

G Marius Clore

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.

561
papers

60,146
citations

116
h-index

232
g-index

592
ext. papers

62,995
ext. citations

8.8
avg, IF

7.49
L-index

#	Paper	IF	Citations
561	Time-resolved DEER EPR and solid-state NMR afford kinetic and structural elucidation of substrate binding to Ca-ligated calmodulin.. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022 , 119,	11.5	2
560	NMR methods for exploring 'dark' states in ligand binding and protein-protein interactions.. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2022 , 128, 1-24	10.4	2
559	Visualization of Sparsely-populated Lower-order Oligomeric States of Human Mitochondrial Hsp60 by Cryo-electron Microscopy. <i>Journal of Molecular Biology</i> , 2021 , 433, 167322	6.5	0
558	Probing Side-Chain Dynamics in Proteins by NMR Relaxation of Isolated C Magnetization Modes in CH Methyl Groups. <i>Journal of Physical Chemistry B</i> , 2021 , 125, 3343-3352	3.4	2
557	Probing the Interaction of Huntingtin Exon-1 Polypeptides with the Chaperonin Nanomachine GroEL. <i>ChemBioChem</i> , 2021 , 22, 1985-1991	3.8	2
556	Quantitative Exchange NMR-Based Analysis of Huntingtin-SH3 Interactions Suggests an Allosteric Mechanism of Inhibition of Huntingtin Aggregation. <i>Journal of the American Chemical Society</i> , 2021 , 143, 9672-9681	16.4	1
555	A simple and cost-effective protocol for high-yield expression of deuterated and selectively isoleucine/leucine/valine methyl protonated proteins in Escherichia coli grown in shaker flasks. <i>Journal of Biomolecular NMR</i> , 2021 , 75, 83-87	3	1
554	The measurement of relaxation rates of degenerate H transitions in methyl groups of proteins using acute angle radiofrequency pulses. <i>Journal of Magnetic Resonance</i> , 2021 , 330, 107034	3	1
553	Submillisecond Freezing Permits Cryoprotectant-Free EPR Double Electron-Electron Resonance Spectroscopy. <i>ChemPhysChem</i> , 2020 , 21, 1224-1229	3.2	12
552	Probing the Interaction between HIV-1 Protease and the Homodimeric p66/p66' Reverse Transcriptase Precursor by Double Electron-Electron Resonance EPR Spectroscopy. <i>ChemBioChem</i> , 2020 , 21, 3051-3055	3.8	2
551	Abrogation of prenucleation, transient oligomerization of the Huntingtin exon 1 protein by human profilin I. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020 , 117, 5844-5852 ¹¹	11.5	11
550	Kinetics of Fast Tetramerization of the Huntingtin Exon 1 Protein Probed by Concentration-Dependent On-Resonance Measurements. <i>Journal of Physical Chemistry Letters</i> , 2020 , 11, 5643-5648	6.4	3
549	Determining methyl sidechain conformations in a CS-ROSETTA model using methyl H-C residual dipolar couplings. <i>Journal of Biomolecular NMR</i> , 2020 , 74, 111-118	3	2
548	Magic-Angle-Pulse Driven Separation of Degenerate H Transitions in Methyl Groups of Proteins: Application to Studies of Methyl Axis Dynamics. <i>ChemPhysChem</i> , 2020 , 21, 1087-1091	3.2	7
547	Optimized NMR Experiments for the Isolation of I=1/2 Manifold Transitions in Methyl Groups of Proteins. <i>ChemPhysChem</i> , 2020 , 21, 13-19	3.2	8
546	XIPP: multi-dimensional NMR analysis software. <i>Journal of Biomolecular NMR</i> , 2020 , 74, 9-25	3	1
545	Optimized selection of slow-relaxing C transitions in methyl groups of proteins: application to relaxation dispersion. <i>Journal of Biomolecular NMR</i> , 2020 , 74, 673-680	3	6

