

# Alexander G Stepanov

## List of Publications by Year in descending order

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185  
papers

5,804  
citations

71102

41  
h-index

110387

64  
g-index

190  
all docs

190  
docs citations

190  
times ranked

4597  
citing authors

#	ARTICLE	IF	CITATIONS
1	Structure, hydrogen bond dynamics and phase transition in a model ionic liquid electrolyte. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 6064-6071.	2.8	8
2	Property–activity relations of multifunctional reactive ensembles in cation-exchanged zeolites: a case study of methane activation on Zn <sup>2+</sup> -modified zeolite BEA. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 6492-6504.	2.8	5
3	High-Temperature Quantum Tunneling and Hydrogen Bonding Rearrangements Characterize the Solid–Solid Phase Transitions in a Phosphonium-Based Protic Ionic Liquid. <i>Chemistry - A European Journal</i> , 2022, , .	3.3	1
4	Selective Dimerization of Ethene to 2-Butene on Zn <sup>2+</sup> -Modified ZSM-5 Zeolite. <i>Journal of Physical Chemistry C</i> , 2022, 126, 6570-6577.	3.1	8
5	Butane isomers mobility and framework dynamics in UiO-66 (Zr) MOF: Impact of the hydroxyl groups in zirconia cluster. <i>Solid State Nuclear Magnetic Resonance</i> , 2022, 118, 101784.	2.3	4
6	METAL-ORGANIC FRAMEWORKS IN RUSSIA: FROM THE SYNTHESIS AND STRUCTURE TO FUNCTIONAL PROPERTIES AND MATERIALS. <i>Journal of Structural Chemistry</i> , 2022, 63, 671-843.	1.0	35
7	Dynamics in nanoporous materials probed by <sup>2</sup> H solid state NMR: estimation of self-diffusion coefficients. <i>Adsorption</i> , 2021, 27, 841-855.	3.0	5
8	Effect of Copper State in Cu/H-ZSM-5 on Methane Activation by Brønsted Acid Sites, Studied by <sup>1</sup> H MAS NMR In Situ Monitoring the H/D Hydrogen Exchange of the Alkane with Brønsted Acid Sites. <i>Journal of Physical Chemistry C</i> , 2021, 125, 2182-2193.	3.1	16
9	Dissecting the effects of water guest adsorption and framework breathing on the AlO <sub>4</sub> (OH) <sub>2</sub> centres of metal–organic framework MIL-53 (Al) by solid state NMR and structural analysis. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 18925-18929.	2.8	9
10	Selective Gas Uptake and Rotational Dynamics in a (3,24)-Connected Metal–Organic Framework Material. <i>Journal of the American Chemical Society</i> , 2021, 143, 3348-3358.	13.7	39
11	Unraveling the Mechanism of Methane Activation on Zn-Modified Zeolites by Solid-State NMR. <i>Chemistry Methods</i> , 2021, 1, 224-230.	3.8	6
12	UiO-66 (Zr) MOF as a Promising Material for Butane Isomers Separation: Evidence Based on the Analysis of the Adsorbed Alkanes Mobility by <sup>2</sup> H NMR and Molecular Dynamics Simulation. <i>Journal of Physical Chemistry C</i> , 2021, 125, 13391-13400.	3.1	15
13	Molecular Insight into the Slow Dynamics of C <sub>4</sub> Hydrocarbons in the Zeolitic–Imidazole Framework (ZIF-8). <i>ACS Applied Materials &amp; Interfaces</i> , 2021, 13, 33685-33692.	8.0	7
14	Isobutene Transformation to Aromatics on Zn-Modified Zeolite: Particular Effects of Zn <sup>2+</sup> and ZnO Species on the Reaction Occurrence Revealed with Solid-State NMR and FTIR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2021, 125, 15343-15353.	3.1	17
15	Isobutane Transformation to Aromatics on Zn-Modified Zeolites: Intermediates and the Effect of Zn <sup>2+</sup> and ZnO Species on the Reaction Occurrence Revealed by <sup>13</sup> C MAS NMR. <i>ChemPhysChem</i> , 2021, , .	2.1	5
16	Does the Zn <sup>2+</sup> Species Introduced into H-ZSM-5 Zeolite Affect the Strength of Brønsted Acid Sites?. <i>ChemCatChem</i> , 2020, 12, 478-487.	3.7	12
17	The accuracy challenge of the DFT-based molecular assignment of <sup>13</sup> C MAS NMR characterization of surface intermediates in zeolite catalysis. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 24004-24013.	2.8	11
18	Freezing the Motion in Hydroxy-Functionalized Ionic Liquids—Temperature Dependent NMR Deuteron Quadrupole Coupling Constants for Two Types of Hydrogen Bonds Far below the Glass Transition. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 6000-6006.	4.6	10

