <i>Biomolecular NMR Assignments</i> , 2016 , 10, 269-76	0.7	17
186	Mechanism of TAp73 inhibition by Bp63 and structural basis of p63/p73 hetero-tetramerization. <i>Cell Death and Differentiation</i> , 2016 , 23, 1930-1940	12.7	22
185	Increased reliability of nuclear magnetic resonance protein structures by consensus structure bundles. <i>Structure</i> , 2015 , 23, 425-34	5.2	20
184	The RING domain of human promyelocytic leukemia protein (PML). <i>Journal of Biomolecular NMR</i> , 2015 , 61, 173-80	3	4
183	Systematic evaluation of combined automated NOE assignment and structure calculation with CYANA. <i>Journal of Biomolecular NMR</i> , 2015 , 62, 81-95	3	23
182	Combined automated NOE assignment and structure calculation with CYANA. <i>Journal of Biomolecular NMR</i> , 2015 , 62, 453-71	3	202
181	Complementarity and congruence between exact NOEs and traditional NMR probes for spatial decoding of protein dynamics. <i>Journal of Structural Biology</i> , 2015 , 191, 306-17	3.4	14
180	NMR structure calculation for all small molecule ligands and non-standard residues from the PDB Chemical Component Dictionary. <i>Journal of Biomolecular NMR</i> , 2015 , 63, 21-37	3	13
179	Atomic-resolution three-dimensional structure of amyloid β -fibrils bearing the Osaka mutation. <i>Angewandte Chemie - International Edition</i> , 2015 , 54, 331-5	16.4	211
178	Compiled data set of exact NOE distance limits, residual dipolar couplings and scalar couplings for the protein GB3. <i>Data in Brief</i> , 2015 , 5, 99-106	1.2	10
177	Die atomare dreidimensionale Struktur von Amyloid- β -Fibrillen mit der Osaka-Mutation. <i>Angewandte Chemie</i> , 2015 , 127, 337-342	3.6	12
176	A Structural Ensemble for the Enzyme Cyclophilin Reveals an Orchestrated Mode of Action at Atomic Resolution. <i>Angewandte Chemie - International Edition</i> , 2015 , 54, 11657-61	16.4	23
175	A Structural Ensemble for the Enzyme Cyclophilin Reveals an Orchestrated Mode of Action at Atomic Resolution. <i>Angewandte Chemie</i> , 2015 , 127, 11823-11827	3.6	1
174	NMR Exchange Format: a unified and open standard for representation of NMR restraint data. <i>Nature Structural and Molecular Biology</i> , 2015 , 22, 433-4	17.6	26
173	Evaluation of the reliability of the maximum entropy method for reconstructing 3D and 4D NOESY-type NMR spectra of proteins. <i>Biochemical and Biophysical Research Communications</i> , 2015 , 457, 200-5	3.4	4
172	Automated structure determination from NMR spectra. <i>Methods in Molecular Biology</i> , 2015 , 1261, 303-29.4	29.4	9
171	Peak picking NMR spectral data using non-negative matrix factorization. <i>BMC Bioinformatics</i> , 2014 , 15, 46	3.6	13
170	Automated NMR resonance assignment strategy for RNA via the phosphodiester backbone based on high-dimensional through-bond APSY experiments. <i>Journal of Biomolecular NMR</i> , 2014 , 59, 87-93	3	15

169	Towards a true protein movie: a perspective on the potential impact of the ensemble-based structure determination using exact NOEs. <i>Journal of Magnetic Resonance</i> , 2014 , 241, 53-9	3	27
168	Automated resonance assignment of the 21kDa stereo-array isotope labeled thioldisulfide oxidoreductase DsbA. <i>Journal of Magnetic Resonance</i> , 2014 , 249, 88-93	3	7
167	Spontaneous self-assembly of engineered armadillo repeat protein fragments into a folded structure. <i>Structure</i> , 2014 , 22, 985-95	5.2	18
166	Structural features of peptoid-peptide hybrids in lipid-water interfaces. <i>FEBS Letters</i> , 2014 , 588, 3291-7	3.8	7
165	RBFOX and SUP-12 sandwich a G base to cooperatively regulate tissue-specific splicing. <i>Nature Structural and Molecular Biology</i> , 2014 , 21, 778-86	17.6	24
