

Sharon E Ashbrook

List of Publications by Year in descending order

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167
papers

7,907
citations

44069

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179
all docs

179
docs citations

179
times ranked

6439
citing authors

#	ARTICLE	IF	CITATIONS
1	Origin of the temperature dependence of ¹³ C pNMR shifts for copper paddlewheel MOFs. <i>Chemical Science</i> , 2022, 13, 2674-2685.	7.4	2
2	Solid-state NMR spectroscopy. <i>Nature Reviews Methods Primers</i> , 2021, 1, .	21.2	196
3	Thermal Dehydrofluorination of GaPO-34 Revealed by NMR Crystallography. <i>Journal of Physical Chemistry C</i> , 2021, 125, 2537-2545.	3.1	5
4	¹⁷ O NMR spectroscopy of crystalline microporous materials. <i>Chemical Science</i> , 2021, 12, 5016-5036.	7.4	33
5	Exploring cation disorder in mixed-metal pyrochlore ceramics using ¹⁷ O NMR spectroscopy and first-principles calculations. <i>Magnetic Resonance in Chemistry</i> , 2021, 59, 961-974.	1.9	0
6	Formation Mechanism and Porosity Development in Porous Boron Nitride. <i>Journal of Physical Chemistry C</i> , 2021, 125, 27429-27439.	3.1	15
7	Facile, Room-Temperature ¹⁷ O Enrichment of Zeolite Frameworks Revealed by Solid-State NMR Spectroscopy. <i>Journal of the American Chemical Society</i> , 2020, 142, 900-906.	13.7	48
8	Phase Distribution, Composition, and Disorder in Y ₂ (Hf,Sn) ₂ O ₇ Ceramics: Insights from Solid-State NMR Spectroscopy and First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2020, 124, 17073-17084.	3.1	7
9	Application of NMR Crystallography to Highly Disordered Templated Materials: Extensive Local Structural Disorder in the Gallophosphate GaPO-34A. <i>Inorganic Chemistry</i> , 2020, 59, 11616-11626.	4.0	9
10	Site-Specific Iron Substitution in STA-28, a Large Pore Aluminophosphate Zeotype Prepared by Using 1,10-Phenanthrolines as Framework-Bound Templates. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 15186-15190.	13.8	4
11	Mechanochemically assisted hydrolysis in the ADOR process. <i>Chemical Science</i> , 2020, 11, 7060-7069.	7.4	12
12	Site-Specific Iron Substitution in STA-28, a Large Pore Aluminophosphate Zeotype Prepared by Using 1,10-Phenanthrolines as Framework-Bound Templates. <i>Angewandte Chemie</i> , 2020, 132, 15298-15302.	2.0	2
13	Following the unusual breathing behaviour of ¹⁷ O-enriched mixed-metal (Al,Ga)-MIL-53 using NMR crystallography. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 14514-14526.	2.8	16
14	Synthesis and Polymorphism of Mixed Aluminum-Gallium Oxides. <i>Inorganic Chemistry</i> , 2020, 59, 3805-3816.	4.0	28
15	Phosphorus-Bismuth Peri-Substituted Acenaphthenes: A Synthetic, Structural, and Computational Study. <i>Inorganic Chemistry</i> , 2020, 59, 5616-5625.	4.0	13
16	Fast room temperature lability of aluminosilicate zeolites. <i>Nature Communications</i> , 2019, 10, 4690.	12.8	75
17	Ensemble-Based Modeling of the NMR Spectra of Solid Solutions: Cation Disorder in Y ₂ (Sn,Ti) ₂ O ₇ . <i>Journal of the American Chemical Society</i> , 2019, 141, 17838-17846.	13.7	29
18	Visualization of the effect of additives on the nanostructures of individual bio-inspired calcite crystals. <i>Chemical Science</i> , 2019, 10, 1176-1185.	7.4	26

