

# Jerome Boisbouvier

## List of Publications by Year in descending order

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54  
papers

2,413  
citations

201674

27  
h-index

206112

48  
g-index

60  
all docs

60  
docs citations

60  
times ranked

2331  
citing authors

#	ARTICLE	IF	CITATIONS
1	Structural basis for the inhibition of IAPP fibril formation by the co-chaperonin prefoldin. <i>Nature Communications</i> , 2022, 13, 2363.	12.8	5
2	Optimized precursor to simplify assignment transfer between backbone resonances and stereospecifically labelled valine and leucine methyl groups: application to human Hsp90 N-terminal domain. <i>Journal of Biomolecular NMR</i> , 2021, 75, 221-232.	2.8	7
3	Backbone and methyl resonances assignment of the 87 kDa prefoldin from <i>Pyrococcus horikoshii</i> . <i>Biomolecular NMR Assignments</i> , 2021, 15, 351-360.	0.8	5
4	In Vitro Production of Perdeuterated Proteins in H <sub>2</sub> O for Biomolecular NMR Studies. <i>Methods in Molecular Biology</i> , 2021, 2199, 127-149.	0.9	14
5	Spectral editing of intra- and inter-chain methyl <sup>13</sup> C methyl NOEs in protein complexes. <i>Journal of Biomolecular NMR</i> , 2020, 74, 83-94.	2.8	7
6	Asymmetric Synthesis of Methyl Specifically Labelled L- <sup>13</sup> C Threonine and Application to the NMR Studies of High Molecular Weight Proteins. <i>ChemistrySelect</i> , 2020, 5, 5092-5098.	1.5	8
7	Integrated NMR and cryo-EM atomic-resolution structure determination of a half-megadalton enzyme complex. <i>Nature Communications</i> , 2019, 10, 2697.	12.8	80
8	Aromatic Ring Dynamics, Thermal Activation, and Transient Conformations of a 468 kDa Enzyme by Specific <sup>13</sup> C Labeling and Fast Magic-Angle Spinning NMR. <i>Journal of the American Chemical Society</i> , 2019, 141, 11183-11195.	13.7	43
9	Structural investigation of a chaperonin in action reveals how nucleotide binding regulates the functional cycle. <i>Science Advances</i> , 2018, 4, eaau4196.	10.3	44
10	Advanced isotopic labeling for the NMR investigation of challenging proteins and nucleic acids. <i>Journal of Biomolecular NMR</i> , 2018, 71, 115-117.	2.8	15
11	Observation of CH <sup>13</sup> C Interactions between Methyl and Carbonyl Groups in Proteins. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 7564-7567.	13.8	17
12	Observation of CH <sup>13</sup> C Interactions between Methyl and Carbonyl Groups in Proteins. <i>Angewandte Chemie</i> , 2017, 129, 7672-7675.	2.0	5
13	Unraveling self-assembly pathways of the 468-kDa proteolytic machine TET2. <i>Science Advances</i> , 2017, 3, e1601601.	10.3	28
14	Solid-state NMR <sup>13</sup> C- <sup>1</sup> H and <sup>13</sup> C- <sup>13</sup> C 3D/4D Correlation Experiments for Resonance Assignment of Large Proteins. <i>ChemPhysChem</i> , 2017, 18, 2697-2703.	2.1	43
15	Sensitive proton-detected solid-state NMR spectroscopy of large proteins with selective CH <sub>3</sub> labelling: application to the 50S ribosome subunit. <i>Chemical Communications</i> , 2016, 52, 9558-9561.	4.1	30
16	Scrambling free combinatorial labeling of alanine- <sup>13</sup> C, isoleucine- <sup>13</sup> C, leucine-proS and valine-proS methyl groups for the detection of long range NOEs. <i>Journal of Biomolecular NMR</i> , 2015, 61, 73-82.	2.8	37
17	Methyl-specific isotopic labeling: a molecular tool box for solution NMR studies of large proteins. <i>Current Opinion in Structural Biology</i> , 2015, 32, 113-122.	5.7	157
18	CH <sub>3</sub> -specific NMR assignment of alanine, isoleucine, leucine and valine methyl groups in high molecular weight proteins using a single sample. <i>Journal of Biomolecular NMR</i> , 2015, 63, 389-402.	2.8	25