164	Influence of NMR Data Completeness on Structure Determinations of Homodimeric Proteins. <i>Journal of the Chinese Chemical Society</i> , 2014 , 61, 1297-1306	1.5	
163	Identification of residues required for stalled-ribosome rescue in the codon-independent release factor YaeJ. <i>Nucleic Acids Research</i> , 2014 , 42, 3152-63	20.1	28
162	Novel RNA recognition motif domain in the cytoplasmic polyadenylation element binding protein 3. <i>Proteins: Structure, Function and Bioinformatics</i> , 2014 , 82, 2879-86	4.2	2
161	Crystal structure of a PCP/Sfp complex reveals the structural basis for carrier protein posttranslational modification. <i>Chemistry and Biology</i> , 2014 , 21, 552-562		30
160	Multiple-state ensemble structure determination from eNOE spectroscopy. <i>Molecular Physics</i> , 2013 , 111, 437-454	1.7	21
159	Prediction of peak overlap in NMR spectra. <i>Journal of Biomolecular NMR</i> , 2013 , 56, 113-23	3	
158	Peakmatch: a simple and robust method for peak list matching. <i>Journal of Biomolecular NMR</i> , 2013 , 55, 267-77	3	7
157	(1)H, (13)C, and (15)N resonance assignments of the dsRBDs of mouse RNA helicase A. <i>Biomolecular NMR Assignments</i> , 2013 , 7, 69-72	0.7	1
156	Reliability of exclusively NOESY-based automated resonance assignment and structure determination of proteins. <i>Journal of Biomolecular NMR</i> , 2013 , 57, 193-204	3	23
155	Stereospecific assignments in proteins using exact NOEs. <i>Journal of Biomolecular NMR</i> , 2013 , 57, 211-8	3	14
154	Estimating structure quality trends in the Protein Data Bank by equivalent resolution. <i>Computational Biology and Chemistry</i> , 2013 , 46, 8-15	3.6	11
153	Recommendations of the wwPDB NMR Validation Task Force. <i>Structure</i> , 2013 , 21, 1563-70	5.2	117
152	Automated solid-state NMR resonance assignment of protein microcrystals and amyloids. <i>Journal of Biomolecular NMR</i> , 2013 , 56, 243-54	3	32

151	Molecular crowding drives active Pin1 into nonspecific complexes with endogenous proteins prior to substrate recognition. <i>Journal of the American Chemical Society</i> , 2013 , 135, 13796-803	16.4	60
150	Structural basis for phosphorylation-triggered autophagic clearance of Salmonella. <i>Biochemical Journal</i> , 2013 , 454, 459-66	3.8	71
149	Automated and assisted RNA resonance assignment using NMR chemical shift statistics. <i>Nucleic Acids Research</i> , 2013 , 41, e172	20.1	36
148	NMR solution structure of a Chymotrypsin inhibitor from the Taiwan cobra <i>Naja naja atra</i> . <i>Molecules</i> , 2013 , 18, 8906-18	4.8	4
147	Effects of NMR spectral resolution on protein structure calculation. <i>PLoS ONE</i> , 2013 , 8, e68567	3.7	7
146	Blind testing of routine, fully automated determination of protein structures from NMR data. <i>Structure</i> , 2012 , 20, 227-36	5.2	64
145	Requirements on paramagnetic relaxation enhancement data for membrane protein structure determination by NMR. <i>Structure</i> , 2012 , 20, 1019-27	5.2	33
144	The First structure of a lantibiotic immunity protein, SpaI from <i>Bacillus subtilis</i> , reveals a novel fold. <i>Journal of Biological Chemistry</i> , 2012 , 287, 35286-35298	5.4	19
143	WeNMR: Structural Biology on the Grid. <i>Journal of Grid Computing</i> , 2012 , 10, 743-767	4.2	142
142	Structural insight into the interaction of ADP-ribose with the PARP WWE domains. <i>FEBS Letters</i> , 2012 , 586, 3858-64	3.8	32
141	Discrete three-dimensional representation of macromolecular motion from eNOE-based ensemble calculation. <i>Chimia</i> , 2012 , 66, 787-90	1.3	8
140	Solution structure and siRNA-mediated knockdown analysis of the mitochondrial disease-related protein C12orf65. <i>Proteins: Structure, Function and Bioinformatics</i> , 2012 , 80, 2629-42	4.2	26
