

# Sławomir Szymański

## List of Publications by Year in descending order

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22  
papers

163  
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1307594

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1199594

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23  
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23  
docs citations

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times ranked

184  
citing authors

#	ARTICLE	IF	CITATIONS
1	Structure and NMR Spectra of Some [2.2]Paracyclophanes. The Dilemma of [2.2]Paracyclophane Symmetry. <i>Journal of Physical Chemistry A</i> , 2011, 115, 10638-10649.	2.5	30
2	Wave Properties of a Methyl Group under Ambient Conditions. <i>Physical Review Letters</i> , 2002, 89, 023004.	7.8	27
3	Structure and Dynamics of [3.3]Paracyclophane As Studied by Nuclear Magnetic Resonance and Density Functional Theory Calculations. <i>Journal of Physical Chemistry A</i> , 2010, 114, 10467-10473.	2.5	21
4	Insights into the Tautomerism in <i>meso</i> -Substituted Corroles: A Variable-Temperature <sup>1</sup> H, <sup>13</sup> C, <sup>15</sup> N, and <sup>19</sup> F...NMR Spectroscopy Study. <i>Chemistry - A European Journal</i> , 2014, 20, 1720-1730.	3.3	21
5	Structure, NMR and Electronic Spectra of [ <i>m.n</i> ]Paracyclophanes with Varying Bridges Lengths ( <i>m, n</i> = 2-4). <i>Journal of Physical Chemistry A</i> , 2016, 120, 724-736.	2.5	10
6	NMR properties of N <sub>3</sub> <sup>-</sup> . A comparison of theory and experiment. <i>Chemical Physics Letters</i> , 1995, 243, 144-150.	2.6	9
7	High-temperature quantum mechanical exchange of heavy particles. Effective spin Hamiltonian. <i>Journal of Molecular Structure</i> , 1994, 321, 115-124.	3.6	7
8	Dynamics of [n.3]paracyclophanes (n = 2 - 4) as studied by NMR. Obtaining separate Arrhenius parameters for two dynamic processes in [4.3]paracyclophane. <i>Journal of Physical Organic Chemistry</i> , 2013, 26, 596-600.	1.9	6
9	Cross-Correlation of Nuclear Quadrupolar Interactions as a Probe in Structure Elucidation: A Liquid-Phase Nuclear Magnetic Resonance Studies on Bis(hexamethyldisilylamido)mercury(II). <i>Journal of Physical Chemistry A</i> , 2001, 105, 6414-6419.	2.5	5
10	Hindered Rotation of the Silyl Group in Liquid-Phase NMR Spectra of 9-Silyltryptene Derivatives: A Comparison with the Methyl Analogues. <i>Journal of Physical Chemistry A</i> , 2008, 112, 8612-8616.	2.5	5
11	A quantum mechanical alternative to the Arrhenius equation in the interpretation of proton spin-lattice relaxation data for the methyl groups in solids. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 28866-28878.	2.8	5
12	Structure and NMR spectra of cyclophanes with unsaturated bridges (cyclophenes). <i>Magnetic Resonance in Chemistry</i> , 2012, 50, 449-457.	1.9	4
13	Nonclassical dynamics of the methyl group in 1,1,1-triphenylethane. Evidence from powder <sup>1</sup> H NMR spectra. <i>Journal of Chemical Physics</i> , 2017, 146, 104504.	3.0	4
14	Proton Tunnelling Effects in Metal Hydride NMR. <i>Annual Reports on NMR Spectroscopy</i> , 1998, , 1-54.	1.5	3
15	Synthesis and structural characterization of exemplary silyl triptycenes. <i>New Journal of Chemistry</i> , 2019, 43, 7567-7573.	2.8	2
16	Cross correlations of quadrupolar interactions and scalar relaxation of the second kind. <i>Journal of Magnetic Resonance</i> , 1990, 87, 166-168.	0.5	1
17	Blue-Shift Hydrogen Bonds in Silyltryptene Derivatives: Antibonding $\sigma^*$ Orbitals of the Si-C Bond as Effective Acceptors of Electron Density. <i>ChemPhysChem</i> , 2020, 21, 540-545.	2.1	1
18	Principles of NMR Spectroscopy. , 2018, , 7-83.		0

#	ARTICLE	IF	CITATIONS
19	Quantum Molecular Dynamics in Liquid-Phase Stick NMR Spectra. , 2018, , 333-348.		0
20	Quantum Mechanical Rate Processes in NMR Spectra. , 2018, , 349-389.		0
21	Rotational Tunneling in Stick NMR Spectra of Solids. , 2018, , 305-331.		0
22	Discrete Molecular Dynamics and NMR Line Shape Effects. Intramolecular Exchange. , 2018, , 195-248.		0