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19	Molecular Mobility of Tert-butyl Alcohol Confined in a Breathing MIL-53 (Al) Metal-Organic Framework. <i>ChemPhysChem</i> , 2020, 21, 1951-1956.	2.1	3
20	Which Species, Zn <sup>2+</sup> Cations or ZnO Clusters, Are More Efficient for Olefin Aromatization? <sup>13</sup> C Solid-State NMR Investigation of <i>n</i> -But-1-ene Transformation on Zn-Modified Zeolite. <i>ACS Catalysis</i> , 2020, 10, 14224-14233.	11.2	29
21	Dynamics of isobutane is a sensitive probe for framework breathing in MIL-53 (Al) MOF. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 18695-18702.	2.8	8
22	<i>n</i> -Butane transformation on Zn/H-BEA. The effect of different Zn species (Zn <sup>2+</sup> and ZnO) on the reaction performance. <i>Journal of Catalysis</i> , 2020, 391, 69-79.	6.2	12
23	Mechanism of H/D Hydrogen Exchange of <i>n</i> -Butane with Brønsted Acid Sites on Zn-Modified Zeolite: The Effect of Different Zn Species (Zn <sup>2+</sup> and ZnO) on the Activation of Alkane C-H Bonds. <i>Journal of Physical Chemistry C</i> , 2020, 124, 20270-20279.	3.1	15
24	Heterogeneous epoxidation of menadione with hydrogen peroxide over the zeolite imidazolate framework ZIF-8. <i>Dalton Transactions</i> , 2020, 49, 12546-12549.	3.3	7
25	Counting cations involved in cationic clusters of hydroxy-functionalized ionic liquids by means of infrared and solid-state NMR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 6861-6867.	2.8	17
26	Superprotonic Conductivity in Metal-Organic Framework via Solvent-Free Coordinative Urea Insertion. <i>Journal of the American Chemical Society</i> , 2020, 142, 6861-6865.	13.7	65
27	Transformation of a proton insulator to a conductor <i>via</i> reversible amorphous to crystalline structure transformation of MOFs. <i>Chemical Communications</i> , 2020, 56, 4468-4471.	4.1	11
28	Dynamics of propene and propane in ZIF-8 probed by solid-state <sup>2</sup> H NMR. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 5976-5984.	2.8	15
29	Nature of the Surface Intermediates Formed from Methane on Cu-ZSM-5 Zeolite: A Combined Solid-State Nuclear Magnetic Resonance and Density Functional Theory Study. <i>Journal of Physical Chemistry C</i> , 2020, 124, 6242-6252.	3.1	38
30	Methane Activation on H-ZSM-5 Zeolite with Low Copper Loading. The Nature of Active Sites and Intermediates Identified with the Combination of Spectroscopic Methods. <i>Inorganic Chemistry</i> , 2020, 59, 2037-2050.	4.0	25
31	Dynamics of xylene isomers in MIL-53 (Al) MOF probed by solid state <sup>2</sup> H NMR. <i>Microporous and Mesoporous Materials</i> , 2020, 300, 110155.	4.4	15
32	Hydrogen Bonding Between Ions of Like Charge in Ionic Liquids Characterized by NMR Deuteron Quadrupole Coupling Constants—Comparison with Salt Bridges and Molecular Systems. <i>Angewandte Chemie</i> , 2019, 131, 18027-18035.	2.0	7
33	Hydrogen Bonding Between Ions of Like Charge in Ionic Liquids Characterized by NMR Deuteron Quadrupole Coupling Constants—Comparison with Salt Bridges and Molecular Systems. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 17863-17871.	13.8	41
34	Propylene Transformation on Zn-Modified Zeolite: Is There Any Difference in the Effect of Zn <sup>2+</sup> Cations or ZnO Species on the Reaction Occurrence?. <i>Journal of Physical Chemistry C</i> , 2019, 123, 27573-27583.	3.1	23
35	Propane activation on Zn-modified zeolite. The effect of the nature of Zn-species on the mechanism of H/D hydrogen exchange of the alkane with Brønsted acid sites. <i>Journal of Catalysis</i> , 2019, 378, 341-352.	6.2	23
36	Post-synthetic modulation of the charge distribution in a metal-organic framework for optimal binding of carbon dioxide and sulfur dioxide. <i>Chemical Science</i> , 2019, 10, 1472-1482.	7.4	62