544	An S/T motif controls reversible oligomerization of the Hsp40 chaperone DNAJB6b through subtle reorganization of a β sheet backbone. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020 , 117, 30441-30450	11.5	7
543	filtering by H-methyl labeling in a deuterated protein for pulsed double electron-electron resonance EPR. <i>Chemical Communications</i> , 2020 , 56, 10890-10893	5.8	2
542	A three-dimensional potential of mean force to improve backbone and sidechain hydrogen bond geometry in Xplor-NIH protein structure determination. <i>Protein Science</i> , 2020 , 29, 100-110	6.3	7
541	Quantitative Interpretation of Solvent Paramagnetic Relaxation for Probing Protein-Cosolute Interactions. <i>Journal of the American Chemical Society</i> , 2020 , 142, 8281-8290	16.4	4
540	Spatial domain organization in the HIV-1 reverse transcriptase p66 homodimer precursor probed by double electron-electron resonance EPR. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019 , 116, 17809-17816	11.5	10
539	Inhibition of HIV Maturation via Selective Unfolding and Cross-Linking of Gag Polyprotein by a Mercaptobenzamide Acetylator. <i>Journal of the American Chemical Society</i> , 2019 , 141, 8327-8338	16.4	1
538	Probing initial transient oligomerization events facilitating Huntingtin fibril nucleation at atomic resolution by relaxation-based NMR. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019 , 116, 3562-3571	11.5	33
537	Exchange saturation transfer and associated NMR techniques for studies of protein interactions involving high-molecular-weight systems. <i>Journal of Biomolecular NMR</i> , 2019 , 73, 461-469	3	8
536	Protein NMR: Boundless opportunities. <i>Journal of Magnetic Resonance</i> , 2019 , 306, 187-191	3	21
535	Unraveling the structure and dynamics of the human DNAJB6b chaperone by NMR reveals insights into Hsp40-mediated proteostasis. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019 , 116, 21529-21538	11.5	21
534	A simple protocol for expression of isotope-labeled proteins in Escherichia coli grown in shaker flasks at high cell density. <i>Journal of Biomolecular NMR</i> , 2019 , 73, 743-748	3	3
533	Probing transient excited states of the bacterial cell division regulator MinE by relaxation dispersion NMR spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019 , 116, 25446-25455	11.5	8
532	TiO Nanoparticles Catalyze Oxidation of Huntingtin Exon 1-Derived Peptides Impeding Aggregation: A Quantitative NMR Study of Binding and Kinetics. <i>Journal of the American Chemical Society</i> , 2019 , 141, 94-97	16.4	23
531	Targeting a Dark Excited State of HIV-1 Nucleocapsid by Antiretroviral Thioesters Revealed by NMR Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2018 , 57, 2687-2691	16.4	16
530	Targeting a Dark Excited State of HIV-1 Nucleocapsid by Antiretroviral Thioesters Revealed by NMR Spectroscopy. <i>Angewandte Chemie</i> , 2018 , 130, 2717-2721	3.6	2
529	Probing Conformational States of the Finger and Thumb Subdomains of HIV-1 Reverse Transcriptase Using Double Electron-Electron Resonance Electron Paramagnetic Resonance Spectroscopy. <i>Biochemistry</i> , 2018 , 57, 489-493	3.2	6
528	Interaction of Huntingtin Exon-1 Peptides with Lipid-Based Micellar Nanoparticles Probed by Solution NMR and Q-Band Pulsed EPR. <i>Journal of the American Chemical Society</i> , 2018 , 140, 6199-6202	16.4	22
527	Disassembly/reassembly strategy for the production of highly pure GroEL, a tetradecameric supramolecular machine, suitable for quantitative NMR, EPR and mutational studies. <i>Protein Expression and Purification</i> , 2018 , 142, 8-15	2	2