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19	A Picture of Disorder in Hydrated Wadsleyite Under the Combined Microscope of Solid-State NMR Spectroscopy and Ab Initio Random Structure Searching. <i>Journal of the American Chemical Society</i> , 2019, 141, 3024-3036.	13.7	13
20	STA-27, a porous Lewis acidic scandium MOF with an unexpected topology type prepared with 2,3,5,6-tetrakis(4-carboxyphenyl)pyrazine. <i>Journal of Materials Chemistry A</i> , 2019, 7, 5685-5701.	10.3	22
21	A procedure for identifying possible products in the assembly–disassembly–organization–reassembly (ADOR) synthesis of zeolites. <i>Nature Protocols</i> , 2019, 14, 781-794.	12.0	22
22	¹³ C pNMR of α -crumple zone Cu(II) isophthalate metal-organic frameworks. <i>Solid State Nuclear Magnetic Resonance</i> , 2019, 101, 44-50.	2.3	11
23	NMR chemical shifts of urea loaded copper benzoate. A joint solid-state NMR and DFT study. <i>Solid State Nuclear Magnetic Resonance</i> , 2019, 101, 31-37.	2.3	17
24	Sensitivity improvement in 5QMAS NMR experiments using FAM-N pulses. <i>Solid State Nuclear Magnetic Resonance</i> , 2019, 100, 1-10.	2.3	3
25	Nuclear Magnetic Resonance Spectroscopy as a Dynamical Structural Probe of Hydrogen under High Pressure. <i>Physical Review Letters</i> , 2019, 122, 135501.	7.8	9
26	Rationalization of solid-state NMR multi-pulse decoupling strategies: Coupling of spin $I = 1/2$ and half-integer quadrupolar nuclei. <i>Journal of Magnetic Resonance</i> , 2019, 303, 48-56.	2.1	3
27	Kinetics and Mechanism of the Hydrolysis and Rearrangement Processes within the Assembly–Disassembly–Organization–Reassembly Synthesis of Zeolites. <i>Journal of the American Chemical Society</i> , 2019, 141, 4453-4459.	13.7	21
28	Is the ³¹ P chemical shift anisotropy of aluminophosphates a useful parameter for NMR crystallography?. <i>Magnetic Resonance in Chemistry</i> , 2019, 57, 176-190.	1.9	6
29	¹⁷ O solid-state NMR spectroscopy of A ₂ B ₂ O ₇ oxides: quantitative isotopic enrichment and spectral acquisition?. <i>RSC Advances</i> , 2018, 8, 7089-7101.	3.6	13
30	Recent Advances in Solid-State Nuclear Magnetic Resonance Spectroscopy. <i>Annual Review of Analytical Chemistry</i> , 2018, 11, 485-508.	5.4	45
31	Cost-effective ¹⁷ O enrichment and NMR spectroscopy of mixed-metal terephthalate metal–organic frameworks. <i>Chemical Science</i> , 2018, 9, 850-859.	7.4	49
32	Pressure-induced chemistry for the 2D to 3D transformation of zeolites. <i>Journal of Materials Chemistry A</i> , 2018, 6, 5255-5259.	10.3	21
33	Perspective: Current advances in solid-state NMR spectroscopy. <i>Journal of Chemical Physics</i> , 2018, 149, 040901.	3.0	28
34	Polymorphism, Weak Interactions and Phase Transitions in Chalcogen–Phosphorus Heterocycles. <i>Chemistry - A European Journal</i> , 2018, 24, 11067-11081.	3.3	4
35	Hydrolytic stability in hemilabile metal–organic frameworks. <i>Nature Chemistry</i> , 2018, 10, 1096-1102.	13.6	134
36	Investigating FAM-N pulses for signal enhancement in MQMAS NMR of quadrupolar nuclei. <i>Solid State Nuclear Magnetic Resonance</i> , 2017, 84, 89-102.	2.3	9

