

# Peter A Beckmann

## List of Publications by Year in descending order

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54  
papers

1,033  
citations

516710

16  
h-index

454955

30  
g-index

55  
all docs

55  
docs citations

55  
times ranked

785  
citing authors

#	ARTICLE	IF	CITATIONS
1	Spectral densities and nuclear spin relaxation in solids. <i>Physics Reports</i> , 1988, 171, 85-128.	25.6	242
2	A Thermometer for Nonspinning Solid-State NMR Spectroscopy. <i>Journal of Magnetic Resonance</i> , 2000, 146, 379-380.	2.1	140
3	Nuclear-spin relaxation in molecular solids with reorienting methyl and n-butyl groups: The spectral density and the state of the solid. <i>Physical Review B</i> , 1988, 38, 11098-11111.	3.2	31
4	<sup>1</sup> H nuclear magnetic resonance spin-lattice relaxation, <sup>13</sup> C magic-angle-spinning nuclear magnetic resonance spectroscopy, differential scanning calorimetry, and x-ray diffraction of two polymorphs of 2,6-di-tert-butyl-naphthalene. <i>Journal of Chemical Physics</i> , 2000, 113, 1958-1965.	3.0	30
5	<sup>127</sup> I and <sup>207</sup> Pb Solid-State NMR Spectroscopy and Nuclear Spin Relaxation in PbI <sub>2</sub> : A Preliminary Study. <i>Journal of Physical Chemistry C</i> , 2014, 118, 9143-9153.	3.1	29
6	Solid state phase transitions and molecular reorientation in ortho- and para-carborane: An isomer effect. <i>Journal of Chemical Physics</i> , 1980, 72, 4600-4607.	3.0	28
7	A New Mechanism for Spin-Lattice Relaxation of Heavy Nuclei in the Solid State: <sup>207</sup> Pb Relaxation in Lead Nitrate. <i>Journal of the American Chemical Society</i> , 2001, 123, 7094-7100.	13.7	28
8	CF <sub>3</sub> Rotation in 3-(Trifluoromethyl)phenanthrene. X-ray Diffraction and ab Initio Electronic Structure Calculations. <i>Journal of Physical Chemistry A</i> , 2006, 110, 3954-3960.	2.5	27
9	Methyl group rotation, <sup>1</sup> H spin-lattice relaxation in an organic solid, and the analysis of nonexponential relaxation. <i>Journal of Chemical Physics</i> , 2012, 136, 054508.	3.0	27
10	Intramolecular and Intermolecular Contributions to the Barriers for Rotation of Methyl Groups in Crystalline Solids: Electronic Structure Calculations and Solid-State NMR Relaxation Measurements. <i>Journal of Organic Chemistry</i> , 2011, 76, 5170-5176.	3.2	26
11	Solid state proton spin relaxation in ethylbenzenes: Methyl reorientation barriers and molecular structure. <i>Journal of Chemical Physics</i> , 1991, 95, 828-835.	3.0	25
12	Methyl and n-butyl group reorientation in planar aromatic solids: Low-frequency nuclear magnetic resonance relaxometry and x-ray diffraction. <i>Journal of Chemical Physics</i> , 2003, 118, 11129-11138.	3.0	20
13	CF <sub>3</sub> Rotation in 3-(Trifluoromethyl)phenanthrene: Solid State <sup>19</sup> F and <sup>1</sup> H NMR Relaxation and Bloch-Wangsness-Redfield Theory. <i>Journal of Physical Chemistry A</i> , 2006, 110, 3947-3953.	2.5	19
14	Superpositions of intramolecular reorientations and nuclear spin relaxation. <i>Molecular Physics</i> , 1980, 41, 1227-1238.	1.7	17
15	Solid state proton spin relaxation and methyl and n-butyl reorientation. <i>Journal of Chemical Physics</i> , 1994, 100, 752-753.	3.0	17
16	Methyl reorientation in solid 3-ethylchrysene and 3-isopropylchrysene. <i>Solid State Nuclear Magnetic Resonance</i> , 1998, 12, 251-256.	2.3	17
17	Proton spin-lattice relaxation in meta-carborane. <i>Journal of Chemical Physics</i> , 1980, 73, 3514-3515.	3.0	16
18	<sup>111</sup> Cd and <sup>113</sup> Cd spin-lattice relaxation in CdMoO <sub>4</sub> by paramagnetic centers in the absence of spin diffusion. <i>Physical Review B</i> , 2005, 71, .	3.2	16