526	Extensive Sampling of the Cavity of the GroEL Nanomachine by Protein Substrates Probed by Paramagnetic Relaxation Enhancement. <i>Journal of Physical Chemistry Letters</i> , 2018 , 9, 3368-3371	6.4	8
525	Xplor-NIH for molecular structure determination from NMR and other data sources. <i>Protein Science</i> , 2018 , 27, 26-40	6.3	96
524	Probing the mechanism of inhibition of amyloid- β (1-42)-induced neurotoxicity by the chaperonin GroEL. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018 , 115, E11924-E11932	11.5	19
523	Decorrelating Kinetic and Relaxation Parameters in Exchange Saturation Transfer NMR: A Case Study of N-Terminal Huntingtin Peptides Binding to Unilamellar Lipid Vesicles. <i>Journal of Physical Chemistry B</i> , 2018 , 122, 11271-11278	3.4	12
522	Confinement and Stabilization of Fyn SH3 Folding Intermediate Mimetics within the Cavity of the Chaperonin GroEL Demonstrated by Relaxation-Based NMR. <i>Biochemistry</i> , 2017 , 56, 903-906	3.2	16
521	Probing the Binding Modes of a Multidomain Protein to Lipid-based Nanoparticles by Relaxation-based NMR. <i>Journal of Physical Chemistry Letters</i> , 2017 , 8, 2535-2540	6.4	7
520	Binding kinetics and substrate selectivity in HIV-1 protease-Gag interactions probed at atomic resolution by chemical exchange NMR. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017 , 114, E9855-E9862	11.5	23
519	Structure Determination of Large Macromolecular Complexes Using NMR 2017 , 316-318		
518	Chaperonin GroEL accelerates protofibril formation and decorates fibrils of the Het-s prion protein. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017 , 114, 9104-9109	11.5	16
517	Towards interpretation of intermolecular paramagnetic relaxation enhancement outside the fast exchange limit. <i>Journal of Biomolecular NMR</i> , 2016 , 66, 1-7	3	5
516	NMR in Structural and Cell Biology 2016 , 98-107		1
515	Long Distance Measurements up to 160 Å in the GroEL Tetradecamer Using Q-Band DEER EPR Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2016 , 55, 15905-15909	16.4	92
514	Long Distance Measurements up to 160 Å in the GroEL Tetradecamer Using Q-Band DEER EPR Spectroscopy. <i>Angewandte Chemie</i> , 2016 , 128, 16137-16141	3.6	26
513	Transient HIV-1 Gag-protease interactions revealed by paramagnetic NMR suggest origins of compensatory drug resistance mutations. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016 , 113, 12456-12461	11.5	19
512	Quantitative Characterization of Configurational Space Sampled by HIV-1 Nucleocapsid Using Solution NMR, X-ray Scattering and Protein Engineering. <i>ChemPhysChem</i> , 2016 , 17, 1548-52	3.2	10
511	Hybrid Approaches to Structural Characterization of Conformational Ensembles of Complex Macromolecular Systems Combining NMR Residual Dipolar Couplings and Solution X-ray Scattering. <i>Chemical Reviews</i> , 2016 , 116, 6305-22	68.1	30
510	Dynamics of SOX2 Interactions with DNA 2016 , 25-41		1
509	Improving NMR Structures of RNA. <i>Structure</i> , 2016 , 24, 806-815	5.2	30