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37	Porous zinc and cobalt 2-nitroimidazolate frameworks with six-membered ring windows and a layered cobalt 2-nitroimidazolate polymorph. <i>CrystEngComm</i> , 2017, 19, 1377-1388.	2.6	6
38	A Multinuclear NMR Study of Six Forms of AlPO-34: Structure and Motional Broadening. <i>Journal of Physical Chemistry C</i> , 2017, 121, 1781-1793.	3.1	25
39	Exploiting NMR spectroscopy for the study of disorder in solids. <i>International Reviews in Physical Chemistry</i> , 2017, 36, 39-115.	2.3	65
40	Selective Oxidation and Functionalization of 6-Diphenylphosphinoacenaphthyl-5-tellurenyl Species 6-Ph ₂ P-Ace-5-TeX (X = Mes, Cl, O ₃ SCF ₃). Various Types of P ^{IV} -Te(II,IV) Bonding Situations (E = O, S, Se). <i>Organometallics</i> , 2017, 36, 1566-1579.	2.3	18
41	In situ solid-state NMR and XRD studies of the ADOR process and the unusual structure of zeolite IPC-6. <i>Nature Chemistry</i> , 2017, 9, 1012-1018.	13.6	63
42	An NMR Crystallographic Investigation of the Relationships between the Crystal Structure and ²⁹ Si Isotropic Chemical Shift in Silica Zeolites. <i>Journal of Physical Chemistry C</i> , 2017, 121, 15198-15210.	3.1	28
43	Solid-State NMR Spectroscopy Proves the Presence of Penta-coordinated Sc Sites in MIL-100(Sc). <i>Chemistry - A European Journal</i> , 2017, 23, 9525-9534.	3.3	19
44	Synthesis, Isotopic Enrichment, and Solid-State NMR Characterization of Zeolites Derived from the Assembly, Disassembly, Organization, Reassembly Process. <i>Journal of the American Chemical Society</i> , 2017, 139, 5140-5148.	13.7	42
45	Determining the Surface Structure of Silicated Alumina Catalysts via Isotopic Enrichment and Dynamic Nuclear Polarization Surface-Enhanced NMR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2017, 121, 22977-22984.	3.1	34
46	Calculation and experimental measurement of paramagnetic NMR parameters of phenolic oximate Cu(II) complexes. <i>Chemical Communications</i> , 2017, 53, 10512-10515.	4.1	11
47	Investigation of zeolitic imidazolate frameworks using ¹³ C and ¹⁵ N solid-state NMR spectroscopy. <i>Solid State Nuclear Magnetic Resonance</i> , 2017, 87, 54-64.	2.3	21
48	Effects of Extraframework Species on the Structure-Based Prediction of ³¹ P Isotropic Chemical Shifts of Aluminophosphates. <i>Journal of Physical Chemistry C</i> , 2017, 121, 28065-28076.	3.1	12
49	Ionothermal synthesis and characterization of CoAPO-34 molecular sieve. <i>Microporous and Mesoporous Materials</i> , 2017, 239, 336-341.	4.4	17
50	A gel aging effect in the synthesis of open-framework gallium phosphates: structure solution and solid-state NMR of a large-pore, open-framework material. <i>Dalton Transactions</i> , 2017, 46, 16895-16904.	3.3	4
51	The ambient hydration of the aluminophosphate JDF-2 to AlPO-53(A): insights from NMR crystallography. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2017, 73, 191-201.	0.5	6
52	Combining solid-state NMR spectroscopy with first-principles calculations – a guide to NMR crystallography. <i>Chemical Communications</i> , 2016, 52, 7186-7204.	4.1	202
53	Paramagnetic NMR of Phenolic Oxime Copper Complexes: A Joint Experimental and Density Functional Study. <i>Chemistry - A European Journal</i> , 2016, 22, 15328-15339.	3.3	22
54	Phase Composition and Disorder in La ₂ (Sn,Ti) ₂ O ₇ Ceramics: New Insights from NMR Crystallography. <i>Journal of Physical Chemistry C</i> , 2016, 120, 20288-20296.	3.1	15

