

Sharon Ashbrook

List of Publications by Year in descending order

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178
papers

8,238
citations

41344

49
h-index

62596

80
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186
all docs

186
docs citations

186
times ranked

7030
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	Recent advances in solid-state nuclear magnetic resonance spectroscopy of quadrupolar nuclei. <i>Magnetic Resonance in Chemistry</i> , 2021, 59, 851-852.	1.9	1
7	Single-step synthesis and interface tuning of core-shell metal-organic framework nanoparticles. <i>Chemical Science</i> , 2021, 12, 4494-4502.	7.4	11
8	Formation Mechanism and Porosity Development in Porous Boron Nitride. <i>Journal of Physical Chemistry C</i> , 2021, 125, 27429-27439.	3.1	15
9	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
10	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
11	Solid-state host-guest influences on a BODIPY dye hosted within a crystalline sponge. <i>New Journal of Chemistry</i> , 2020, 44, 14108-14115.	2.8	6
12	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
13	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
14	Synthesis of Chiral MOF-74 Frameworks by Post-synthetic Modification by Using an Amino Acid. <i>Chemistry - A European Journal</i> , 2020, 26, 13957-13965.	3.3	35
15	Mechanochemically assisted hydrolysis in the ADOR process. <i>Chemical Science</i> , 2020, 11, 7060-7069.	7.4	12
16	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
17	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
18	Synthesis and Polymorphism of Mixed Aluminum-Gallium Oxides. <i>Inorganic Chemistry</i> , 2020, 59, 3805-3816.	4.0	28

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19	Phosphorus- and Bismuth-Substituted Acenaphthenes: A Synthetic, Structural, and Computational Study. <i>Inorganic Chemistry</i> , 2020, 59, 5616-5625.	4.0	13
20	Fast room temperature lability of aluminosilicate zeolites. <i>Nature Communications</i> , 2019, 10, 4690.	12.8	75
21	Ensemble-Based Modeling of the NMR Spectra of Solid Solutions: Cation Disorder in $Y_2(Sn,Ti)_7O_{17}$. <i>Journal of the American Chemical Society</i> , 2019, 141, 17838-17846.	13.7	29
22	NMR spectroscopy of paramagnetic solids. <i>Solid State Nuclear Magnetic Resonance</i> , 2019, 104, 101625.	2.3	1
23	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
24	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
25	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
26	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
27	^{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
28	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
29	Sensitivity improvement in 5QMAS NMR experiments using FAM-N pulses. <i>Solid State Nuclear Magnetic Resonance</i> , 2019, 100, 1-10.	2.3	3
30	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
31	Rationalization of solid-state NMR multi-pulse decoupling strategies: Coupling of spin $l = \frac{1}{2}$ and half-integer quadrupolar nuclei. <i>Journal of Magnetic Resonance</i> , 2019, 303, 48-56.	2.1	3
32	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
33	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
34	^{17}O solid-state NMR spectroscopy of $A_2B_7O_{17}$ oxides: quantitative isotopic enrichment and spectral acquisition?. <i>RSC Advances</i> , 2018, 8, 7089-7101.	3.6	13
35	Modulator-Controlled Synthesis of Microporous STA-26, an Interpenetrated 8,3-Connected Zirconium MOF with the <i>the</i> Topology, and its Reversible Lattice Shift. <i>Chemistry - A European Journal</i> , 2018, 24, 6115-6126.	3.3	23
36	Recent Advances in Solid-State Nuclear Magnetic Resonance Spectroscopy. <i>Annual Review of Analytical Chemistry</i> , 2018, 11, 485-508.	5.4	45