139	Spatial elucidation of motion in proteins by ensemble-based structure calculation using exact NOEs. <i>Nature Structural and Molecular Biology</i> , 2012 , 19, 1053-7	17.6	73
138	Influence of ¹ H chemical shift assignments of the interface residues on structure determinations of homodimeric proteins. <i>Journal of Magnetic Resonance</i> , 2012 , 222, 96-104	3	6
137	A new algorithm for reliable and general NMR resonance assignment. <i>Journal of the American Chemical Society</i> , 2012 , 134, 12817-29	16.4	114
136	Characterization of molecular interactions between ACP and halogenase domains in the Curacin A polyketide synthase. <i>ACS Chemical Biology</i> , 2012 , 7, 378-86	4.9	29
135	Protein structure validation by generalized linear model root-mean-square deviation prediction. <i>Protein Science</i> , 2012 , 21, 229-38	6.3	38
134	Solution structure of the splicing factor motif of the human Prp18 protein. <i>Proteins: Structure, Function and Bioinformatics</i> , 2012 , 80, 968-74	4.2	3

133	Solution structures of the double-stranded RNA-binding domains from RNA helicase A. <i>Proteins: Structure, Function and Bioinformatics</i> , 2012 , 80, 1699-706	4.2	5
132	Fast automated NMR spectroscopy of short-lived biological samples. <i>ChemBioChem</i> , 2012 , 13, 964-7	3.8	2
131	A universal expression tag for structural and functional studies of proteins. <i>ChemBioChem</i> , 2012 , 13, 959-63	3.8	29
130	Simultaneous single-structure and bundle representation of protein NMR structures in torsion angle space. <i>Journal of Biomolecular NMR</i> , 2012 , 52, 351-64	3	19
129	Comparative NMR analysis of an 80-residue G protein-coupled receptor fragment in two membrane mimetic environments. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2011 , 1808, 2674-84	3.8	11
128	Characterization of the interaction of GABARAPL-1 with the LIR motif of NBR1. <i>Journal of Molecular Biology</i> , 2011 , 410, 477-87	6.5	68
127	Structural insights into Rcs phosphotransfer: the newly identified RcsD-ABL domain enhances interaction with the response regulator RcsB. <i>Structure</i> , 2011 , 19, 577-87	5.2	14
126	Structure-guided fragment-based in silico drug design of dengue protease inhibitors. <i>Journal of Computer-Aided Molecular Design</i> , 2011 , 25, 263-74	4.2	46
125	Optimization of amino acid type-specific ¹³ C and ¹⁵ N labeling for the backbone assignment of membrane proteins by solution- and solid-state NMR with the UPLABEL algorithm. <i>Journal of Biomolecular NMR</i> , 2011 , 49, 75-84	3	40
124	Exclusively NOESY-based automated NMR assignment and structure determination of proteins. <i>Journal of Biomolecular NMR</i> , 2011 , 50, 137-46	3	25
123	Objective identification of residue ranges for the superposition of protein structures. <i>BMC Bioinformatics</i> , 2011 , 12, 170	3.6	36
122	Structures of the first and second double-stranded RNA-binding domains of human TAR RNA-binding protein. <i>Protein Science</i> , 2011 , 20, 118-30	6.3	45
121	Calculation of Structures from NMR Restraints 2011 , 159-192		
120	Solution NMR Structure of Proteorhodopsin. <i>Angewandte Chemie</i> , 2011 , 123, 12148-12152	3.6	15
119	Solution NMR structure of proteorhodopsin. <i>Angewandte Chemie - International Edition</i> , 2011 , 50, 11942-6	16.4	150
118	Structural basis for the dual RNA-recognition modes of human Tra2-IRRM. <i>Nucleic Acids Research</i> , 2011 , 39, 1538-53	20.1	54
117	NMR protein structure determination in living E. coli cells using nonlinear sampling. <i>Nature Protocols</i> , 2010 , 5, 1051-60	18.8	35
116	Structural investigation of the C-terminal catalytic fragment of presenilin 1. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010 , 107, 9644-9	11.5	68