508	Global Dynamics and Exchange Kinetics of a Protein on the Surface of Nanoparticles Revealed by Relaxation-Based Solution NMR Spectroscopy. <i>Journal of the American Chemical Society</i> , 2016 , 138, 5789-92	16.4	49
507	A simple and robust protocol for high-yield expression of perdeuterated proteins in <i>Escherichia coli</i> grown in shaker flasks. <i>Journal of Biomolecular NMR</i> , 2016 , 66, 85-91	3	30
506	Quantitative Resolution of Monomer-Dimer Populations by Inversion Modulated DEER EPR Spectroscopy. <i>ChemPhysChem</i> , 2016 , 17, 2987-2991	3.2	11
505	Visualizing transient dark states by NMR spectroscopy. <i>Quarterly Reviews of Biophysics</i> , 2015 , 48, 35-1167		152
504	Conformation and dynamics of the Gag polyprotein of the human immunodeficiency virus 1 studied by NMR spectroscopy. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015 , 112, 3374-9	11.5	35
503	Intrinsic unfoldase/foldase activity of the chaperonin GroEL directly demonstrated using multinuclear relaxation-based NMR. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015 , 112, 8817-23	11.5	68
502	Dynamic equilibrium between closed and partially closed states of the bacterial Enzyme I unveiled by solution NMR and X-ray scattering. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015 , 112, 11565-70	11.5	24
501	Mutations Proximal to Sites of Autoproteolysis and the β -Helix That Co-evolve under Drug Pressure Modulate the Autoprocessing and Vitality of HIV-1 Protease. <i>Biochemistry</i> , 2015 , 54, 5414-24	3.2	8
500	Reply to Marchenko et al.: Flux analysis of GroEL-assisted protein folding/unfolding. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2015 , 112, E6833-4	11.5	1
499	Dependence of Distance Distributions Derived from Double Electron-Electron Resonance Pulsed EPR Spectroscopy on Pulse-Sequence Time. <i>Angewandte Chemie</i> , 2015 , 127, 5426-5429	3.6	2
498	The energetics of a three-state protein folding system probed by high-pressure relaxation dispersion NMR spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2015 , 54, 11157-61	16.4	21
497	The Energetics of a Three-State Protein Folding System Probed by High-Pressure Relaxation Dispersion NMR Spectroscopy. <i>Angewandte Chemie</i> , 2015 , 127, 11309-11313	3.6	
496	Dependence of distance distributions derived from double electron-electron resonance pulsed EPR spectroscopy on pulse-sequence time. <i>Angewandte Chemie - International Edition</i> , 2015 , 54, 5336-9	16.4	28
495	Practical Aspects of Paramagnetic Relaxation Enhancement in Biological Macromolecules. <i>Methods in Enzymology</i> , 2015 , 564, 485-97	1.7	53
494	The C34 Peptide Fusion Inhibitor Binds to the Six-Helix Bundle Core Domain of HIV-1 gp41 by Displacement of the C-Terminal Helical Repeat Region. <i>Biochemistry</i> , 2015 , 54, 6796-805	3.2	5
493	Large interdomain rearrangement triggered by suppression of micro- to millisecond dynamics in bacterial Enzyme I. <i>Nature Communications</i> , 2015 , 6, 5960	17.4	27
492	Interplay between conformational selection and induced fit in multidomain protein-ligand binding probed by paramagnetic relaxation enhancement. <i>Biophysical Chemistry</i> , 2014 , 186, 3-12	3.5	37
491	A NMR experiment for simultaneous correlations of valine and leucine/isoleucine methyls with carbonyl chemical shifts in proteins. <i>Journal of Biomolecular NMR</i> , 2014 , 58, 1-8	3	15