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19	Spin-lattice relaxation of heavy spin-1/2 nuclei in diamagnetic solids: A Raman process mediated by spin-rotation interaction. <i>Physical Review B</i> , 2006, 74, .	3.2	16
20	Methoxy and Methyl Group Rotation: Solid-state NMR <sup>1</sup> H Spin-Lattice Relaxation, Electronic Structure Calculations, X-ray Diffraction, and Scanning Electron Microscopy. <i>ChemPhysChem</i> , 2015, 16, 1509-1519.	2.1	15
21	The relationship between crystal structure and methyl and t-butyl group dynamics in van der Waals organic solids. <i>Journal of Chemical Physics</i> , 2004, 120, 5309-5314.	3.0	14
22	Distributions of methyl group rotational barriers in polycrystalline organic solids. <i>Journal of Chemical Physics</i> , 2013, 139, 204501.	3.0	14
23	Superlattices, polymorphs and solid-state NMR spin-lattice relaxation (T <sub>1</sub> ) measurements of 2,6-di-tert-butyl-naphthalene. <i>Chemical Communications</i> , 2000, , 651-652.	4.1	13
24	The quenching of isopropyl group rotation in van der Waals molecular solids. <i>Journal of Chemical Physics</i> , 2008, 128, 124502.	3.0	13
25	Dipole-dipole spin relaxation in solids. <i>Physica B: Condensed Matter</i> , 1993, 190, 267-284.	2.7	12
26	Single-Crystal X-ray Diffraction, Isolated-Molecule and Cluster Electronic Structure Calculations, and Scanning Electron Microscopy in an Organic solid: Models for Intramolecular Motion in 4,4'-Dimethoxybiphenyl. <i>ChemPhysChem</i> , 2012, 13, 2082-2089.	2.1	12
27	Nonexponential <sup>1</sup> H spin-lattice relaxation and methyl group rotation in molecular solids. <i>Solid State Nuclear Magnetic Resonance</i> , 2015, 71, 91-95.	2.3	12
28	The electron-methyl group spin-spin interaction. <i>Molecular Physics</i> , 1977, 34, 665-680.	1.7	11
29	Solid state proton spin-lattice relaxation in four structurally related organic molecules. <i>Chemical Physics</i> , 2003, 290, 241-250.	1.9	10
30	<sup>119</sup> Sn spin-lattice relaxation in $\text{SnF}_2$ . <i>Physical Review B</i> , 2009, 79, .	3.2	10
31	A proton spin-lattice relaxation rate study of methyl and t-butyl group reorientation in the solid state. <i>Solid State Nuclear Magnetic Resonance</i> , 2010, 38, 31-35.	2.3	9
32	Nonexponential Solid State <sup>1</sup> H and <sup>19</sup> F Spin-Lattice Relaxation, Single-crystal X-ray Diffraction, and Isolated-Molecule and Cluster Electronic Structure Calculations in an Organic Solid: Coupled Methyl Group Rotation and Methoxy Group Libration in 4,4'-Dimethoxyoctafluorobiphenyl. <i>Journal of Physical Chemistry A</i> , 2012, 116, 11946-11956.	2.5	9
33	Proton spin-lattice relaxation in MBDBP. <i>Molecular Physics</i> , 1980, 41, 1239-1258.	1.7	8
34	Proton spin relaxation, internal motion and structure in solid 1,2,4,5-tetraisopropylbenzene. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1993, 89, 3801.	1.7	8
35	<sup>207</sup> Pb spin-lattice relaxation in solid $\text{PbMoO}_4$ and $\text{PbCl}_2$ . <i>Physical Review B</i> , 2006, 74, .	3.2	8
36	Methyl and t-butyl reorientation in an organic molecular solid. <i>Solid State Nuclear Magnetic Resonance</i> , 2009, 36, 86-91.	2.3	8

