

Piotr Paluch

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Study of Intermolecular Interactions in the Corrole Matrix by Solid-State NMR under 100 kHz MAS and Theoretical Calculations. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 14108-14111.	13.8	86
2	Simple and accurate determination of ^{13}C - ^1H distances under ultra-fast MAS NMR. <i>Journal of Magnetic Resonance</i> , 2013, 233, 56-63.	2.1	59
3	^1H -Detected Biomolecular NMR under Fast Magic-Angle Spinning. <i>Chemical Reviews</i> , 2022, 122, 9943-10018.	47.7	51
4	Ibuprofen in Mesopores of Mobil Crystalline Material 41 (MCM-41): A Deeper Understanding. <i>Molecular Pharmaceutics</i> , 2014, 11, 1512-1519.	4.6	45
5	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.	13.7	41
6	Dynamic Motion of Organic Spacer Cations in Ruddlesden-Popper Lead Iodide Perovskites Probed by Solid-State NMR Spectroscopy. <i>Chemistry of Materials</i> , 2021, 33, 642-656.	6.7	33
7	NMR Study of BA/FBA Cocrystal Confined Within Mesoporous Silica Nanoparticles Employing Thermal Solid Phase Transformation. <i>Journal of Physical Chemistry C</i> , 2015, 119, 8652-8661.	3.1	27
8	Fine refinement of solid state structure of racemic form of phospho-tyrosine employing NMR Crystallography approach. <i>Solid State Nuclear Magnetic Resonance</i> , 2015, 65, 2-11.	2.3	24
9	Insights into the Tautomerism in <i>meso</i> -Substituted Corroles: A Variable Temperature ^1H , ^{13}C , ^{15}N , and ^{19}F -NMR Spectroscopy Study. <i>Chemistry - A European Journal</i> , 2014, 20, 1720-1730.	3.3	21
10	A DFT Study of the Kinetic Isotope Effects on the Competing $\text{S}_{\text{N}}2$ and E2 Reactions between Hypochlorite Anion and Ethyl Chloride. <i>Journal of Chemical Theory and Computation</i> , 2009, 5, 33-36.	5.3	20
11	Analysis of local molecular motions of aromatic sidechains in proteins by 2D and 3D fast MAS NMR spectroscopy and quantum mechanical calculations. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 28789-28801.	2.8	19
12	Approach toward the Understanding of Coupling Mechanism for EDC Reagent in Solvent-Free Mechanosynthesis. <i>Organic Letters</i> , 2017, 19, 5360-5363.	4.6	19
13	Crystal structure determination of an elusive methanol solvate hydrate of catechin using crystal structure prediction and NMR crystallography. <i>CrystEngComm</i> , 2020, 22, 4969-4981.	2.6	19
14	Imaging the spatial distribution of radiofrequency field, sample and temperature in MAS NMR rotor. <i>Solid State Nuclear Magnetic Resonance</i> , 2017, 87, 137-142.	2.3	17
15	Application of ^1H and ^{27}Al magic angle spinning solid state NMR at 60 kHz for studies of Au and Au-Ni catalysts supported on boehmite/alumina. <i>Solid State Nuclear Magnetic Resonance</i> , 2017, 84, 111-117.	2.3	15
16	Evaluation of excitation schemes for indirect detection of ^{14}N via solid-state HMQC NMR experiments. <i>Journal of Magnetic Resonance</i> , 2019, 303, 28-41.	2.1	15
17	Understanding the formation of apremilast cocrystals. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2019, 75, 803-814.	1.1	15
18	Study of host-guest interactions in benzodiazacoronands by means of solid state NMR spectroscopy, X-ray diffraction and quantum mechanical computations. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 6423.	2.8	14