490	Characterizing methyl-bearing side chain contacts and dynamics mediating amyloid β -protofibril interactions using ^{13}C (methyl)-DEST and lifetime line broadening. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 10345-9	16.4	40
489	Probing the rate-limiting step for intramolecular transfer of a transcription factor between specific sites on the same DNA molecule by $(15)\text{N}$ -exchange NMR spectroscopy. <i>Journal of the American Chemical Society</i> , 2014 , 136, 14369-72	16.4	6
488	Investigation of the Structure and Dynamics of the Capsid-Spacer Peptide 1-Nucleocapsid Fragment of the HIV-1 Gag Polyprotein by Solution NMR Spectroscopy. <i>Angewandte Chemie</i> , 2014 , 126, 1043-1046	3.6	1
487	Investigation of the structure and dynamics of the capsid-spacer peptide 1-nucleocapsid fragment of the HIV-1 gag polyprotein by solution NMR spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 1025-8	16.4	14
486	Binding of HIV-1 gp41-directed neutralizing and non-neutralizing fragment antibody binding domain (Fab) and single chain variable fragment (ScFv) antibodies to the ectodomain of gp41 in the pre-hairpin and six-helix bundle conformations. <i>PLoS ONE</i> , 2014 , 9, e104683	3.7	6
485	Characterizing Methyl-Bearing Side Chain Contacts and Dynamics Mediating Amyloid β -Protofibril Interactions Using ^{13}C -methyl-DEST and Lifetime Line Broadening. <i>Angewandte Chemie</i> , 2014 , 126, 10513-10517	3.6	17
484	Using small angle solution scattering data in Xplor-NIH structure calculations. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2014 , 80, 1-11	10.4	43
483	Encounter complexes and dimensionality reduction in protein-protein association. <i>ELife</i> , 2014 , 3, e013708.9	3.9	42
482	Generating accurate contact maps of transient long-range interactions in intrinsically disordered proteins by paramagnetic relaxation enhancement. <i>Biophysical Journal</i> , 2013 , 104, 1635-6	2.9	6
481	Structure, dynamics and biophysics of the cytoplasmic protein-protein complexes of the bacterial phosphoenolpyruvate: sugar phosphotransferase system. <i>Trends in Biochemical Sciences</i> , 2013 , 38, 515-30	10.3	46
480	Structure and dynamics of full-length HIV-1 capsid protein in solution. <i>Journal of the American Chemical Society</i> , 2013 , 135, 16133-47	16.4	95
479	Sequence-specific determination of protein and peptide concentrations by absorbance at 205 nm. <i>Protein Science</i> , 2013 , 22, 851-8	6.3	219
478	The length of the calmodulin linker determines the extent of transient interdomain association and target affinity. <i>Journal of the American Chemical Society</i> , 2013 , 135, 9648-51	16.4	21
477	Structural basis for enzyme I inhibition by β -ketoglutarate. <i>ACS Chemical Biology</i> , 2013 , 8, 1232-40	4.9	24
476	Seeing the invisible by paramagnetic and diamagnetic NMR. <i>Biochemical Society Transactions</i> , 2013 , 41, 1343-54	5.1	16
475	Probing the transient dark state of substrate binding to GroEL by relaxation-based solution NMR. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2013 , 110, 11361-6	11.5	66
474	Complexes of neutralizing and non-neutralizing affinity matured Fabs with a mimetic of the internal trimeric coiled-coil of HIV-1 gp41. <i>PLoS ONE</i> , 2013 , 8, e78187	3.7	16
473	Smooth statistical torsion angle potential derived from a large conformational database via adaptive kernel density estimation improves the quality of NMR protein structures. <i>Protein Science</i> , 2012 , 21, 1824-36	6.3	46

472	Peptides from second extracellular loop of C-C chemokine receptor type 5 (CCR5) inhibit diverse strains of HIV-1. <i>Journal of Biological Chemistry</i> , 2012 , 287, 15076-86	5.4	20
471	Contrast-matched small-angle X-ray scattering from a heavy-atom-labeled protein in structure determination: application to a lead-substituted calmodulin-peptide complex. <i>Journal of the American Chemical Society</i> , 2012 , 134, 14686-9	16.4	32
470	Coupling between internal dynamics and rotational diffusion in the presence of exchange between discrete molecular conformations. <i>Journal of Chemical Physics</i> , 2012 , 136, 034108	3.9	18
469	Probing exchange kinetics and atomic resolution dynamics in high-molecular-weight complexes using dark-state exchange saturation transfer NMR spectroscopy. <i>Nature Protocols</i> , 2012 , 7, 1523-33	18.8	81
468	An efficient protocol for incorporation of an unnatural amino acid in perdeuterated recombinant proteins using glucose-based media. <i>Journal of Biomolecular NMR</i> , 2012 , 52, 191-5	3	13
467	Structure of the Plasmodium 6-cysteine s48/45 domain. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2012 , 109, 6692-7	11.5	51
466	Solution structure of the IIAChitobiose-HPr complex of the N,N'-diacetylchitobiose branch of the Escherichia coli phosphotransferase system. <i>Journal of Biological Chemistry</i> , 2012 , 287, 23819-29	5.4	11
465	Interplay between minor and major groove-binding transcription factors Sox2 and Oct1 in translocation on DNA studied by paramagnetic and diamagnetic NMR. <i>Journal of Biological Chemistry</i> , 2012 , 287, 14349-63	5.4	24
464	Impact of protein/protein interactions on global intermolecular translocation rates of the transcription factors Sox2 and Oct1 between DNA cognate sites analyzed by z-exchange NMR spectroscopy. <i>Journal of Biological Chemistry</i> , 2012 , 287, 26962-70	5.4	14
463	Conformational selection and substrate binding regulate the monomer/dimer equilibrium of the C-terminal domain of Escherichia coli enzyme I. <i>Journal of Biological Chemistry</i> , 2012 , 287, 26989-98	5.4	25
462	Recent Developments in Biomolecular NMR. <i>RSC Biomolecular Sciences</i> , 2012 ,		5
461	No interaction of barrier-to-autointegration factor (BAF) with HIV-1 MA, cone-rod homeobox (Crx) or MAN1-C in absence of DNA. <i>PLoS ONE</i> , 2011 , 6, e25123	3.7	11
460	A short recollection on the paper entitled "A common sense approach to peak picking in two-, three-, and four-dimensional spectra using automatic computer analysis of contour diagrams" by D.S. Garrett, R. Powers, A.M. Gronenborn, and G.M. Clore [J. Magn. Reson. 95 (1991) 214-220]. <i>Journal of Magnetic Resonance</i> , 2011 , 213, 364-5	3	13
459	A common sense approach to peak picking in two-, three-, and four-dimensional spectra using automatic computer analysis of contour diagrams. 1991. <i>Journal of Magnetic Resonance</i> , 2011 , 213, 357-63		126
458	A rigid disulfide-linked nitroxide side chain simplifies the quantitative analysis of PRE data. <i>Journal of Biomolecular NMR</i> , 2011 , 51, 105-14	3	49
457	Exploring translocation of proteins on DNA by NMR. <i>Journal of Biomolecular NMR</i> , 2011 , 51, 209-19	3	31
456	Automated sequence- and stereo-specific assignment of methyl-labeled proteins by paramagnetic relaxation and methyl-methyl nuclear Overhauser enhancement spectroscopy. <i>Journal of Biomolecular NMR</i> , 2011 , 51, 319-28	3	44
455	Exploring sparsely populated states of macromolecules by diamagnetic and paramagnetic NMR relaxation. <i>Protein Science</i> , 2011 , 20, 229-46	6.3	55