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55	Investigating Unusual Homonuclear Intermolecular δ -Through-Space J Couplings in Organochalcogen Systems. <i>Inorganic Chemistry</i> , 2016, 55, 10881-10887.	4.0	15
56	Hunting for hydrogen: random structure searching and prediction of NMR parameters of hydrous wadsleyite. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 10173-10181.	2.8	19
57	NMR spectroscopy of minerals and allied materials. <i>Nuclear Magnetic Resonance</i> , 2016, , 1-52.	0.2	21
58	New insights into phase distribution, phase composition and disorder in $Y_{2-x}(Zr,Sn)_xO_7$ ceramics from NMR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 9049-9059.	2.8	22
59	Conformational Dependence of Through-Space Tellurium Spin-Spin Coupling in β -Substituted Bis(Tellurides). <i>Chemistry - A European Journal</i> , 2015, 21, 3613-3627.	3.3	19
60	β -Substituted Phosphorus-Tellurium Systems: An Experimental and Theoretical Investigation of the P-Te through-Space Interaction. <i>Inorganic Chemistry</i> , 2015, 54, 2435-2446.	4.0	30
61	Unusual Intermolecular δ -Through-Space J Couplings in Se Heterocycles. <i>Journal of the American Chemical Society</i> , 2015, 137, 6172-6175.	13.7	24
62	Exploiting Synthetic Conditions to Promote Structural Diversity within the Scandium(III)/Pyrimidine-4,6-dicarboxylate System. <i>Crystal Growth and Design</i> , 2015, 15, 2352-2363.	3.0	31
63	An NMR crystallographic approach to monitoring cation substitution in the aluminophosphate STA-2. <i>Solid State Nuclear Magnetic Resonance</i> , 2015, 65, 64-74.	2.3	14
64	Solid-state NMR measurements and DFT calculations of the magnetic shielding tensors of protons of water trapped in barium chlorate monohydrate. <i>RSC Advances</i> , 2014, 4, 56248-56258.	3.6	17
65	Mixed-Metal MIL-100(Sc,M) (M=Al, Cr, Fe) for Lewis Acid Catalysis and Tandem $C-C$ Bond Formation and Alcohol Oxidation. <i>Chemistry - A European Journal</i> , 2014, 20, 17185-17197.	3.3	104
66	Sterically Restricted Tin Phosphines, Stabilized by Weak Intramolecular Donor-Acceptor Interactions. <i>Organometallics</i> , 2014, 33, 2424-2433.	2.3	18
67	Probing interactions through space using spin-spin coupling. <i>Dalton Transactions</i> , 2014, 43, 6548-6560.	3.3	28
68	New Methods and Applications in Solid-State NMR Spectroscopy of Quadrupolar Nuclei. <i>Journal of the American Chemical Society</i> , 2014, 136, 15440-15456.	13.7	120
69	Calculating NMR parameters in aluminophosphates: evaluation of dispersion correction schemes. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 2660.	2.8	32
70	Recent developments in solid-state NMR spectroscopy of crystalline microporous materials. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 8223-8242.	2.8	69
71	Investigating Relationships between the Crystal Structure and ^{31}P Isotropic Chemical Shifts in Calcined Aluminophosphates. <i>Journal of Physical Chemistry C</i> , 2014, 118, 23285-23296.	3.1	23
72	Efficient Amplitude-Modulated Pulses for Triple- to Single-Quantum Coherence Conversion in MQMAS NMR. <i>Journal of Physical Chemistry A</i> , 2014, 118, 6018-6025.	2.5	19