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37	Solid state <sup>1</sup> H spin-lattice relaxation and isolated-molecule and cluster electronic structure calculations in organic molecular solids: The relationship between structure and methyl group and t-butyl group rotation. <i>Journal of Chemical Physics</i> , 2014, 140, 194304.	3.0	8
38	Solid-State Proton Spin Relaxation and Methyl Reorientation in Isopropylbenzene. <i>The Journal of Physical Chemistry</i> , 1995, 99, 391-394.	2.9	7
39	Unusual proton Zeeman spin relaxation in an organic solid: several crystal polymorphs or different glass structures?. <i>Solid State Nuclear Magnetic Resonance</i> , 2000, 16, 239-244.	2.3	7
40	The relationship between crystal structure and NMR relaxation in molecular solids with tert-butyl groups. <i>Chemical Physics</i> , 2008, 345, 116-118.	1.9	7
41	<sup>1</sup> H and <sup>19</sup> F spin-lattice relaxation and CH <sub>3</sub> or CF <sub>3</sub> reorientation in molecular solids containing both H and F atoms. <i>Journal of Chemical Physics</i> , 2016, 144, 154308.	3.0	7
42	Solid state proton spin relaxation and t-butyl and methyl group reorientation in 1-bromo-2,4,6-trimethylbenzene. <i>Journal of Chemical Physics</i> , 1991, 95, 4778-4782.	3.0	6
43	Methyl Group Rotation and <sup>1</sup> H and <sup>2</sup> H Zeeman Relaxation in Organic Solids. <i>Journal of Physical Chemistry A</i> , 2001, 105, 7350-7355.	2.5	6
44	Methyl and t-butyl group rotation in a molecular solid: <sup>1</sup> H NMR spin-lattice relaxation and X-ray diffraction. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 1720-1726.	2.8	6
45	Solid Phase Transitions and tert-Butyl and Methyl Group Rotation in an Organic Solid: X-ray Diffractometry, Differential Scanning Calorimetry, and Solid-State <sup>1</sup> H Nuclear Spin Relaxation. <i>Journal of Physical Chemistry A</i> , 2017, 121, 6220-6230.	2.5	6
46	Proton spin-lattice relaxation and intramolecular reorientation in solids. <i>Journal of Magnetic Resonance</i> , 1984, 59, 63-70.	0.5	5
47	Proton Spin Lattice Relaxation in Organic Molecular Solids: Polymorphism and the Dependence on Sample Preparation. <i>ChemPhysChem</i> , 2018, 19, 2423-2436.	2.1	3
48	Concomitant Polymorphism in an Organic Solid: Molecular and Crystal Structure and Intra- and Intermolecular Potential Contributions to tert-Butyl and Methyl Group Rotation. <i>ChemPhysChem</i> , 2019, 20, 2887-2894.	2.1	3
49	Proton Spin Relaxation and Thermal History Effects in Organic Molecular Solids. , 1992, , 357-362.		3
50	Physics in Elementary School. <i>MRS Bulletin</i> , 1992, 17, 47-48.	3.5	2
51	More on g?. <i>American Journal of Physics</i> , 1987, 55, 969-969.	0.7	0
52	Monitoring a simple hydrolysis process in an organic solid by observing methyl group rotation. <i>Solid State Nuclear Magnetic Resonance</i> , 2017, 85-86, 1-11.	2.3	0
53	Note: Methyl and t-butyl group rotation in van der Waals solids. <i>Journal of Chemical Physics</i> , 2018, 148, 106101.	3.0	0
54	Solid state proton spin-lattice relaxation in polycrystalline methylphenanthrenes. IV. 1,4-dimethylphenanthrene. <i>Journal of Chemical Physics</i> , 2019, 150, 124508.	3.0	0