115	Solution structure of the catalytic domain of the mitochondrial protein ICT1 that is essential for cell vitality. <i>Journal of Molecular Biology</i> , 2010 , 404, 260-73	6.5	44
114	Solution structure of polytheonamide B, a highly cytotoxic nonribosomal polypeptide from marine sponge. <i>Journal of the American Chemical Society</i> , 2010 , 132, 12941-5	16.4	71
113	Determinants of activity in glutaredoxins: an in vitro evolved Grx1-like variant of Escherichia coli Grx3. <i>Biochemical Journal</i> , 2010 , 430, 487-95	3.8	7
112	Structural insight into the zinc finger CW domain as a histone modification reader. <i>Structure</i> , 2010 , 18, 1127-39	5.2	93
111	Structural basis for the sequence-specific RNA-recognition mechanism of human CUG-BP1 RRM3. <i>Nucleic Acids Research</i> , 2009 , 37, 5151-66	20.1	62
110	Automated assignment of NMR chemical shifts using peak-particle dynamics simulation with the DYNASSIGN algorithm. <i>Journal of Biomolecular NMR</i> , 2009 , 43, 97-109	3	24
109	Automated NMR structure determination of stereo-array isotope labeled ubiquitin from minimal sets of spectra using the SAIL-FLYA system. <i>Journal of Biomolecular NMR</i> , 2009 , 44, 261-72	3	25
108	Automated structure determination from NMR spectra. <i>European Biophysics Journal</i> , 2009 , 38, 129-43	1.9	189
107	Solution structure of the GUCT domain from human RNA helicase II/Gu beta reveals the RRM fold, but implausible RNA interactions. <i>Proteins: Structure, Function and Bioinformatics</i> , 2009 , 74, 133-44	4.2	9
106	Structure of the Cdt1 C-terminal domain: conservation of the winged helix fold in replication licensing factors. <i>Protein Science</i> , 2009 , 18, 2252-64	6.3	28
105	Solution structure of the cysteine-rich domain in Fn14, a member of the tumor necrosis factor receptor superfamily. <i>Protein Science</i> , 2009 , 18, 650-6	6.3	24
104	Protein structure determination in living cells by in-cell NMR spectroscopy. <i>Nature</i> , 2009 , 458, 102-5	50.4	267
103	Conformational stability and activity of p73 require a second helix in the tetramerization domain. <i>Cell Death and Differentiation</i> , 2009 , 16, 1582-9	12.7	46
102	CASD-NMR: critical assessment of automated structure determination by NMR. <i>Nature Methods</i> , 2009 , 6, 625-6	21.6	51
101	The beta(E)-domain of wheat E(c)-1 metallothionein: a metal-binding domain with a distinctive structure. <i>Journal of Molecular Biology</i> , 2009 , 387, 207-18	6.5	80
100	Structural and functional characterization of the NHR1 domain of the Drosophila neuralized E3 ligase in the notch signaling pathway. <i>Journal of Molecular Biology</i> , 2009 , 393, 478-95	6.5	26
99	SAIL--stereo-array isotope labeling. <i>Quarterly Reviews of Biophysics</i> , 2009 , 42, 247-300	7	57
98	Solution structure of the RNA binding domain in the human muscleblind-like protein 2. <i>Protein Science</i> , 2009 , 18, 80-91	6.3	19

97	Structural basis for the selectivity of the external thioesterase of the surfactin synthetase. <i>Nature</i> , 2008 , 454, 907-11	50.4	103
96	Solution structure of the extraterminal domain of the bromodomain-containing protein BRD4. <i>Protein Science</i> , 2008 , 17, 2174-9	6.3	38
95	Solution structure of the c-terminal dimerization domain of SARS coronavirus nucleocapsid protein solved by the SAIL-NMR method. <i>Journal of Molecular Biology</i> , 2008 , 380, 608-22	6.5	87
94	Structural basis of the role of the NikA ribbon-helix-helix domain in initiating bacterial conjugation. <i>Journal of Molecular Biology</i> , 2008 , 384, 690-701	6.5	18
93	Solution structure of the second RNA recognition motif (RRM) domain of murine T cell intracellular antigen-1 (TIA-1) and its RNA recognition mode. <i>Biochemistry</i> , 2008 , 47, 6437-50	3.2	15