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73	Zeolites with Continuously Tuneable Porosity. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 13210-13214.	13.8	104
74	Characterization of Structural Disorder in β -Ga ₂ O ₃ . <i>Journal of Physical Chemistry C</i> , 2014, 118, 16188-16198.	3.1	107
75	Multirate delivery of multiple therapeutic agents from metal-organic frameworks. <i>APL Materials</i> , 2014, 2, .	5.1	58
76	Solid-State NMR of High-Pressure Silicates in the Earth's Mantle. <i>Annual Reports on NMR Spectroscopy</i> , 2013, 79, 241-332.	1.5	11
77	Unusual Phase Behavior in the Piezoelectric Perovskite System, $\text{Li}_x\text{Na}_{1-x}\text{NbO}_3$. <i>Inorganic Chemistry</i> , 2013, 52, 8872-8880.	4.0	31
78	Investigation of the hydrothermal crystallisation of the perovskite solid solution $\text{NaCe}_{1-x}\text{La}_x\text{Ti}_2\text{O}_6$ and its defect chemistry. <i>Journal of Solid State Chemistry</i> , 2013, 207, 117-125.	2.9	8
79	Application of NMR crystallography to the determination of the mechanism of charge-balancing in organocation-templated $\text{AlPO}_4\text{STA-2}$. <i>CrystEngComm</i> , 2013, 15, 8668.	2.6	28
80	Structural Study of $\text{La}_x\text{Y}_{1-x}\text{ScO}_3$, Combining Neutron Diffraction, Solid-State NMR, and First-Principles DFT Calculations. <i>Journal of Physical Chemistry C</i> , 2013, 117, 2252-2265.	3.1	39
81	Exploiting Periodic First-Principles Calculations in NMR Spectroscopy of Disordered Solids. <i>Accounts of Chemical Research</i> , 2013, 46, 1964-1974.	15.6	53
82	High-resolution solid-state ¹³ C NMR spectroscopy of the paramagnetic metal-organic frameworks, STAM-1 and HKUST-1. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 919-929.	2.8	64
83	The pyrochlore to defect fluorite phase transition in $\text{Y}_2\text{Sn}_2\text{Zr}_2\text{O}_7$. <i>RSC Advances</i> , 2013, 3, 5090.	3.6	55
84	A family of zeolites with controlled pore size prepared using a top-down method. <i>Nature Chemistry</i> , 2013, 5, 628-633.	13.6	355
85	Color and Brightness Tuning in Heteronuclear Lanthanide Terephthalate Coordination Polymers. <i>European Journal of Inorganic Chemistry</i> , 2013, 2013, 3464-3476.	2.0	76
86	Water in the Earth's mantle: a solid-state NMR study of hydrous wadsleyite. <i>Chemical Science</i> , 2013, 4, 1523.	7.4	41
87	First-Principles Calculation of NMR Parameters Using the Gauge Including Projector Augmented Wave Method: A Chemist's Point of View. <i>Chemical Reviews</i> , 2012, 112, 5733-5779.	47.7	446
88	A novel structural form of MIL-53 observed for the scandium analogue and its response to temperature variation and CO ₂ adsorption. <i>Dalton Transactions</i> , 2012, 41, 3937-3941.	3.3	95
89	Applications of NMR Crystallography to Problems in Biomineralization: Refinement of the Crystal Structure and ³¹ P Solid-State NMR Spectral Assignment of Octacalcium Phosphate. <i>Journal of the American Chemical Society</i> , 2012, 134, 12508-12515.	13.7	80
90	A Multinuclear Solid-State NMR Study of Templated and Calcined Chabazite-Type GaPO-34. <i>Journal of Physical Chemistry C</i> , 2012, 116, 15048-15057.	3.1	24