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37	Diffusion in Nanoporous Materials: from Paradigm Shift by Zhdanov Zeolites Till Recent Insight. <i>Petroleum Chemistry</i> , 2019, 59, 275-296.	1.4	5
38	<sup>2</sup> H Solid-State NMR Spectroscopy Reveals the Dynamics of a Pyridine Probe Interacting with Coordinatively Unsaturated Metal Sites of MIL-100(Al) Metal-Organic Frameworks. <i>Chemistry - A European Journal</i> , 2019, 25, 10808-10812.	3.3	22
39	Mobility of Aromatic Guests and Isobutane in ZIF-8 Metal-Organic Framework Studied by <sup>2</sup> H Solid State NMR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2019, 123, 13765-13774.	3.1	23
40	The effect of amorphization on the molecular motion of the 2-methylimidazolate linkers in ZIF-8. <i>Chemical Communications</i> , 2019, 55, 5906-5909.	4.1	14
41	Guests Like Gear Levers: Donor Binding to Coordinatively Unsaturated Metal Sites in MIL-101 Controls the Linker's Rotation. <i>Chemistry - A European Journal</i> , 2019, 25, 5163-5168.	3.3	8
42	Simultaneous determination of deuteron quadrupole coupling constants and rotational correlation times: the model case of hydrogen bonded ionic liquids. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 25597-25605.	2.8	8
43	Propane Transformation on Zn-Modified Zeolite. Effect of the Nature of Zn Species on Alkane Aromatization and Hydrogenolysis. <i>Journal of Physical Chemistry C</i> , 2019, 123, 30473-30485.	3.1	29
44	NMR Study of the Host Structure and Guest Dynamics Investigated with Alkane/Alkene Mixtures in Metal Organic Frameworks ZIF-8. <i>Journal of Physical Chemistry C</i> , 2019, 123, 1904-1912.	3.1	22
45	Dynamical heterogeneities in ionic liquids as revealed from deuteron NMR. <i>Chemical Communications</i> , 2018, 54, 3098-3101.	4.1	21
46	Pulse EPR Study of Gas Adsorption in Cu <sup>2+</sup> -Doped Metal-Organic Framework [Zn <sub>2</sub> (1,4-bdc) <sub>2</sub> (dabco)]. <i>Applied Magnetic Resonance</i> , 2018, 49, 255-264.	1.2	20
47	Alkane/alkene mixture diffusion in silicalite-1 studied by MAS PFG NMR. <i>Microporous and Mesoporous Materials</i> , 2018, 257, 128-134.	4.4	23
48	Which Activation Energy Do We Measure? Analysis of the Kinetics of Propene-3- <sup>13</sup> C Double-Bond-Shift Reaction on Silicalite-1 by <sup>1</sup> H MAS NMR In Situ. <i>Journal of Physical Chemistry C</i> , 2018, 122, 23432-23440.	3.1	9
49	Enhancement of Proton Conductivity in Nonporous Metal-Organic Frameworks: The Role of Framework Proton Density and Humidity. <i>Chemistry of Materials</i> , 2018, 30, 7593-7602.	6.7	55
50	Direct Measurement of Zeolite Brønsted Acidity by FTIR Spectroscopy: Solid-State <sup>1</sup> H MAS NMR Approach for Reliable Determination of the Integrated Molar Absorption Coefficients. <i>Journal of Physical Chemistry C</i> , 2018, 122, 25386-25395.	3.1	69
51	Characterization of Fast Restricted Librations of Terephthalate Linkers in MOF UiO-66(Zr) by <sup>2</sup> H NMR Spin-Lattice Relaxation Analysis. <i>Journal of Physical Chemistry C</i> , 2018, 122, 12956-12962.	3.1	19
52	Uncovering the Rotation and Translational Mobility of Benzene Confined in UiO-66 (Zr) Metal-Organic Framework by the <sup>2</sup> H NMR-QENS Experimental Toolbox. <i>Journal of Physical Chemistry C</i> , 2017, 121, 2844-2857.	3.1	35
53	Different Efficiency of Zn <sup>2+</sup> and ZnO Species for Methane Activation on Zn-Modified Zeolite. <i>ACS Catalysis</i> , 2017, 7, 1818-1830.	11.2	151
54	Tailoring porosity and rotational dynamics in a series of octacarboxylate metal-organic frameworks. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, 3056-3061.	7.1	73