92	The RRM domain of poly(A)-specific ribonuclease has a noncanonical binding site for mRNA cap analog recognition. <i>Nucleic Acids Research</i> , 2008 , 36, 4754-67	20.1	39
91	Transmembrane segment enhanced labeling as a tool for the backbone assignment of alpha-helical membrane proteins. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008 , 105, 8262-7	11.5	35
90	Complex assembly mechanism and an RNA-binding mode of the human p14-SF3b155 spliceosomal protein complex identified by NMR solution structure and functional analyses. <i>Proteins: Structure, Function and Bioinformatics</i> , 2008 , 71, 1617-36	4.2	13
89	Structure of the putative 32 kDa myrosinase-binding protein from Arabidopsis (At3g16450.1) determined by SAIL-NMR. <i>FEBS Journal</i> , 2008 , 275, 5873-84	5.7	25
88	Automated structure determination of proteins with the SAIL-FLYA NMR method. <i>Nature Protocols</i> , 2007 , 2, 2896-902	18.8	45
87	KUJIRA, a package of integrated modules for systematic and interactive analysis of NMR data directed to high-throughput NMR structure studies. <i>Journal of Biomolecular NMR</i> , 2007 , 39, 31-52	3	137
86	Structural insight into dimeric interaction of the SARAH domains from Mst1 and RASSF family proteins in the apoptosis pathway. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2007 , 104, 9236-41	11.5	109
85	Solution structure of an atypical WW domain in a novel beta-clam-like dimeric form. <i>FEBS Letters</i> , 2007 , 581, 462-8	3.8	30
84	Solution structures of the SURP domains and the subunit-assembly mechanism within the splicing factor SF3a complex in 17S U2 snRNP. <i>Structure</i> , 2006 , 14, 1677-89	5.2	19
83	A photoswitchable miniprotein based on the sequence of avian pancreatic polypeptide. <i>Angewandte Chemie - International Edition</i> , 2006 , 45, 6297-300	16.4	35
82	Fully automated structure determinations of the Fes SH2 domain using different sets of NMR spectra. <i>Magnetic Resonance in Chemistry</i> , 2006 , 44 Spec No, S83-8	2.1	9
81	Evaluation of stereo-array isotope labeling (SAIL) patterns for automated structural analysis of proteins with CYANA. <i>Magnetic Resonance in Chemistry</i> , 2006 , 44 Spec No, S152-7	2.1	28
80	A Photoswitchable Miniprotein Based on the Sequence of Avian Pancreatic Polypeptide. <i>Angewandte Chemie</i> , 2006 , 118, 6445-6448	3.6	18

79	NMR structure of the R-module: a parallel beta-roll subunit from an <i>Azotobacter vinelandii</i> mannuronan C-5 epimerase. <i>Journal of Biological Chemistry</i> , 2006 , 281, 7350-6	5.4	31
78	Automated protein structure determination from NMR spectra. <i>Journal of the American Chemical Society</i> , 2006 , 128, 13112-22	16.4	136
77	Symbolic NMR product operator calculations. <i>International Journal of Quantum Chemistry</i> , 2006 , 106, 344-350	2.1	14
76	Solution structures of the first and fourth TSR domains of F-spondin. <i>Proteins: Structure, Function and Bioinformatics</i> , 2006 , 64, 665-72	4.2	31
75	Optimal isotope labelling for NMR protein structure determinations. <i>Nature</i> , 2006 , 440, 52-7	50.4	389
74	Solution structure of the antifreeze-like domain of human sialic acid synthase. <i>Protein Science</i> , 2006 , 15, 1010-6	6.3	18
73	Solution structure of the rhodanese homology domain At4g01050(175-295) from <i>Arabidopsis thaliana</i> . <i>Protein Science</i> , 2005 , 14, 224-30	6.3	22
72	Solution structure of the PWWP domain of the hepatoma-derived growth factor family. <i>Protein Science</i> , 2005 , 14, 756-64	6.3	44
71	NMR solution structure of the monomeric form of the bacteriophage lambda capsid stabilizing protein gpD. <i>Journal of Biomolecular NMR</i> , 2005 , 31, 351-6	3	15
70	Solution structure of the Src homology 2 domain from the human feline sarcoma oncogene Fes. <i>Journal of Biomolecular NMR</i> , 2005 , 31, 357-61	3	15