454	Impact of ¹⁵ N R2/R1 relaxation restraints on molecular size, shape, and bond vector orientation for NMR protein structure determination with sparse distance restraints. <i>Journal of the American Chemical Society</i> , 2011 , 133, 6154-7	16.4	7
453	Transient, sparsely populated compact states of apo and calcium-loaded calmodulin probed by paramagnetic relaxation enhancement: interplay of conformational selection and induced fit. <i>Journal of the American Chemical Society</i> , 2011 , 133, 18966-74	16.4	99
452	Combined use of residual dipolar couplings and solution X-ray scattering to rapidly probe rigid-body conformational transitions in a non-phosphorylatable active-site mutant of the 128 kDa enzyme I dimer. <i>Journal of the American Chemical Society</i> , 2011 , 133, 424-7	16.4	31
451	Atomic-resolution dynamics on the surface of amyloid- β protofibrils probed by solution NMR. <i>Nature</i> , 2011 , 480, 268-72	50.4	320
450	Solution structure of the monovalent lectin microvirin in complex with Man(α)(1-2)Man provides a basis for anti-HIV activity with low toxicity. <i>Journal of Biological Chemistry</i> , 2011 , 286, 20788-96	5.4	55
449	Intra- and intermolecular translocation of the bi-domain transcription factor Oct1 characterized by liquid crystal and paramagnetic NMR. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2011 , 108, E169-76	11.5	35
448	Mechanistic details of a protein-protein association pathway revealed by paramagnetic relaxation enhancement titration measurements. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2010 , 107, 1379-84	11.5	71
447	Solution structure of the IIAChitobiose-IIBChitobiose complex of the N,N'-diacetylchitobiose branch of the Escherichia coli phosphotransferase system. <i>Journal of Biological Chemistry</i> , 2010 , 285, 4173-4184	5.4	8
446	Structural basis of HIV-1 neutralization by affinity matured Fabs directed against the internal trimeric coiled-coil of gp41. <i>PLoS Pathogens</i> , 2010 , 6, e1001182	7.6	37
445	Structure-independent analysis of the breadth of the positional distribution of disordered groups in macromolecules from order parameters for long, variable-length vectors using NMR paramagnetic relaxation enhancement. <i>Journal of the American Chemical Society</i> , 2010 , 132, 13346-56	16.4	31
444	Kinetics of amyloid beta monomer-to-oligomer exchange by NMR relaxation. <i>Journal of the American Chemical Society</i> , 2010 , 132, 9948-51	16.4	152
443	Direct use of ¹⁵ N relaxation rates as experimental restraints on molecular shape and orientation for docking of protein-protein complexes. <i>Journal of the American Chemical Society</i> , 2010 , 132, 5987-9	16.4	15
442	Solution structure of the 128 kDa enzyme I dimer from Escherichia coli and its 146 kDa complex with HPr using residual dipolar couplings and small- and wide-angle X-ray scattering. <i>Journal of the American Chemical Society</i> , 2010 , 132, 13026-45	16.4	100
441	Water in the polar and nonpolar cavities of the protein interleukin-1 β . <i>Journal of Physical Chemistry B</i> , 2010 , 114, 16290-7	3.4	37
440	Structural basis of the association of HIV-1 matrix protein with DNA. <i>PLoS ONE</i> , 2010 , 5, e15675	3.7	19
439	Affinity maturation by targeted diversification of the CDR-H2 loop of a monoclonal Fab derived from a synthetic naive human antibody library and directed against the internal trimeric coiled-coil of gp41 yields a set of Fabs with improved HIV-1 neutralization potency and breadth. <i>Virology</i> , 2009 , 388, 112-8	3.6	19
438	Using multiple quantum coherence to increase the ¹⁵ N resolution in a three-dimensional TROSY HNC0 experiment for accurate PRE and RDC measurements. <i>Journal of Magnetic Resonance</i> , 2009 , 200, 173-7	3	8
437	Using the experimentally determined components of the overall rotational diffusion tensor to restrain molecular shape and size in NMR structure determination of globular proteins and protein-protein complexes. <i>Journal of the American Chemical Society</i> , 2009 , 131, 9522-31	16.4	24