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37	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
38	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
39	A Bifunctional MOF Catalyst Containing Metal-Phosphine and Lewis Acidic Active Sites. <i>Chemistry - A European Journal</i> , 2018, 24, 15309-15318.	3.3	40
40	SERS of Trititanate Nanotubes: Selective Enhancement of Catechol Compounds. <i>ChemistrySelect</i> , 2018, 3, 8338-8343.	1.5	3
41	Perspective: Current advances in solid-state NMR spectroscopy. <i>Journal of Chemical Physics</i> , 2018, 149, 040901.	3.0	28
42	Polymorphism, Weak Interactions and Phase Transitions in Chalcogen-Phosphorus Heterocycles. <i>Chemistry - A European Journal</i> , 2018, 24, 11067-11081.	3.3	4
43	An expanded MIL-53-type coordination polymer with a reactive pendant ligand. <i>CrystEngComm</i> , 2018, 20, 4355-4358.	2.6	5
44	Hydrolytic stability in hemilabile metal-organic frameworks. <i>Nature Chemistry</i> , 2018, 10, 1096-1102.	13.6	134
45	Alkaline-Earth Rhodium Hydroxides: Synthesis, Structures, and Thermal Decomposition to Complex Oxides. <i>Inorganic Chemistry</i> , 2018, 57, 11217-11224.	4.0	8
46	Synthesis of ZIF-93/11 Hybrid Nanoparticles via Post-Synthetic Modification of ZIF-93 and Their Use for H ₂ /CO ₂ Separation. <i>Chemistry - A European Journal</i> , 2018, 24, 11211-11219.	3.3	27
47	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
48	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
49	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
50	Exploiting NMR spectroscopy for the study of disorder in solids. <i>International Reviews in Physical Chemistry</i> , 2017, 36, 39-115.	2.3	65
51	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 E-Te(II,IV) Bonding Situations (E = O, S, Se). <i>Organometallics</i> , 2017, 36, 1566-1579.	2.3	18
52	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
53	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
54	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

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55	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
56	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
57	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
58	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
59	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
60	Ionothermal synthesis and characterization of CoAPO-34 molecular sieve. <i>Microporous and Mesoporous Materials</i> , 2017, 239, 336-341.	4.4	17
61	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
62	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
63	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
64	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
65	Exploring the self-assembly and energy transfer of dynamic supramolecular iridium-porphyrin systems. <i>Dalton Transactions</i> , 2016, 45, 17195-17205.	3.3	23
66	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
67	Investigating Unusual Homonuclear Intermolecular Through-Space J Couplings in Organochalcogen Systems. <i>Inorganic Chemistry</i> , 2016, 55, 10881-10887.	4.0	15
68	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
69	Inside Cover: Conformational Dependence of Through-Space Tellurium–Tellurium Spin–Spin Coupling in <i>Peri</i> -Substituted Bis(Tellurides) (<i>Chem. Eur. J.</i> 9/2015). <i>Chemistry - A European Journal</i> , 2015, 21, 3506-3506.	3.3	0
70	[1,2,5]Selenadiazolo[3,4- <i>b</i>]pyrazines: Synthesis from 3,4-Diamino-1,2,5-selenadiazole and Generation of Persistent Radical Anions. <i>European Journal of Organic Chemistry</i> , 2015, 2015, 5585-5593.	2.4	18
71	New insights into phase distribution, phase composition and disorder in Y ₂ (Zr,Sn) ₂ O ₇ ceramics from NMR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 9049-9059.	2.8	22
72	Conformational Dependence of Through-Space Tellurium–Tellurium Spin–Spin Coupling in <i>Peri</i> -Substituted Bis(Tellurides). <i>Chemistry - A European Journal</i> , 2015, 21, 3613-3627.	3.3	19