69	Solution structure of the mouse enhancer of rudimentary protein reveals a novel fold. <i>Journal of Biomolecular NMR</i> , 2005 , 32, 329-34	3	13
68	RECOORD: a recalculated coordinate database of 500+ proteins from the PDB using restraints from the BioMagResBank. <i>Proteins: Structure, Function and Bioinformatics</i> , 2005 , 59, 662-72	4.2	285
67	Prion protein NMR structures of chickens, turtles, and frogs. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005 , 102, 651-5	11.5	154
66	Prion protein NMR structures of cats, dogs, pigs, and sheep. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005 , 102, 640-5	11.5	178
65	Solution structure of the RWD domain of the mouse GCN2 protein. <i>Protein Science</i> , 2004 , 13, 2089-100	6.3	56
64	Assignments of ¹ H and ¹⁵ N resonances of the bacteriophage lambda capsid stabilizing protein gpD. <i>Journal of Biomolecular NMR</i> , 2004 , 28, 89-90	3	3
63	NMR assignment of the hypothetical ENTH-VHS domain At3g16270 from <i>Arabidopsis thaliana</i> . <i>Journal of Biomolecular NMR</i> , 2004 , 29, 205-6	3	18
62	NMR assignment of the hypothetical rhodanese domain At4g01050 from <i>Arabidopsis thaliana</i> . <i>Journal of Biomolecular NMR</i> , 2004 , 29, 207-8	3	12

61	NMR assignment of the SH2 domain from the human feline sarcoma oncogene FES. <i>Journal of Biomolecular NMR</i> , 2004 , 30, 463-4	3	13
60	Automated NMR structure calculation with CYANA. <i>Methods in Molecular Biology</i> , 2004 , 278, 353-78	1.4	1018
59	NMR structure of the integral membrane protein OmpX. <i>Journal of Molecular Biology</i> , 2004 , 336, 1211-21	1.5	152
58	Solution structure of the 162 residue C-terminal domain of human elongation factor 1Bgamma. <i>Journal of Biological Chemistry</i> , 2003 , 278, 43443-51	5.4	10
57	Influence of the completeness of chemical shift assignments on NMR structures obtained with automated NOE assignment. <i>Journal of Structural and Functional Genomics</i> , 2003 , 4, 179-89		81
56	ATP-induced conformational changes of the nucleotide-binding domain of Na,K-ATPase. <i>Nature Structural and Molecular Biology</i> , 2003 , 10, 468-74	17.6	88
55	Automated NMR protein structure calculation. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2003 , 43, 105-125	10.4	210
54	NMR structure of a variant human prion protein with two disulfide bridges. <i>Journal of Molecular Biology</i> , 2003 , 326, 225-34	6.5	35
53	NMR structure of the human doppel protein. <i>Journal of Molecular Biology</i> , 2003 , 326, 1549-57	6.5	58
52	NMR structure of the heme chaperone CcmE reveals a novel functional motif. <i>Structure</i> , 2002 , 10, 1551-7	5.2	57
51	The NMR structure of the class I human ubiquitin-conjugating enzyme 2b. <i>Journal of Biomolecular NMR</i> , 2002 , 22, 89-92	3	13
50	Protein NMR structure determination with automated NOE-identification in the NOESY spectra using the new software ATNOS. <i>Journal of Biomolecular NMR</i> , 2002 , 24, 171-89	3	410
49	NMR structure of the unliganded Bombyx mori pheromone-binding protein at physiological pH. <i>FEBS Letters</i> , 2002 , 531, 314-8	3.8	76
48	Protein NMR structure determination with automated NOE assignment using the new software CANDID and the torsion angle dynamics algorithm DYANA. <i>Journal of Molecular Biology</i> , 2002 , 319, 209-27	6.5	1330
47	NMR structures of 36 and 73-residue fragments of the calreticulin P-domain. <i>Journal of Molecular Biology</i> , 2002 , 322, 773-84	6.5	49
46	NMR studies in aqueous solution fail to identify significant conformational differences between the monomeric forms of two Alzheimer peptides with widely different plaque-competence, A beta(1-40)(ox) and A beta(1-42)(ox). <i>FEBS Journal</i> , 2001 , 268, 5930-6		190