436	Theory, practice, and applications of paramagnetic relaxation enhancement for the characterization of transient low-population states of biological macromolecules and their complexes. <i>Chemical Reviews</i> , 2009 , 109, 4108-39	68.1	575
435	Visualizing transient events in amino-terminal autoprocessing of HIV-1 protease. <i>Nature</i> , 2008 , 455, 693-6	50.4	111
434	Antibody elicited against the gp41 N-heptad repeat (NHR) coiled-coil can neutralize HIV-1 with modest potency but non-neutralizing antibodies also bind to NHR mimetics. <i>Virology</i> , 2008 , 377, 170-83	3.6	45
433	Visualization of transient ultra-weak protein self-association in solution using paramagnetic relaxation enhancement. <i>Journal of the American Chemical Society</i> , 2008 , 130, 4048-56	16.4	73
432	Visualizing lowly-populated regions of the free energy landscape of macromolecular complexes by paramagnetic relaxation enhancement. <i>Molecular BioSystems</i> , 2008 , 4, 1058-69		58
431	A pseudopotential for improving the packing of ellipsoidal protein structures determined from NMR data. <i>Journal of Physical Chemistry B</i> , 2008 , 112, 6070-3	3.4	21
430	Sequestering of the prehairpin intermediate of gp41 by peptide N36Mut(e,g) potentiates the human immunodeficiency virus type 1 neutralizing activity of monoclonal antibodies directed against the N-terminal helical repeat of gp41. <i>Journal of Virology</i> , 2008 , 82, 10032-41	6.6	28
429	Replica exchange simulations of transient encounter complexes in protein-protein association. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008 , 105, 12855-60	11.5	94
428	Global jumping and domain-specific intersegment transfer between DNA cognate sites of the multidomain transcription factor Oct-1. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2008 , 105, 13871-6	11.5	75
427	Impact of phosphorylation on structure and thermodynamics of the interaction between the N-terminal domain of enzyme I and the histidine phosphocarrier protein of the bacterial phosphotransferase system. <i>Journal of Biological Chemistry</i> , 2008 , 283, 18980-9	5.4	28
426	Solution NMR structures of productive and non-productive complexes between the A and B domains of the cytoplasmic subunit of the mannose transporter of the Escherichia coli phosphotransferase system. <i>Journal of Biological Chemistry</i> , 2008 , 283, 11024-37	5.4	22
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