Lidong Zhang

List of Publications by Year
in descending order

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129 papers	7,546 citations	66234 42 h-index	54797 84 g-index
130 all docs	130 docs citations	130 times ranked	7823 citing authors

#	Article	IF	Citations
1	Mesoporous Coâ€“Mn Spinel Oxides as Efficient Catalysts for Low Temperature Propane Oxidation. Catalysis Letters, 2022, 152, 2695-2704.	1.4	3
2	Experimental and kinetic modeling studies of 2-acetylfuran pyrolysis at atmospheric pressure. Combustion and Flame, 2022, 236, 111824.	2.8	6
3	Converting CO₂ Hydrogenation Products from Paraffins to Olefins: Modification of Zeolite Surface Properties by a UIO-n Membrane. ACS Catalysis, 2022, 12, 5894-5902.	5.5	10
4	A theoretical kinetics study on low-temperature oxidation of n-C4H9 radicals. Proceedings of the Combustion Institute, 2021, 38, 681-689.	2.4	12
5	Experimental and kinetic modeling study of methyl heptanoate low-temperature oxidation in a jet-stirred reactor. Fuel, 2021, 283, 118885.	3.4	10
6	Catalytic combustion of methyl butanoate over HZSM-5 zeolites. Chemical Communications, 2021, 57, 2233-2244.	2.2	11
7	Unique Coordination Structure of Cobalt Single-Atom Catalyst Supported on Dopant-Free Carbon. Journal of Physical Chemistry C, 2021, 125, 6735-6742.	1.5	1
8	Optimized rapid flame synthesis of morphology-controlled Î±-MoO3 layered nanoflakes. Journal of Nanoparticle Research, 2021, 23, 1.	0.8	5
9	Toward the Mechanism Study of Pd/Î³-Al₂O₃-Assisted Bioalcohol Combustion in a Flow Reactor. Energy & Fuels, 2021, 35, 14954-14962.	2.5	6
10	Unraveling the low-temperature oxidation mechanism between methyl crotonate radicals and O2. Combustion and Flame, 2021, 231, 111473.	2.8	9
11	Toward high-level theoretical studies on the reaction kinetics of PAHs growth based on HACA pathway: An ONIOM[G3(MP2,CC)//B3LYP:DFT] method developed. Fuel, 2021, 301, 121052.	3.4	15
12	Theoretical investigation of chemical reaction kinetics of CO2 and vinyl radical under catalytic combustion. Fuel, 2021, 305, 121566.	3.4	3
13	Flame In Situ Synthesis of Metal-Anchored CuO Nanowires for CO Catalytic Oxidation and Kinetic Analysis. ACS Applied Energy Materials, 2021, 4, 13226-13238.	2.5	9
14	Highly active and stable Co3O4 catalyst for the Low-temperature oxidative dehydrogenation of propane. Inorganic Chemistry Communication, 2020, 112, 107725.	1.8	9
15	Theoretical study of transition metals supported on g-C3N4 as electrochemical catalysts for CO2 reduction to CH3OH and CH4. Journal of CO2 Utilization, 2020, 36, 116-123.	3.3	59
16	Influence of Torsional Anharmonicity on the Reactions of Methyl Butanoate with Hydroperoxyl Radical. Journal of Physical Chemistry A, 2020, 124, 8643-8652.	1.1	8
17	Dynamic Surface Reconstruction of Single-Atom Bimetallic Alloy under Operando Electrochemical Conditions. Nano Letters, 2020, 20, 8319-8325.	4.5	28
18	Theoretical investigations of electrochemical CO₂ reduction by transition metals anchored on CNTs. Sustainable Energy and Fuels, 2020, 4, 6156-6164.	2.5	13