45	Sampling of conformation space in torsion angle dynamics calculations. <i>Computer Physics Communications</i> , 2001 , 138, 155-169	4.2	7
44	NMR structure of the calreticulin P-domain. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2001 , 98, 3133-8	11.5	154

43	NMR structure reveals intramolecular regulation mechanism for pheromone binding and release. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2001 , 98, 14374-9	11.5	237
42	Point-centered domain decomposition for parallel molecular dynamics simulation. <i>Computer Physics Communications</i> , 2000 , 124, 139-147	4.2	142
41	Sequence-specific NMR assignment of proteins by global fragment mapping with the program MAPPER. <i>Journal of Biomolecular NMR</i> , 2000 , 18, 129-37	3	78
40	NMR structures of three single-residue variants of the human prion protein. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2000 , 97, 8340-5	11.5	92
39	NMR structure of the sea urchin (<i>Strongylocentrotus purpuratus</i>) metallothionein MTA. <i>Journal of Molecular Biology</i> , 1999 , 291, 417-28	6.5	79
38	Sequence-specific ¹ H, ¹⁵ N, and ¹³ C assignments of the periplasmic chaperone FimC from <i>Escherichia coli</i> . <i>Journal of Biomolecular NMR</i> , 1998 , 11, 229-30	3	6
37	PSEUDYANA for NMR structure calculation of paramagnetic metalloproteins using torsion angle molecular dynamics. <i>Journal of Biomolecular NMR</i> , 1998 , 12, 553-7	3	62
36	Conformational analysis of protein and nucleic acid fragments with the new grid search algorithm FOUND. <i>Journal of Biomolecular NMR</i> , 1998 , 12, 543-8	3	49
35	NMR solution structure of the periplasmic chaperone FimC. <i>Nature Structural Biology</i> , 1998 , 5, 885-90		37
34	Automated peak picking and peak integration in macromolecular NMR spectra using AUTOPSY. <i>Journal of Magnetic Resonance</i> , 1998 , 135, 288-97	3	108
33	Structure calculation of biological macromolecules from NMR data. <i>Quarterly Reviews of Biophysics</i> , 1998 , 31, 145-237	7	133
32	Calculating protein structures from NMR data. <i>Methods in Molecular Biology</i> , 1997 , 60, 157-94	1.4	15
31	The NMR solution conformation of unligated human cyclophilin A. <i>Journal of Molecular Biology</i> , 1997 , 272, 64-81	6.5	66
30	Torsion angle dynamics for NMR structure calculation with the new program DYANA. <i>Journal of Molecular Biology</i> , 1997 , 273, 283-98	6.5	2561
29	Automated combined assignment of NOESY spectra and three-dimensional protein structure determination. <i>Journal of Biomolecular NMR</i> , 1997 , 10, 351-62	3	128
28	GARANT-a general algorithm for resonance assignment of multidimensional nuclear magnetic resonance spectra 1997 , 18, 139-149		112
27	Hydration and DNA recognition by homeodomains. <i>Cell</i> , 1996 , 85, 1057-65	56.2	104
26	NMR studies of the hydration of biological macromolecules. <i>Faraday Discussions</i> , 1996 , 103, 245-253	3.6	42

25	Molecular dynamics simulations on Cray clusters using the SCIDDLE-PVM environment. <i>Lecture Notes in Computer Science</i> , 1996 , 142-149	0.9	3
24	Conformational sampling by NMR solution structures calculated with the program DIANA evaluated by comparison with long-time molecular dynamics calculations in explicit water. <i>Proteins: Structure, Function and Bioinformatics</i> , 1996 , 24, 304-13	4.2	23
23	Automated sequence-specific NMR assignment of homologous proteins using the program GARANT. <i>Journal of Biomolecular NMR</i> , 1996 , 7, 207-13	3	113
22	The new program OPAL for molecular dynamics simulations and energy refinements of biological macromolecules. <i>Journal of Biomolecular NMR</i> , 1996 , 8, 136-46	3	174
21	Ancestral beta gamma-crystallin precursor structure in a yeast killer toxin. <i>Nature Structural and Molecular Biology</i> , 1996 , 3, 662-5	17.6	70