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19	Probing Transient Conformational States of Proteins by Solid-State Relaxation-Dispersion NMR Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 4312-4317.	13.8	81
20	A Cost-Effective Protocol for the Parallel Production of Libraries of <sup>13</sup> CH <sub>3</sub> -Specifically Labeled Mutants for NMR Studies of High Molecular Weight Proteins. <i>Methods in Molecular Biology</i> , 2014, 1091, 229-244.	0.9	10
21	Specific labeling and assignment strategies of valine methyl groups for NMR studies of high molecular weight proteins. <i>Journal of Biomolecular NMR</i> , 2013, 57, 251-262.	2.8	55
22	An optimized isotopic labelling strategy of isoleucine- <sup>13</sup> C <sub>2</sub> methyl groups for solution NMR studies of high molecular weight proteins. <i>Chemical Communications</i> , 2012, 48, 1434-1436.	4.1	37
23	A simple biosynthetic method for stereospecific resonance assignment of prochiral methyl groups in proteins. <i>Journal of Biomolecular NMR</i> , 2011, 49, 61-67.	2.8	17
24	A systematic mutagenesis-driven strategy for site-resolved NMR studies of supramolecular assemblies. <i>Journal of Biomolecular NMR</i> , 2011, 50, 229-236.	2.8	70
25	Rapid measurement of residual dipolar couplings for fast fold elucidation of proteins. <i>Journal of Biomolecular NMR</i> , 2011, 51, 369-378.	2.8	18
26	Solid-State NMR Measurements of Asymmetric Dipolar Couplings Provide Insight into Protein Side-Chain Motion. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 11005-11009.	13.8	53
27	NMR structure of the A730 loop of the <i>Neurospora VS</i> ribozyme: insights into the formation of the active site. <i>Nucleic Acids Research</i> , 2011, 39, 4427-4437.	14.5	20
28	Stereospecific Isotopic Labeling of Methyl Groups for NMR Spectroscopic Studies of High-Molecular-Weight Proteins. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 1958-1962.	13.8	193
29	Inside Cover: Stereospecific Isotopic Labeling of Methyl Groups for NMR Spectroscopic Studies of High-Molecular-Weight Proteins ( <i>Angew. Chem. Int. Ed.</i> 11/2010). <i>Angewandte Chemie - International Edition</i> , 2010, 49, 1896-1896.	13.8	1
30	Direct detection of CH/π interactions in proteins. <i>Nature Chemistry</i> , 2010, 2, 466-471.	13.6	247
31	An efficient protocol for the complete incorporation of methyl-protonated alanine in perdeuterated protein. <i>Journal of Biomolecular NMR</i> , 2009, 43, 111-119.	2.8	140
32	Parallel screening and optimization of protein constructs for structural studies. <i>Protein Science</i> , 2009, 18, 434-439.	7.6	7
33	Longitudinal-Relaxation-Enhanced NMR Experiments for the Study of Nucleic Acids in Solution. <i>Journal of the American Chemical Society</i> , 2009, 131, 8571-8577.	13.7	90
34	Fast Two-Dimensional NMR Spectroscopy of High Molecular Weight Protein Assemblies. <i>Journal of the American Chemical Society</i> , 2009, 131, 3448-3449.	13.7	99
35	Liquid-crystal NMR structure of HIV TAR RNA bound to its SELEX RNA aptamer reveals the origins of the high stability of the complex. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008, 105, 9210-9215.	7.1	44
36	High-Accuracy Distance Measurement between Remote Methyls in Specifically Protonated Proteins. <i>Journal of the American Chemical Society</i> , 2007, 129, 472-473.	13.7	43