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55	Probing the Guest-Mediated Structural Mobility in the UiO-66(Zr) Framework by $^2\text{H}$ NMR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2017, 121, 11593-11600.	3.1	20
56	Charakterisierung von Wasserstoffbrücken zwischen Ionen in protischen ionischen Flüssigkeiten mittels NMR-Deuteron-Quadrupol-Kopplungskonstanten – Unterschiede zu Brücken in Amiden, Peptiden und Proteinen. <i>Angewandte Chemie</i> , 2017, 129, 14500-14505.	2.0	5
57	Defibrillation of soft porous metal-organic frameworks with electric fields. <i>Science</i> , 2017, 358, 347-351.	12.6	352
58	Probing Gas Adsorption in Metal-Organic Framework ZIF-8 by EPR of Embedded Nitroxides. <i>Journal of Physical Chemistry C</i> , 2017, 121, 19880-19886.	3.1	19
59	Characterization of Doubly Ionic Hydrogen Bonds in Protic Ionic Liquids by NMR Deuteron Quadrupole Coupling Constants: Differences to H-bonds in Amides, Peptides, and Proteins. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 14310-14314.	13.8	35
60	Porous Metal-Organic Polyhedral Frameworks with Optimal Molecular Dynamics and Pore Geometry for Methane Storage. <i>Journal of the American Chemical Society</i> , 2017, 139, 13349-13360.	13.7	99
61	Monitoring the Diffusivity of Light Hydrocarbons in a Mixture by Magic Angle Spinning Pulsed Field Gradient NMR: Methane/Ethane/Ethene in ZIF-8. <i>Journal of Physical Chemistry C</i> , 2017, 121, 25372-25376.	3.1	17
62	Mobility and Reactivity of 4-Substituted TEMPO Derivatives in Metal-Organic Framework MIL-53(Al). <i>Journal of Physical Chemistry C</i> , 2016, 120, 10698-10704.	3.1	23
63	Ultraslow Dynamics of a Framework Linker in MIL-53 (Al) as a Sensor for Different Isomers of Xylene. <i>Journal of Physical Chemistry C</i> , 2016, 120, 21704-21709.	3.1	27
64	Competitive pathways of methane activation on $\text{Zn}^{2+}$ -modified ZSM-5 zeolite: H/D hydrogen exchange with Brønsted acid sites versus dissociative adsorption to form Zn-methyl species. <i>Catalysis Science and Technology</i> , 2016, 6, 6381-6388.	4.1	28
65	Mobility of Stable $\pi$ -Complexes of Ethylene with $\text{Ag}^+$ Cations in Ag/H-ZSM-5 Zeolite: A $^{2+}\text{H}$ Solid-State NMR Study. <i>Journal of Physical Chemistry C</i> , 2016, 120, 4993-5000.	3.1	12
66	Methane Interaction with $\text{Zn}^{2+}$ -Exchanged Zeolite H-ZSM-5: Study of Adsorption and Mobility by One- and Two-Dimensional Variable-Temperature $^1\text{H}$ Solid-State NMR. <i>Journal of Physical Chemistry C</i> , 2015, 119, 14255-14261.	3.1	15
67	Metal-Cation-Independent Dynamics of Phenylene Ring in Microporous MOFs: A $^{2+}\text{H}$ Solid-State NMR Study. <i>Journal of Physical Chemistry C</i> , 2015, 119, 28038-28045.	3.1	36
68	Mobility of the 2-Methylimidazolate Linkers in ZIF-8 Probed by $^{2+}\text{H}$ NMR: Saloon Doors for the Guests. <i>Journal of Physical Chemistry C</i> , 2015, 119, 27512-27520.	3.1	97
69	Diffusion of $\text{CH}_4$ in ZIF-8 Studied by Quasi-Elastic Neutron Scattering. <i>Journal of Physical Chemistry C</i> , 2015, 119, 16115-16120.	3.1	30
70	Diffusion of Benzene in the Breathing Metal-Organic Framework MIL-53(Cr): A Joint Experimental-Computational Investigation. <i>Journal of Physical Chemistry C</i> , 2015, 119, 8217-8225.	3.1	38
71	Methane Activation on $\text{Zn}^{2+}$ -Exchanged ZSM-5 Zeolites. The Effect of Molecular Oxygen Addition. <i>Journal of Physical Chemistry C</i> , 2015, 119, 24910-24918.	3.1	67
72	Methane Mobility in Ag/H-ZSM-5 Zeolite in the Presence of Ethene: A View Based on PFG $^1\text{H}$ MAS NMR Analysis of Methane Diffusivity. <i>Journal of Physical Chemistry C</i> , 2015, 119, 18481-18486.	3.1	11