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19	Abatement of n-Butane by Catalytic Combustion over Co-ZSM-5 Catalysts. Energy & Fuels, 2020, 34, 12880-12890.	2.5	20
20	Theoretical calculation of low-temperature oxidation of heptyl radicals and O2. Combustion and Flame, 2020, 217, 274-284.	2.8	14
21	Au@Pt Nanotubes within CoZn-Based Metal-Organic Framework for Highly Efficient Semi-hydrogenation of Acetylene. IScience, 2020, 23, 101233.	1.9	12
22	Theoretical chemical kinetics for catalytic pyrolysis of methyl acetate over H-ZSM-5 zeolites. Fuel, 2020, 277, 118101.	3.4	7
23	A thorough theoretical mechanistic study of OH-initiated oxidative degradation mechanism for large polycyclic aromatic hydrocarbons. Computational and Theoretical Chemistry, 2020, 1175, 112730.	1.1	5
24	A theoretical investigation on Bell-Evans-Polanyi correlations for hydrogen abstraction reactions of large biodiesel molecules by H and OH radicals. Combustion and Flame, 2020, 214, 394-406.	2.8	18
25	Coordination-Engineered Cuâ€“N_x Single-Site Catalyst for Enhancing Oxygen Reduction Reaction. ACS Applied Energy Materials, 2019, 2, 6497-6504.	2.5	58
26	Interfacial engineering of cobalt sulfide/graphene hybrids for highly efficient ammonia electrosynthesis. Proceedings of the National Academy of Sciences of the United States of America, 2019, 116, 6635-6640.	3.3	242
27	Towards high-level theoretical studies of large biodiesel molecules: an ONIOM/RRKM/Master-equation approach to the isomerization and dissociation kinetics of methyl decanoate radicals. Physical Chemistry Chemical Physics, 2019, 21, 5232-5242.	1.3	13
28	Theoretical Study on Criegee Intermediateâ€™s Role in Ozonolysis of Acrylic Acid. Journal of Physical Chemistry A, 2019, 123, 1929-1936.	1.1	7
29	Experimental and Theoretical Investigation of the Pyrolysis of Furfural. Journal of Physical Chemistry A, 2019, 123, 103-110.	1.1	21
30	Utilization of generalized energy-based fragmentation method on the study of hydrogen abstraction reactions of large methyl esters. Combustion and Flame, 2018, 190, 467-476.	2.8	17
31	Experimental and modeling studies of small typical methyl esters pyrolysis: Methyl butanoate and methyl crotonate. Combustion and Flame, 2018, 191, 160-174.	2.8	48
32	Ab initio kinetics on low temperature oxidation of iso-pentane: The first oxygen addition. Combustion and Flame, 2018, 190, 119-132.	2.8	25
33	Theoretical studies on the reaction kinetics of methyl crotonate with hydroxyl radical. Sustainable Energy and Fuels, 2018, 2, 392-402.	2.5	24
34	Oxygen Vacancies Confined in Nickel Molybdenum Oxide Porous Nanosheets for Promoted Electrocatalytic Urea Oxidation. ACS Catalysis, 2018, 8, 1-7.	5.5	372
35	Toward High-Level Theoretical Studies of Large Biodiesel Molecules: An ONIOM [QCISD(T)/CBS:DFT] Study of the Reactions between Unsaturated Methyl Esters (C_nH_{2nâ€“1}COOCH₃) and Hydrogen Radical. Journal of Physical Chemistry A, 2018, 122, 4882-4893.	1.1	18
36	A theoretical kinetics study on low-temperature reactions of methyl acetate radicals with molecular oxygen. Combustion and Flame, 2018, 196, 45-53.	2.8	20

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37	Experimental and kinetic modeling investigation on methyl decanoate pyrolysis at low and atmospheric pressures. Fuel, 2018, 232, 333-340.	3.4	27
38	Theoretical Studies on Isomerization and Decomposition Reactions of 2-Methyl-1-butanol Radicals. Energy & Fuels, 2018, 32, 7652-7659.	2.5	3
39	Predictive kinetics on the formation and decomposition of ethylbenzene. Proceedings of the Combustion Institute, 2017, 36, 533-542.	2.4	15
40	Hydroperoxide Measurements During Low-Temperature Gas-Phase Oxidation of n-Heptane and n-Decane. Journal of Physical Chemistry A, 2017, 121, 1861-1876.	1.1	31
41	Theoretical kinetic studies for low temperature oxidation of two typical methylcyclohexyl radicals. Combustion and Flame, 2017, 182, 216-224.	2.8	32
42	Atomically Dispersed Ironâ€“Nitrogen Species as Electrocatalysts for Bifunctional Oxygen Evolution and Reduction Reactions. Angewandte Chemie - International Edition, 2017, 56, 610-614.	7.2	950
43	Experimental and kinetic modeling studies of furan pyrolysis: Fuel decomposition and aromatic ring formation. Fuel, 2017, 206, 239-247.	3.4	38
44	Regulating Waterâ€Reduction Kinetics in Cobalt Phosphide for Enhancing HER Catalytic Activity in Alkaline Solution. Advanced Materials, 2017, 29, 1606980.	11.1	220
45	3D Nitrogenâ€Anionâ€Decorated Nickel Sulfides for Highly Efficient Overall Water Splitting. Advanced Materials, 2017, 29, 1701584.	11.1	478
46	Pyrolysis of n-butane investigated using synchrotron threshold photoelectron photoion coincidence spectroscopy. RSC Advances, 2017, 7, 28746-28753.	1.7	7
47	Spin-State Regulation of Perovskite Cobaltite to Realize Enhanced Oxygen Evolution Activity. CheM, 2017, 3, 812-821.	5.8	225
48	Roaming-Mediated CH₂NH Elimination from the Ionization of Aromatic Ethylamines. ChemistryOpen, 2017, 6, 40-45.	0.9	2
49	Theoretical studies for reaction kinetics of cy-C6H11CH2 radical with O2. Proceedings of the Combustion Institute, 2017, 36, 179-186.	2.4	25
50	New insights into the low-temperature oxidation of 2-methylhexane. Proceedings of the Combustion Institute, 2017, 36, 373-382.	2.4	36
51	Experimental and kinetic modeling study of laminar premixed decalin flames. Proceedings of the Combustion Institute, 2017, 36, 1193-1202.	2.4	14
52	Phaseâ€Transformation Engineering in Cobalt Diselenide Realizing Enhanced Catalytic Activity for Hydrogen Evolution in an Alkaline Medium. Advanced Materials, 2016, 28, 7527-7532.	11.1	307
53	Experimental and Kinetic Modeling Study of Nitroethane Pyrolysis at a Low Pressure: Competition Reactions in the Primary Decomposition. Energy & Fuels, 2016, 30, 7738-7745.	2.5	12
54	Metallic Nickel Hydroxide Nanosheets Give Superior Electrocatalytic Oxidation of Urea for Fuel Cells. Angewandte Chemie - International Edition, 2016, 55, 12465-12469.	7.2	356