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19	Synthesis, characterization, and catalytic activity for thioanisole oxidation of homogeneous and heterogeneous binuclear manganese(II) complexes with amino acid-based ligands. <i>Transition Metal Chemistry</i> , 2013, 38, 511-521.	1.4	14
20	Study of the thermal processes in molecular crystals of peptides by means of NMR crystallography. <i>CrystEngComm</i> , 2013, 15, 8680.	2.6	14
21	Analysis of HMQC experiments applied to a spin $\hat{A}^{1/2}$ nucleus subject to very large CSA. <i>Solid State Nuclear Magnetic Resonance</i> , 2019, 100, 11-25.	2.3	14
22	Crystal structures of two furazidin polymorphs revealed by a joint effort of crystal structure prediction and NMR crystallography. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2020, 76, 322-335.	1.1	13
23	Photosensitive nanocapsules for use in imaging from poly(styrene-co-divinylbenzene) cross-linked with coumarin derivatives. <i>Colloids and Surfaces B: Biointerfaces</i> , 2013, 111, 571-578.	5.0	12
24	The structure of cyclolinopeptide A in chloroform refined by RDC measurements. <i>Journal of Peptide Science</i> , 2014, 20, 901-907.	1.4	11
25	In Depth Analysis of Chiroptical Properties of Enones Derived from Abietic Acid. <i>Journal of Organic Chemistry</i> , 2018, 83, 3547-3561.	3.2	8
26	Combined solid state NMR and ONIOM studies of reversible crystalline phase reaction for nickel coordination compounds. <i>Solid State Nuclear Magnetic Resonance</i> , 2009, 36, 103-109.	2.3	7
27	Stereoselective cyclopropyl phosphonate formation using (S)-dimethylsulfonium-(p-tolylsulfanyl)methylide. Unusual phosphoryl group migration. <i>Tetrahedron Letters</i> , 2013, 54, 223-226.	1.4	7
28	Recent Progress in the Solid-State NMR Studies of Short Peptides. <i>Annual Reports on NMR Spectroscopy</i> , 2014, , 67-143.	1.5	7
29	Full Characterization of Linezolid and Its Synthetic Precursors by Solid-State Nuclear Magnetic Resonance Spectroscopy and Mass Spectrometry. <i>Journal of Pharmaceutical Sciences</i> , 2015, 104, 3883-3892.	3.3	7
30	A Multi-Technique Experimental and Computational Approach To Study the Dehydration Processes in the Crystals of Endomorphin Opioid Peptide Derivative. <i>Crystal Growth and Design</i> , 2016, 16, 5312-5322.	3.0	7
31	$\text{I}^{\ominus}\text{E}$ -Phylic Molecular Recognition in the Solid State as a Driving Force for Mechanochemical Formation of Apremilast Solvates and Cocrystals. <i>Crystal Growth and Design</i> , 2018, 18, 3959-3970.	3.0	7
32	Chiral crystals from porphyrinoids possessing a very low racemization barrier. <i>CrystEngComm</i> , 2016, 18, 3561-3565.	2.6	6
33	^1H - ^{31}P CPVC NMR method under Very Fast Magic Angle Spinning for analysis of dipolar interactions and dynamics processes in the crystalline phosphonium tetrafluoroborate salts. <i>Solid State Nuclear Magnetic Resonance</i> , 2017, 87, 96-103.	2.3	6
34	New synthetic pathway leading to oxospirochlorins. <i>RSC Advances</i> , 2018, 8, 21354-21362.	3.6	6
35	Magic angle spinning NMR study of interaction of N-terminal sequence of dermorphin (Tyr-d-Ala-Phe-Gly) with phospholipids. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2012, 1818, 2579-2587.	2.6	5
36	NMR Assignment of Methyl Groups in Immobilized Proteins Using Multiple-Bond ^{13}C Homonuclear Transfers, Proton Detection, and Very Fast MAS. <i>Frontiers in Molecular Biosciences</i> , 2022, 9, 828785.	3.5	5

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37	Spontaneous Keto↔Enol Tautomerization in the Crystal Lattice Visualized with the Help of Water Encapsulated in Hydrophilic Reservoirs. <i>ChemPhysChem</i> , 2017, 18, 2850-2854.	2.1	4
38	Study of the Mechanism of Thermal Chemical Processes in the Crystals of YAF Tripeptides by Means of Mass Spectrometry and Solid State NMR. <i>Journal of Physical Chemistry B</i> , 2013, 117, 13481-13489.	2.6	3
39	Simple and Robust Study of Backbone Dynamics of Crystalline Proteins Employing ¹ H- ¹⁵ N Dipolar Coupling Dispersion. <i>Journal of Physical Chemistry B</i> , 2018, 122, 8146-8156.	2.6	3
40	Influence of Environmental Humidity on Organization and Molecular Dynamics of Heteromacrocyclic Assemblies. <i>Journal of Physical Chemistry B</i> , 2013, 117, 14420-14431.	2.6	2
41	Slow and Very Fast MAS Solid State NMR Study of Biopolymers. <i>Macromolecular Symposia</i> , 2014, 339, 60-69.	0.7	2
42	The ¹ H, ¹³ C, ¹⁵ N, and ¹⁹ F NMR chemical shifts assignments in 5,10,15-tris (pentafluorophenyl)tetra- ¹⁵ N corrole at 191‰K. <i>Magnetic Resonance in Chemistry</i> , 2015, 53, 167-171.	1.9	1
43	Hadamard acquisition of 13 C- ¹³ C 2D correlation NMR spectra. <i>Magnetic Resonance in Chemistry</i> , 2021, 59, 247-256.	1.9	1
44	The influence of the stereochemistry and C-end chemical modification of dermorphin derivatives on the peptide-phospholipid interactions. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2020, 1862, 183066.	2.6	0