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73	Characterization and Dynamics of the Different Protonic Species in Hydrated 12-Tungstophosphoric Acid Studied by $^2\text{H}$ NMR. <i>Journal of Physical Chemistry C</i> , 2014, 118, 30023-30033.	3.1	25
74	Structural Dynamics in a "Breathing" Metal-Organic Framework Studied by Electron Paramagnetic Resonance of Nitroxide Spin Probes. <i>Journal of Physical Chemistry Letters</i> , 2014, 5, 20-24.	4.6	48
75	Guest Controlled Rotational Dynamics of Terephthalate Phenylenes in Metal-Organic Framework MIL-53(Al): Effect of Different Xylene Loadings. <i>Journal of Physical Chemistry C</i> , 2014, 118, 15978-15984.	3.1	42
76	Rotational and Translational Motion of Benzene in ZIF-8 Studied by $^2\text{H}$ NMR: Estimation of Microscopic Self-Diffusivity and Its Comparison with Macroscopic Measurements. <i>Journal of Physical Chemistry C</i> , 2014, 118, 12873-12879.	3.1	39
77	Solid-State NMR Characterization of Brønsted Acid Sites of Cesium Salt of 12-Tungstophosphoric Acid. <i>Journal of Physical Chemistry C</i> , 2014, 118, 21042-21048.	3.1	9
78	Methane Activation on In-Modified ZSM-5: The State of Indium in the Zeolite and Pathways of Methane Transformation to Surface Species. <i>Journal of Physical Chemistry C</i> , 2014, 118, 8034-8043.	3.1	47
79	Methane Activation on In-Modified ZSM-5 Zeolite. H/D Hydrogen Exchange of the Alkane with Brønsted Acid Sites. <i>Journal of Physical Chemistry C</i> , 2014, 118, 14427-14432.	3.1	25
80	Solid-state NMR monitoring of a double bond isomerization in propene on ZnO. <i>Chemical Physics Letters</i> , 2014, 607, 21-24.	2.6	4
81	Coaromatization of Methane with Propane on Mo-Containing Zeolite H-BEA: A Solid-State NMR and GC-MS Study. <i>Journal of Physical Chemistry C</i> , 2013, 117, 22867-22873.	3.1	20
82	Solid-state NMR study of the kinetics and mechanism of dimethyl ether carbonylation on cesium salt of 12-tungstophosphoric acid modified with Ag, Pt, and Rh. <i>Journal of Catalysis</i> , 2013, 308, 250-257.	6.2	20
83	Oxidation of methane to methanol on the surface of FeZSM-5 zeolite. <i>Journal of Catalysis</i> , 2013, 300, 47-54.	6.2	160
84	Diffusion of Xylene Isomers in the MIL-47(V) MOF Material: A Synergic Combination of Computational and Experimental Tools. <i>Journal of Physical Chemistry C</i> , 2013, 117, 6293-6302.	3.1	44
85	Parahydrogen-Induced Polarization Detected with Continuous Flow Magic Angle Spinning NMR. <i>Journal of Physical Chemistry C</i> , 2013, 117, 2888-2892.	3.1	25
86	Methane Activation and Transformation on Ag/H-ZSM-5 Zeolite Studied with Solid-State NMR. <i>Journal of Physical Chemistry C</i> , 2013, 117, 7690-7702.	3.1	72
87	Carbonylation of Dimethyl Ether with CO on Solid 12-Tungstophosphoric Acid: In Situ Magic Angle Spinning NMR Monitoring of the Reaction Kinetics. <i>Journal of Physical Chemistry C</i> , 2013, 117, 11168-11175.	3.1	9
88	Direct $^2\text{H}$ NMR Observation of the Proton Mobility of the Acidic Sites of Anhydrous 12-Tungstophosphoric Acid. <i>ChemPhysChem</i> , 2013, 14, 1783-1786.	2.1	16
89	Solid-State NMR Characterization of the Structure of Intermediates Formed from Olefins on Metal Oxides ( $\text{Al}_2\text{O}_3$ and $\text{Ga}_2\text{O}_3$ ). <i>Journal of Physical Chemistry C</i> , 2012, 116, 21430-21438.	3.1	30
90	Propene disproportionation on ZnO. <i>Chemical Physics Letters</i> , 2012, 552, 88-91.	2.6	2