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91	Isothermal ^{17}O enrichment of oxides using microlitre quantities of labelled water. <i>Chemical Science</i> , 2012, 3, 2293.	7.4	57
92	Exploiting the Chemical Shielding Anisotropy to Probe Structure and Disorder in Ceramics: ^{89}Y MAS NMR and First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2012, 116, 4273-4286.	3.1	41
93	Noncovalent Interactions in Peri-Substituted Chalconium Acenaphthene and Naphthalene Salts: A Combined Experimental, Crystallographic, Computational, and Solid-State NMR Study. <i>Inorganic Chemistry</i> , 2012, 51, 11087-11097.	4.0	38
94	New Twists on the Perovskite Theme: Crystal Structures of the Elusive Phases R and S of NaNbO_3 . <i>Inorganic Chemistry</i> , 2012, 51, 6876-6889.	4.0	78
95	Towards homonuclear J solid-state NMR correlation experiments for half-integer quadrupolar nuclei: experimental and simulated ^{11}B MAS spin-echo dephasing and calculated $2J_{\text{BB}}$ coupling constants for lithium diborate. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 5778.	2.8	34
96	^{77}Se Solid-State NMR of Inorganic and Organoselenium Systems: A Combined Experimental and Computational Study. <i>Journal of Physical Chemistry C</i> , 2011, 115, 10859-10872.	3.1	25
97	^{93}Nb NMR and DFT investigation of the polymorphs of NaNbO_3 . <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 7565.	2.8	50
98	^{119}Sn MAS NMR and first-principles calculations for the investigation of disorder in stannate pyrochlores. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 488-497.	2.8	49
99	Structural Chemistry, Monoclinic-to-Orthorhombic Phase Transition, and CO_2 Adsorption Behavior of the Small Pore Scandium Terephthalate, $\text{Sc}_2(\text{O}_2\text{CC}_6\text{H}_4\text{CO}_2)_3$, and Its Nitro- And Amino-Functionalized Derivatives. <i>Inorganic Chemistry</i> , 2011, 50, 10844-10858.	4.0	75
100	Protecting group and switchable pore-discriminating adsorption properties of a hydrophilic-hydrophobic metal-organic framework. <i>Nature Chemistry</i> , 2011, 3, 304-310.	13.6	141
101	Observation of "hidden" magnesium: First-principles calculations and ^{25}Mg solid-state NMR of enstatite. <i>Solid State Nuclear Magnetic Resonance</i> , 2011, 40, 91-99.	2.3	25
102	A co-templating route to the synthesis of Cu SAPO STA-7, giving an active catalyst for the selective catalytic reduction of NO. <i>Microporous and Mesoporous Materials</i> , 2011, 146, 36-47.	4.4	44
103	Octaselenocyclododecane. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 4123-4126.	13.8	23
104	Synthesis, characterisation and adsorption properties of microporous scandium carboxylates with rigid and flexible frameworks. <i>Microporous and Mesoporous Materials</i> , 2011, 142, 322-333.	4.4	170
105	Detecting solid-state reactivity in 10-hydroxy-10,9-boroxophenanthrene using NMR spectroscopy. <i>Tetrahedron</i> , 2010, 66, 6238-6250.	1.9	21
106	Molecular Modeling, Multinuclear NMR, and Diffraction Studies in the Templated Synthesis and Characterization of the Aluminophosphate Molecular Sieve STA-2. <i>Journal of Physical Chemistry C</i> , 2010, 114, 12698-12710.	3.1	44
107	The Polar Phase of NaNbO_3 : A Combined Study by Powder Diffraction, Solid-State NMR, and First-Principles Calculations. <i>Journal of the American Chemical Society</i> , 2010, 132, 8732-8746.	13.7	178
108	High-Resolution ^{19}F MAS NMR Spectroscopy: Structural Disorder and Unusual C-F Couplings in a Fluorinated Hydroxy-Silicate. <i>Journal of the American Chemical Society</i> , 2010, 132, 15651-15660.	13.7	83