20	The program XEASY for computer-supported NMR spectral analysis of biological macromolecules. <i>Journal of Biomolecular NMR</i> , 1995 , 6, 1-10	3	1482
19	Structure and internal dynamics of the bovine pancreatic trypsin inhibitor in aqueous solution from long-time molecular dynamics simulations. <i>Proteins: Structure, Function and Bioinformatics</i> , 1995 , 23, 49-62	4.2	40
18	NMR solution structure of the recombinant tick anticoagulant protein (rTAP), a factor Xa inhibitor from the tick <i>Ornithodoros moubata</i> . <i>FEBS Letters</i> , 1994 , 352, 251-7	3.8	40
17	Computer-Supported Protein Structure Determination by NMR. <i>NATO ASI Series Series B: Physics</i> , 1994 , 197-207		
16	Nuclear magnetic resonance solution structure of dendrotoxin K from the venom of <i>Dendroaspis polylepis polylepis</i> . <i>Journal of Molecular Biology</i> , 1993 , 234, 735-50	6.5	68
15	POMA: A Complete Mathematica Implementation of the NMR Product-Operator Formalism. <i>Journal of Magnetic Resonance Series A</i> , 1993 , 101, 103-105		56
14	The program ASNO for computer-supported collection of NOE upper distance constraints as input for protein structure determination. <i>Journal of Biomolecular NMR</i> , 1993 , 3, 601	3	65
13	Protein dynamics studied by rotating frame ¹⁵ N spin relaxation times. <i>Journal of Biomolecular NMR</i> , 1993 , 3, 151-64	3	150
12	Determination of a high-quality nuclear magnetic resonance solution structure of the bovine pancreatic trypsin inhibitor and comparison with three crystal structures. <i>Journal of Molecular Biology</i> , 1992 , 227, 757-75	6.5	168
11	Nuclear magnetic resonance solution structure of hirudin(1-51) and comparison with corresponding three-dimensional structures determined using the complete 65-residue hirudin polypeptide chain. <i>Journal of Molecular Biology</i> , 1992 , 228, 1193-205	6.5	50
10	Impact of protein-protein contacts on the conformation of thrombin-bound hirudin studied by comparison with the nuclear magnetic resonance solution structure of hirudin(1-51). <i>Journal of Molecular Biology</i> , 1992 , 228, 1206-11	6.5	27
9	Determination of scalar coupling constants by inverse Fourier transformation of in-phase multiplets. <i>Journal of Magnetic Resonance</i> , 1992 , 99, 552-560		16
8	Processing of multi-dimensional NMR data with the new software PROSA. <i>Journal of Biomolecular NMR</i> , 1992 , 2, 619-629	3	267

7	FLATTA new procedure for high-quality baseline correction of multidimensional NMR spectra. <i>Journal of Magnetic Resonance</i> , 1992 , 96, 403-407		6
6	Complete relaxation matrix refinement of NMR structures of proteins using analytically calculated dihedral angle derivatives of NOE intensities. <i>Journal of Biomolecular NMR</i> , 1991 , 1, 257-69	3	37
5	Efficient analysis of protein 2D NMR spectra using the software package EASY. <i>Journal of Biomolecular NMR</i> , 1991 , 1, 111-30	3	253
4	Improved efficiency of protein structure calculations from NMR data using the program DIANA with redundant dihedral angle constraints. <i>Journal of Biomolecular NMR</i> , 1991 , 1, 447-56	3	328
3	Efficient computation of three-dimensional protein structures in solution from nuclear magnetic resonance data using the program DIANA and the supporting programs CALIBA, HABAS and GLOMSA. <i>Journal of Molecular Biology</i> , 1991 , 217, 517-30	6.5	864
2	Structure determination of the Antp (C39----S) homeodomain from nuclear magnetic resonance data in solution using a novel strategy for the structure calculation with the programs DIANA, CALIBA, HABAS and GLOMSA. <i>Journal of Molecular Biology</i> , 1991 , 217, 531-40	6.5	124
1	Structure-based methyl resonance assignment with MethylFLYA		1