

# JÃ³zef R Lewandowski

## List of Publications by Year in descending order

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55  
papers

3,287  
citations

147801

31  
h-index

155660

55  
g-index

59  
all docs

59  
docs citations

59  
times ranked

3051  
citing authors

#	ARTICLE	IF	CITATIONS
1	Dynamic Nuclear Polarization of Amyloidogenic Peptide Nanocrystals: GNNQQNY, a Core Segment of the Yeast Prion Protein Sup35p. <i>Journal of the American Chemical Society</i> , 2006, 128, 10840-10846.	13.7	255
2	Direct observation of hierarchical protein dynamics. <i>Science</i> , 2015, 348, 578-581.	12.6	222
3	The SARS-COV-2 Spike Protein Binds Sialic Acids and Enables Rapid Detection in a Lateral Flow Point of Care Diagnostic Device. <i>ACS Central Science</i> , 2020, 6, 2046-2052.	11.3	222
4	Proton assisted recoupling and protein structure determination. <i>Journal of Chemical Physics</i> , 2008, 129, 245101.	3.0	183
5	Solid-State NMR Study of Amyloid Nanocrystals and Fibrils Formed by the Peptide GNNQQNY from Yeast Prion Protein Sup35p. <i>Journal of the American Chemical Society</i> , 2007, 129, 5117-5130.	13.7	177
6	Proton Assisted Insensitive Nuclei Cross Polarization. <i>Journal of the American Chemical Society</i> , 2007, 129, 728-729.	13.7	163
7	Huntingtin exon 1 fibrils feature an interdigitated $\beta^2$ -hairpin-based polyglutamine core. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2016, 113, 1546-1551.	7.1	143
8	Enhanced Resolution and Coherence Lifetimes in the Solid-State NMR Spectroscopy of Perdeuterated Proteins under Ultrafast Magic-Angle Spinning. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2205-2211.	4.6	123
9	High-Resolution Solid-State NMR Structure of a 17.6 kDa Protein. <i>Journal of the American Chemical Society</i> , 2010, 132, 1032-1040.	13.7	117
10	Site-Specific Measurement of Slow Motions in Proteins. <i>Journal of the American Chemical Society</i> , 2011, 133, 16762-16765.	13.7	105
11	Structural Complexity of a Composite Amyloid Fibril. <i>Journal of the American Chemical Society</i> , 2011, 133, 14686-14698.	13.7	88
12	Advances in Solid-State Relaxation Methodology for Probing Site-Specific Protein Dynamics. <i>Accounts of Chemical Research</i> , 2013, 46, 2018-2027.	15.6	88
13	Measurement of Site-Specific $^{13}\text{C}$ Spin Lattice Relaxation in a Crystalline Protein. <i>Journal of the American Chemical Society</i> , 2010, 132, 8252-8254.	13.7	80
14	$^1\text{H}$ line width dependence on MAS speed in solid state NMR – Comparison of experiment and simulation. <i>Journal of Magnetic Resonance</i> , 2018, 291, 32-39.	2.1	80
15	Conformational Dynamics of a Seven Transmembrane Helical Protein Anabaena Sensory Rhodopsin Probed by Solid-State NMR. <i>Journal of the American Chemical Society</i> , 2014, 136, 2833-2842.	13.7	78
16	Solid-State NMR of a Protein in a Precipitated Complex with a Full-Length Antibody. <i>Journal of the American Chemical Society</i> , 2014, 136, 16800-16806.	13.7	73
17	Structural Characterization of GNNQQNY Amyloid Fibrils by Magic Angle Spinning NMR. <i>Biochemistry</i> , 2010, 49, 9457-9469.	2.5	66
18	Proton Assisted Recoupling at High Spinning Frequencies. <i>Journal of Physical Chemistry B</i> , 2009, 113, 9062-9069.	2.6	63

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19	Broadband Homonuclear Correlation Spectroscopy at High Magnetic Fields and MAS Frequencies. <i>Journal of the American Chemical Society</i> , 2006, 128, 1776-1777.	13.7	59
20	<sup>15</sup> Nâ <sup>15</sup> N Proton Assisted Recoupling in Magic Angle Spinning NMR. <i>Journal of the American Chemical Society</i> , 2009, 131, 5769-5776.	13.7	56
21	Spin dynamics in the modulation frame: Application to homonuclear recoupling in magic angle spinning solid-state NMR. <i>Journal of Chemical Physics</i> , 2008, 128, 124503.	3.0	50
22	Intermolecular Interactions and Protein Dynamics by Solid-State NMR Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 15374-15378.	13.8	50
23	Heteronuclear proton assisted recoupling. <i>Journal of Chemical Physics</i> , 2011, 134, 095101.	3.0	48
24	Atomic-Resolution Structural Dynamics in Crystalline Proteins from NMR and Molecular Simulation. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 3657-3662.	4.6	47
25	Anisotropic Collective Motion Contributes to Nuclear Spin Relaxation in Crystalline Proteins. <i>Journal of the American Chemical Society</i> , 2010, 132, 1246-1248.	13.7	43
26	Binding of Distinct Substrate Conformations Enables Hydroxylation of Remote Sites in Thaxtomin D by Cytochrome P450 TxtC. <i>Journal of the American Chemical Society</i> , 2019, 141, 216-222.	13.7	42
27	Structural basis for chain release from the enacyloxin polyketide synthase. <i>Nature Chemistry</i> , 2019, 11, 913-923.	13.6	39
28	High-resolution and sensitivity through-bond correlations in ultra-fast magic angle spinning (MAS) solid-state NMR. <i>Chemical Science</i> , 2011, 2, 345-348.	7.4	38
29	Unraveling the complexity of protein backbone dynamics with combined <sup>13</sup> C and <sup>15</sup> N solid-state NMR relaxation measurements. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 21997-22008.	2.8	37
30	Characterization of Proteinâ€“Protein Interfaces in Large Complexes by Solid-State NMR Solvent Paramagnetic Relaxation Enhancements. <i>Journal of the American Chemical Society</i> , 2017, 139, 12165-12174.	13.7	35
31	Fibrillar vs Crystalline Full-Length Î²-2-Microglobulin Studied by High-Resolution Solid-State NMR Spectroscopy. <i>Journal of the American Chemical Society</i> , 2010, 132, 5556-5557.	13.7	32
32	Multipole-multimode Floquet theory of rotational resonance width experiments: <sup>13</sup> Câ€“ <sup>13</sup> C distance measurements in uniformly labeled solids. <i>Journal of Chemical Physics</i> , 2006, 124, 214107.	3.0	31
33	Mechanism of intersubunit ketosynthaseâ€“dehydratase interaction in polyketide synthases. <i>Nature Chemical Biology</i> , 2018, 14, 270-275.	8.0	31
34	Proteinâ€“protein interactions in <i>trans</i> -AT polyketide synthases. <i>Natural Product Reports</i> , 2018, 35, 1097-1109.	10.3	29
35	Solid-State NMR Provides Evidence for Small-Amplitude Slow Domain Motions in a Multispanning Transmembrane Î±-Helical Protein. <i>Journal of the American Chemical Society</i> , 2017, 139, 9246-9258.	13.7	27
36	Structural studies suggest aggregation as one of the modes of action for teixobactin. <i>Chemical Science</i> , 2018, 9, 8850-8859.	7.4	24