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109	Task specific ionic liquids for the ionothermal synthesis of siliceous zeolites. <i>Chemical Science</i> , 2010, 1, 483.	7.4	81
110	Dynamics on the microsecond timescale in hydrous silicates studied by solid-state ^2H NMR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 2989.	2.8	30
111	Novel Large-Pore Aluminophosphate Molecular Sieve STA-15 Prepared Using the Tetrapropylammonium Cation As a Structure Directing Agent. <i>Chemistry of Materials</i> , 2010, 22, 338-346.	6.7	35
112	Disordered lithium niobate rock-salt materials prepared by hydrothermal synthesis. <i>Dalton Transactions</i> , 2010, 39, 6031.	3.3	14
113	Structure and NMR assignment in $\text{AlPO}_4\text{-15}$: A combined study by diffraction, MAS NMR and first-principles calculations. <i>Solid State Sciences</i> , 2009, 11, 1001-1006.	3.2	38
114	Second-order cross-term interactions in high-resolution MAS NMR of quadrupolar nuclei. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2009, 55, 160-181.	7.5	28
115	Solid-State ^{17}O NMR Spectroscopy of Hydrous Magnesium Silicates: Evidence for Proton Dynamics. <i>Journal of Physical Chemistry C</i> , 2009, 113, 465-471.	3.1	61
116	Spin-locking of half-integer quadrupolar nuclei in nuclear magnetic resonance of solids: Second-order quadrupolar and resonance offset effects. <i>Journal of Chemical Physics</i> , 2009, 131, 194509.	3.0	48
117	Transformation of $\text{AlPO}_4\text{-53}$ to JDF-2 : Reversible Dehydration of a Templated Aluminophosphate Studied by MAS NMR and Diffraction. <i>Journal of Physical Chemistry C</i> , 2009, 113, 10780-10789.	3.1	40
118	Recent advances in solid-state NMR spectroscopy of quadrupolar nuclei. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 6892.	2.8	114
119	Cation Disorder in Pyrochlore Ceramics: ^{89}Y MAS NMR and First-Principles Calculations. <i>Journal of Physical Chemistry C</i> , 2009, 113, 18874-18883.	3.1	62
120	Control of polymorphism in NaNbO_3 by hydrothermal synthesis. <i>Chemical Communications</i> , 2009, , 68-70.	4.1	65
121	DFT calculations of quadrupolar solid-state NMR properties: Some examples in solid-state inorganic chemistry. <i>Journal of Computational Chemistry</i> , 2008, 29, 2279-2287.	3.3	52
122	Structure and NMR assignment in calcined and as-synthesized forms of $\text{AlPO}_4\text{-14}$: a combined study by first-principles calculations and high-resolution ^{27}Al - ^{31}P MAS NMR correlation. <i>Physical Chemistry Chemical Physics</i> , 2008, 10, 5754.	2.8	95
123	First-principles calculations of solid-state ^{17}O and ^{29}Si NMR spectra of Mg_2SiO_4 polymorphs. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 1587-1598.	2.8	65
124	^{17}O and ^{29}Si NMR Parameters of MgSiO_3 Phases from High-Resolution Solid-State NMR Spectroscopy and First-Principles Calculations. <i>Journal of the American Chemical Society</i> , 2007, 129, 13213-13224.	13.7	104
125	The effect of caesium on barium hollandites studied by neutron diffraction and magic-angle spinning (MAS) nuclear magnetic resonance. <i>Journal of Materials Science</i> , 2007, 42, 9379-9391.	3.7	6
126	^{89}Y Magic-Angle Spinning NMR of $\text{Y}_2\text{Ti}_2\text{-xSn}_x\text{O}_7$ Pyrochlores. <i>Journal of Physical Chemistry B</i> , 2006, 110, 10358-10364.	2.6	47

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127	Solid state ^{17}O NMR – an introduction to the background principles and applications to inorganic materials. <i>Chemical Society Reviews</i> , 2006, 35, 718-735.	38.1	203
128	Nuclear Overhauser Effect (NOE) Enhancement of ^{11}B NMR Spectra of Borane Adducts in the Solid State. <i>Journal of the American Chemical Society</i> , 2006, 128, 6782-6783.	13.7	17
129	Characterisation of the $(\text{Y}_{1-x}\text{La}_x)_2\text{Ti}_2\text{O}_7$ system by powder diffraction and nuclear magnetic resonance methods. <i>Journal of Materials Chemistry</i> , 2006, 16, 4665-4674.	6.7	26
130	Dynamics on the Microsecond Timescale in Microporous Aluminophosphate AlPO-14 as Evidenced by ^{27}Al MQMAS and STMAS NMR Spectroscopy. <i>Journal of the American Chemical Society</i> , 2006, 128, 8054-8062.	13.7	72
131	^2H double-quantum MAS NMR spectroscopy as a probe of dynamics on the microsecond timescale in solids. <i>Chemical Physics Letters</i> , 2006, 423, 276-281.	2.6	58
132	STARTMAS: A MAS-based method for acquiring isotropic NMR spectra of spin $I=3/2$ nuclei in real time. <i>Chemical Physics Letters</i> , 2006, 431, 390-396.	2.6	13
133	^{23}Na multiple-quantum MAS NMR of the perovskites NaNbO_3 and NaTaO_3 . <i>Physical Chemistry Chemical Physics</i> , 2006, 8, 3423-3431.	2.8	86
134	Structural information from quadrupolar nuclei in solid state NMR. <i>Concepts in Magnetic Resonance Part A: Bridging Education and Research</i> , 2006, 28A, 183-248.	0.5	136
135	Correlating fast and slow chemical shift spinning sideband patterns in solid-state NMR. <i>Journal of Magnetic Resonance</i> , 2005, 174, 301-309.	2.1	32
136	Disorder and Dynamics in Pollucite from ^{133}Cs and ^{27}Al NMR. <i>Journal of the American Ceramic Society</i> , 2005, 88, 1575-1583.	3.8	20
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