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37	Sensitivity-optimized experiment for the measurement of residual dipolar couplings between amide protons. <i>Journal of Biomolecular NMR</i> , 2007, 38, 47-55.	2.8	6
38	Measurement of eight scalar and dipolar couplings for methine/methylene pairs in proteins and nucleic acids. <i>Journal of Biomolecular NMR</i> , 2005, 31, 201-216.	2.8	32
39	Accurate measurement of $^{15}\text{N}/^{13}\text{C}$ residual dipolar couplings in nucleic acids. <i>Journal of Biomolecular NMR</i> , 2005, 31, 231-241.	2.8	23
40	Resolution-Enhanced Base-Type-Edited HCN Experiment for RNA. <i>Journal of Biomolecular NMR</i> , 2005, 32, 263-271.	2.8	4
41	Resolution-optimized NMR measurement of 1DCH, 1DCC and 2DCH residual dipolar couplings in nucleic acid bases. <i>Journal of Biomolecular NMR</i> , 2004, 30, 287-301.	2.8	28
42	Relaxation-Optimized NMR Spectroscopy of Methylene Groups in Proteins and Nucleic Acids. <i>Journal of the American Chemical Society</i> , 2004, 126, 10560-10570.	13.7	71
43	Experimental and Theoretical Determination of Nucleic Acid Magnetic Susceptibility: Importance for the Study of Dynamics by Field-Induced Residual Dipolar Couplings. <i>Journal of the American Chemical Society</i> , 2004, 126, 10820-10821.	13.7	25
44	Rotational diffusion tensor of nucleic acids from $^{13}\text{C}$ NMR relaxation. <i>Journal of Biomolecular NMR</i> , 2003, 27, 133-142.	2.8	49
45	Direct observation of dipolar couplings between distant protons in weakly aligned nucleic acids. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2003, 100, 11333-11338.	7.1	34
46	Long-Range Magnetization Transfer between Uncoupled Nuclei by Dipole-Dipole Cross-Correlated Relaxation: A Precise Probe of $\beta$ -Sheet Geometry in Proteins. <i>Journal of the American Chemical Society</i> , 2002, 124, 11038-11045.	13.7	10
47	Base-type-selective high-resolution $^{13}\text{C}$ edited NOESY for sequential assignment of large RNAs. <i>Journal of Biomolecular NMR</i> , 2001, 19, 141-151.	2.8	21
48	Simultaneous determination of disulphide bridge topology and three-dimensional structure using ambiguous intersulphur distance restraints: possibilities and limitations. <i>Journal of Biomolecular NMR</i> , 2000, 16, 197-208.	2.8	12
49	NMR Determination of Sugar Puckers in Nucleic Acids from CSA-Dipolar Cross-Correlated Relaxation. <i>Journal of the American Chemical Society</i> , 2000, 122, 6779-6780.	13.7	43
50	Long-Range Structural Information in NMR Studies of Paramagnetic Molecules from Electron Spin-Nuclear Spin Cross-Correlated Relaxation. <i>Journal of the American Chemical Society</i> , 1999, 121, 7700-7701.	13.7	67
51	Improved Sensitivity and Resolution in $^1\text{H}/^{13}\text{C}$ NMR Experiments of RNA. <i>Journal of the American Chemical Society</i> , 1998, 120, 11845-11851.	13.7	101
52	A structural homologue of colipase in black mamba venom revealed by NMR floating disulphide bridge analysis. <i>Journal of Molecular Biology</i> , 1998, 283, 205-219.	4.2	68
53	Chapter 1. Isotope-Labeling of Methyl Groups for NMR Studies of Large Proteins. <i>RSC Biomolecular Sciences</i> , 0, , 1-24.	0.4	5
54	NMR assignment of human HSP90 N-terminal domain bound to a long residence time resorcinol ligand. <i>Biomolecular NMR Assignments</i> , 0, , .	0.8	0