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37	Benchmark calculations of the shielding constants in the water dimer. <i>Chemical Physics Letters</i> , 2001, 333, 139-145.	2.6	23
38	Quantifying Microsecond Exchange in Large Protein Complexes with Accelerated Relaxation Dispersion Experiments in the Solid State. <i>Scientific Reports</i> , 2019, 9, 11082.	3.3	23
39	Revealing Intermolecular Hydrogen Bonding Structure and Dynamics in a Deep Eutectic Pharmaceutical by Magic-Angle Spinning NMR Spectroscopy. <i>Molecular Pharmaceutics</i> , 2020, 17, 622-631.	4.6	22
40	A suite of solid-state NMR experiments to utilize orphaned magnetization for assignment of proteins using parallel high and low gamma detection. <i>Journal of Magnetic Resonance</i> , 2019, 305, 219-231.	2.1	18
41	MAS NMR Investigation of Molecular Order in an Ionic Liquid Crystal. <i>Journal of Physical Chemistry B</i> , 2020, 124, 4975-4988.	2.6	17
42	Docking domain-mediated subunit interactions in natural product megasynth(et)ases. <i>Journal of Industrial Microbiology and Biotechnology</i> , 2021, 48, .	3.0	17
43	Simultaneous acquisition of homonuclear and heteronuclear long-distance contacts with time-shared third spin assisted recoupling. <i>Journal of Magnetic Resonance</i> , 2012, 218, 30-34.	2.1	16
44	Compensated second-order recoupling: application to third spin assisted recoupling. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 7246.	2.8	15
45	Intermolecular Interactions and Protein Dynamics by Solid-€State NMR Spectroscopy. <i>Angewandte Chemie</i> , 2015, 127, 15594-15598.	2.0	15
46	Probing Protein Dynamics Using Multifield Variable Temperature NMR Relaxation and Molecular Dynamics Simulation. <i>Journal of Physical Chemistry B</i> , 2018, 122, 9697-9702.	2.6	15
47	Communication Breakdown: Dissecting the COM Interfaces between the Subunits of Nonribosomal Peptide Synthetases. <i>ACS Catalysis</i> , 2021, 11, 10802-10813.	11.2	14
48	Modulation of Transmembrane Domain Interactions in Neu Receptor Tyrosine Kinase by Membrane Fluidity and Cholesterol. <i>Journal of Membrane Biology</i> , 2019, 252, 357-369.	2.1	10
49	Isolation and structural characterisation of rhodium( <sup>III</sup> )- <sup>2</sup> -fluoroarene complexes: experimental verification of predicted regioselectivity. <i>Dalton Transactions</i> , 2020, 49, 5791-5793.	3.3	7
50	Taming the dynamics in a pharmaceutical by cocrystallization: investigating the impact of the cofomer by solid-state NMR. <i>CrystEngComm</i> , 2021, 23, 6859-6870.	2.6	7
51	Accelerating <sup>15</sup> N and <sup>13</sup> C R1 and R1ρ-relaxation measurements by multiple pathway solid-state NMR experiments. <i>Journal of Magnetic Resonance</i> , 2021, 331, 107049.	2.1	5
52	Molecular basis for acyl carrier protein-€ketoreductase interaction in <i>trans</i> -acyltransferase polyketide synthases. <i>Chemical Science</i> , 2021, 12, 13676-13685.	7.4	3
53	Simultaneous MQMAS NMR Experiments for Two Half-Integer Quadrupolar Nuclei. <i>Journal of Magnetic Resonance</i> , 2020, 320, 106831.	2.1	2
54	Dipolar Order Parameters in Large Systems With Fast Spinning. <i>Frontiers in Molecular Biosciences</i> , 2021, 8, 791026.	3.5	2

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55	Solid State NMR Studies Of Structural And Motional Complexity In Amyloid-Like Fibrils Of The Peptide GNNQQNY. Biophysical Journal, 2009, 96, 219a.	0.5	0