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91	Experimental and Simulation Evidence of a Corkscrew Motion for Benzene in the Metal-Organic Framework MIL-47. <i>Journal of Physical Chemistry C</i> , 2012, 116, 15093-15098.	3.1	40
92	Probing the Dynamics of the Porous Zr Terephthalate UiO-66 Framework Using <sup>2</sup> H NMR and Neutron Scattering. <i>Journal of Physical Chemistry C</i> , 2012, 116, 12131-12136.	3.1	97
93	Mobility of <i>tert</i> -Butyl Alcohol in MFI Framework Type Studied by Deuterium NMR. <i>Journal of Physical Chemistry C</i> , 2012, 116, 8956-8963.	3.1	17
94	Structure of Allylic Intermediate on Zinc Oxide, $\tilde{\epsilon}$ or $\tilde{f}$ ?. <i>Journal of Physical Chemistry C</i> , 2012, 116, 11096-11099.	3.1	16
95	Carbonylation of dimethyl ether on Rh/Cs <sub>2</sub> HPW <sub>12</sub> O <sub>40</sub> : Solid-state NMR study of the mechanism of reaction in the presence of a methyl iodide promoter. <i>Journal of Catalysis</i> , 2012, 291, 9-16.	6.2	16
96	Hydrogen H/D Exchange and Activation of C <sub>1</sub> - <i>n</i> -C <sub>4</sub> Alkanes on Ga-Modified Zeolite BEA Studied with <sup>1</sup> H Magic Angle Spinning Nuclear Magnetic Resonance in Situ. <i>Journal of Physical Chemistry C</i> , 2011, 115, 13877-13886.	3.1	34
97	Mobility of Solid <i>tert</i> -Butyl Alcohol Studied by Deuterium NMR. <i>Journal of Physical Chemistry A</i> , 2011, 115, 7428-7436.	2.5	16
98	Carbonylation of dimethyl ether on solid Rh-promoted Cs-salt of Keggin 12-H <sub>3</sub> PW <sub>12</sub> O <sub>40</sub> : A solid-state NMR study of the reaction mechanism. <i>Journal of Catalysis</i> , 2011, 277, 72-79.	6.2	33
99	Comparison of the dynamics of MIL-53(Cr) and MIL-47(V) frameworks using neutron scattering and DFT methods. <i>European Physical Journal: Special Topics</i> , 2010, 189, 263-271.	2.6	31
100	Dynamics of Benzene Rings in MIL-53(Cr) and MIL-47(V) Frameworks Studied by <sup>2</sup> H-NMR Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 4791-4794.	13.8	127
101	Strong acidity of silanol groups of zeolite beta: Evidence from the studies by IR spectroscopy of adsorbed CO and <sup>1</sup> H MAS NMR. <i>Microporous and Mesoporous Materials</i> , 2010, 131, 210-216.	4.4	111
102	Metal-alkyl species are formed on interaction of small alkanes with gallium oxide: Evidence from solid-state NMR. <i>Chemical Physics Letters</i> , 2010, 496, 148-151.	2.6	24
103	Results of NMR spectroscopic studies of hydrocarbon conversions on solid acid catalysts in the last 25 years. <i>Kinetics and Catalysis</i> , 2010, 51, 854-872.	1.0	13
104	Propane Aromatization on Zn-Modified Zeolite BEA Studied by Solid-State NMR in Situ. <i>Journal of Physical Chemistry C</i> , 2010, 114, 12681-12688.	3.1	64
105	Mobility of <i>n</i> -Butane in ZSM-5 Zeolite Studied by <sup>2</sup> H NMR. <i>Journal of Physical Chemistry C</i> , 2010, 114, 2958-2966.	3.1	17
106	The $\sigma$ -Alkyl and $\sigma$ -Carbenium Pathways of Methane Activation on Ga-Modified Zeolite BEA: <sup>13</sup> C Solid-State NMR and GC-MS Study of Methane Aromatization in the Presence of Higher Alkane. <i>Journal of Physical Chemistry C</i> , 2010, 114, 21555-21561.	3.1	72
107	Hybrid Polyoxotungstate/MIL-101 Materials: Synthesis, Characterization, and Catalysis of H <sub>2</sub> O <sub>2</sub> -Based Alkene Epoxidation. <i>Inorganic Chemistry</i> , 2010, 49, 2920-2930.	4.0	228
108	H/D exchange of molecular hydrogen with Brønsted acid sites of Zn- and Ga-modified zeolite BEA. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 5149.	2.8	21