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73	<i>i>Peri</i>-Substituted Phosphorusâ€“Tellurium Systemsâ€“An Experimental and Theoretical Investigation of the P-Te through-Space Interaction. Inorganic Chemistry, 2015, 54, 2435-2446.</i>	4.0	30
74	Unusual Intermolecular â€œThrough-Spaceâ€•</i> Couplings in Pâ€“Se Heterocycles. Journal of the American Chemical Society, 2015, 137, 6172-6175.	13.7	24
75	Exploiting Synthetic Conditions to Promote Structural Diversity within the Scandium(III)/Pyrimidine-4,6-dicarboxylate System. Crystal Growth and Design, 2015, 15, 2352-2363.	3.0	31
76	Post-synthetic modification of zinc metal-organic frameworks through palladium-catalysed carbonâ€“carbon bond formation. Journal of Organometallic Chemistry, 2015, 792, 134-138.	1.8	4
77	An NMR crystallographic approach to monitoring cation substitution in the aluminophosphate STA-2. Solid State Nuclear Magnetic Resonance, 2015, 65, 64-74.	2.3	14
78	Solid-state NMR measurements and DFT calculations of the magnetic shielding tensors of protons of water trapped in barium chlorate monohydrate. RSC Advances, 2014, 4, 56248-56258.	3.6	17
79	Mixedâ€“Metal MILâ€“100(Sc,M) (M=Al, Cr, Fe) for Lewis Acid Catalysis and Tandem C=C Bond Formation and Alcohol Oxidation. Chemistry - A European Journal, 2014, 20, 17185-17197.	3.3	104
80	Sterically Restricted Tin Phosphines, Stabilized by Weak Intramolecular Donorâ€“Acceptor Interactions. Organometallics, 2014, 33, 2424-2433.	2.3	18
81	Probing interactions through space using spinâ€“spin coupling. Dalton Transactions, 2014, 43, 6548-6560.	3.3	28
82	A Modular Approach for the Synthesis of Nanometer-Sized Polynitroxide Multi-Spin Systems. Journal of Organic Chemistry, 2014, 79, 8313-8323.	3.2	13
83	New Methods and Applications in Solid-State NMR Spectroscopy of Quadrupolar Nuclei. Journal of the American Chemical Society, 2014, 136, 15440-15456.	13.7	120
84	Calculating NMR parameters in aluminophosphates: evaluation of dispersion correction schemes. Physical Chemistry Chemical Physics, 2014, 16, 2660.	2.8	32
85	Recent developments in solid-state NMR spectroscopy of crystalline microporous materials. Physical Chemistry Chemical Physics, 2014, 16, 8223-8242.	2.8	69
86	Investigating Relationships between the Crystal Structure and ³¹ P Isotropic Chemical Shifts in Calcined Aluminophosphates. Journal of Physical Chemistry C, 2014, 118, 23285-23296.	3.1	23
87	Efficient Amplitude-Modulated Pulses for Triple- to Single-Quantum Coherence Conversion in MQMAS NMR. Journal of Physical Chemistry A, 2014, 118, 6018-6025.	2.5	19
88	Zeolites with Continuously Tuneable Porosity. Angewandte Chemie - International Edition, 2014, 53, 13210-13214.	13.8	104
89	Characterization of Structural Disorder in ¹³ Ga ₂ O ₃ . Journal of Physical Chemistry C, 2014, 118, 16188-16198.	3.1	107
90	Multirate delivery of multiple therapeutic agents from metal-organic frameworks. APL Materials, 2014, 2, .	5.1	58

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91	Solid-State NMR of High-Pressure Silicates in the Earth's Mantle. Annual Reports on NMR Spectroscopy, 2013, 79, 241-332.	1.5	11
92	Unusual Phase Behavior in the Piezoelectric Perovskite System, $\text{Li}_x\text{Na}_{1-x}\text{NbO}_3$. Inorganic Chemistry, 2013, 52, 8872-8880.	4.0	31
93	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. Journal of Solid State Chemistry, 2013, 207, 117-125.	2.9	8
94	Application of NMR crystallography to the determination of the mechanism of charge-balancing in organocation-templated AlPO_4 STA-2. CrystEngComm, 2013, 15, 8668.	2.6	28
95	Structural Study of $\text{La}_x\text{Y}_{1-x}\text{ScO}_3$, Combining Neutron Diffraction, Solid-State NMR, and First-Principles DFT Calculations. Journal of Physical Chemistry C, 2013, 117, 2252-2265.	3.1	39
96	Exploiting Periodic First-Principles Calculations in NMR Spectroscopy of Disordered Solids. Accounts of Chemical Research, 2013, 46, 1964-1974.	15.6	53
97	High-resolution solid-state ^{13}C NMR spectroscopy of the paramagnetic metal-organic frameworks, STAM-1 and HKUST-1. Physical Chemistry Chemical Physics, 2013, 15, 919-929.	2.8	64
98	The pyrochlore to defect fluorite phase transition in $\text{Y}_2\text{Sn}_2-x\text{Zr}_x\text{O}_7$. RSC Advances, 2013, 3, 5090.	3.6	55
99	A family of zeolites with controlled pore size prepared using a top-down method. Nature Chemistry, 2013, 5, 628-633.	13.6	355
100	Color and Brightness Tuning in Heteronuclear Lanthanide Terephthalate Coordination Polymers. European Journal of Inorganic Chemistry, 2013, 2013, 3464-3476.	2.0	76
101	Water in the Earth's mantle: a solid-state NMR study of hydrous wadsleyite. Chemical Science, 2013, 4, 1523.	7.4	41
102	First-Principles Calculation of NMR Parameters Using the Gauge Including Projector Augmented Wave Method: A Chemist's Point of View. Chemical Reviews, 2012, 112, 5733-5779.	47.7	446
103	A novel structural form of MIL-53 observed for the scandium analogue and its response to temperature variation and CO_2 adsorption. Dalton Transactions, 2012, 41, 3937-3941.	3.3	95
104	Applications of NMR Crystallography to Problems in Biomineralization: Refinement of the Crystal Structure and ^{31}P Solid-State NMR Spectral Assignment of Octacalcium Phosphate. Journal of the American Chemical Society, 2012, 134, 12508-12515.	13.7	80
105	A Multinuclear Solid-State NMR Study of Templated and Calcined Chabazite-Type GaPO -34. Journal of Physical Chemistry C, 2012, 116, 15048-15057.	3.1	24
106	Ionothermal ^{17}O enrichment of oxides using microlitre quantities of labelled water. Chemical Science, 2012, 3, 2293.	7.4	57
107	Exploiting the Chemical Shielding Anisotropy to Probe Structure and Disorder in Ceramics: ^{89}Y MAS NMR and First-Principles Calculations. Journal of Physical Chemistry C, 2012, 116, 4273-4286.	3.1	41
108	Noncovalent Interactions in Peri-Substituted Chalconium Acenaphthene and Naphthalene Salts: A Combined Experimental, Crystallographic, Computational, and Solid-State NMR Study. Inorganic Chemistry, 2012, 51, 11087-11097.	4.0	38

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109	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
110	Synthesis and crystal chemistry of the STA-12 family of metal N,N'-piperazinebis(methylenephosphonate)s and applications of STA-12(Ni) in the separation of gases. <i>Microporous and Mesoporous Materials</i> , 2012, 157, 3-17.	4.4	49
111	Towards homonuclear J solid-state NMR correlation experiments for half-integer quadrupolar nuclei: experimental and simulated ^{11}B MAS spin-echo dephasing and calculated 2JBB coupling constants for lithium diborate. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 5778.	2.8	34
112	^{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
113	Synthesis and characterization of hybrid organic/inorganic nanotubes of the imogolite type and their behaviour towards methane adsorption. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 744-750.	2.8	102
114	^{93}Nb NMR and DFT investigation of the polymorphs of NaNbO_3 . <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 7565.	2.8	50
115	^{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
116	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
117	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
118	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
119	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
120	Octaselenocyclododecane. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 4123-4126.	13.8	23
121	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
122	Detecting solid-state reactivity in 10-hydroxy-10,9-boroxophenanthrene using NMR spectroscopy. <i>Tetrahedron</i> , 2010, 66, 6238-6250.	1.9	21
123	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
124	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
125	High-Resolution ^{19}F MAS NMR Spectroscopy: Structural Disorder and Unusual J Couplings in a Fluorinated Hydroxy-Silicate. <i>Journal of the American Chemical Society</i> , 2010, 132, 15651-15660.	13.7	83
126	Task specific ionic liquids for the ionothermal synthesis of siliceous zeolites. <i>Chemical Science</i> , 2010, 1, 483.	7.4	81

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127	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
128	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
129	Structure and NMR assignment in AlPO ₄ -15: A combined study by diffraction, MAS NMR and first-principles calculations. <i>Solid State Sciences</i> , 2009, 11, 1001-1006.	3.2	38
130	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
131	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
132	Multinuclear Magnetic Resonance and DFT Studies of the Poly(chlorotrifluoroethylene- <i>i>alt</i>-ethyl vinyl ether) Copolymers. <i>Macromolecules</i>, 2009, 42, 5652-5659.</i>	4.8	42
133	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
134	Transformation of AlPO ₅ 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
135	Early Stage Reversed Crystal Growth of Zeolite A and Its Phase Transformation to Sodalite. <i>Journal of the American Chemical Society</i> , 2009, 131, 17986-17992.	13.7	129
136	Recent advances in solid-state NMR spectroscopy of quadrupolar nuclei. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 6892.	2.8	114
137	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
138	Control of polymorphism in NaNbO ₃ by hydrothermal synthesis. <i>Chemical Communications</i> , 2009, , 68-70.	4.1	65
139	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
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