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109	Efficient generation of terahertz radiation by the method of optical rectification of terawatt laser pulses. Optics and Spectroscopy (English Translation of Optika I Spektroskopiya), 2009, 107, 529-533.	0.6	2
110	In situ high temperature MAS NMR study of the mechanisms of catalysis. Ethane aromatization on Zn-modified zeolite BEA. Solid State Nuclear Magnetic Resonance, 2009, 35, 113-119.	2.3	46
111	Water dynamics in bulk and dispersed in silica CaCl <sub>2</sub> hydrates studied by neutron scattering methods. Microporous and Mesoporous Materials, 2009, 125, 46-50.	4.4	4
112	Methane aromatization on Zn-modified zeolite in the presence of a co-reactant higher alkane: How does it occur?. Catalysis Today, 2009, 144, 265-272.	4.4	87
113	Reactivity of Methoxy Species toward CO on Keggin 12-H <sub>3</sub> PW <sub>12</sub> O <sub>40</sub> : A Study with Solid State NMR. Journal of Physical Chemistry C, 2009, 113, 19639-19644.	3.1	35
114	Significant Influence of Zn on Activation of the C-H Bonds of Small Alkanes by Brønsted Acid Sites of Zeolite. ChemPhysChem, 2008, 9, 2559-2563.	2.1	70
115	Understanding Methane Aromatization on a Zn-Modified High-Silica Zeolite. Angewandte Chemie - International Edition, 2008, 47, 4559-4562.	13.8	143
116	<sup>1</sup> H NMR signal broadening in spectra of alkane molecules adsorbed on MFI-type zeolites. Solid State Nuclear Magnetic Resonance, 2008, 33, 65-71.	2.3	12
117	Zn-promoted hydrogen exchange for methane and ethane on Zn/H-BEA zeolite: In situ <sup>1</sup> H MAS NMR kinetic study. Journal of Catalysis, 2008, 253, 11-21.	6.2	65
118	Water Dynamics in Bulk and Dispersed in Silica CaCl <sub>2</sub> Hydrates Studied by <sup>2</sup> H NMR. Journal of Physical Chemistry C, 2008, 112, 12853-12860.	3.1	21
119	Kinetics of H/D Exchange for n-Butane on Zeolite H-ZSM-5 Studied with <sup>1</sup> H MAS NMR In Situ. Journal of Physical Chemistry C, 2008, 112, 11869-11874.	3.1	29
120	Spectral modifications of femtosecond laser pulses induced by phase-matched optical rectification in LiNbO <sub>3</sub> , 2007, . .		0
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