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55	Manganous oxide nanoparticles encapsulated in few-layer carbon as an efficient electrocatalyst for oxygen reduction in alkaline media. Journal of Materials Chemistry A, 2016, 4, 11775-11781.	5.2	27
56	Experimental and kinetic modeling investigation on decalin pyrolysis at low to atmospheric pressures. Combustion and Flame, 2016, 167, 228-237.	2.8	18
57	Additional chain-branching pathways in the low-temperature oxidation of branched alkanes. Combustion and Flame, 2016, 164, 386-396.	2.8	94
58	Online Study on the Catalytic Pyrolysis of Bituminous Coal over HUSY and HZSM-5 with Photoionization Time-of-Flight Mass Spectrometry. Energy & Fuels, 2016, 30, 1598-1604.	2.5	45
59	Pyrolysis of 2-methyl-1-butanol at low and atmospheric pressures: Mass spectrometry and modeling studies. Proceedings of the Combustion Institute, 2015, 35, 409-417.	2.4	14
60	Experimental and kinetic modeling study of premixed o-xylene flames. Proceedings of the Combustion Institute, 2015, 35, 1745-1752.	2.4	45
61	Investigation on primary decomposition of ethylcyclohexane at atmospheric pressure. Proceedings of the Combustion Institute, 2015, 35, 367-375.	2.4	47
62	Experimental and kinetic modeling study of the low- and intermediate-temperature oxidation of dimethyl ether. Combustion and Flame, 2015, 162, 1113-1125.	2.8	120
63	Kinetics of ethylcyclohexane pyrolysis and oxidation: An experimental and detailed kinetic modeling study. Combustion and Flame, 2015, 162, 2873-2892.	2.8	70
64	Molecular co-catalyst accelerating hole transfer for enhanced photocatalytic H2 evolution. Nature Communications, 2015, 6, 8647.	5.8	172
65	Towards high-level theoretical studies of large biodiesel molecules: an ONIOM [QCISD(T)/CBS:DFT] study of hydrogen abstraction reactions of C_nH_{2n+1}COOC_mH_{2m+1} + H. Physical Chemistry Chemical Physics, 2015, 17, 200-208.	1.3	26
66	Hydrogen dangling bonds induce ferromagnetism in two-dimensional metal-free graphitic-C₃N₄ nanosheets. Chemical Science, 2015, 6, 283-287.	3.7	62
67	Density functional theory study of the reactions of 2-azido- N , N -dimethylethanamine with nitric acid and nitrogen dioxide. Combustion and Flame, 2015, 162, 237-248.	2.8	8
68	A theoretical kinetics study of the reactions of methylbutanoate with hydrogen and hydroxyl radicals. Proceedings of the Combustion Institute, 2015, 35, 481-489.	2.4	40
69	Experimental and kinetic modeling study of i-butanol pyrolysis and combustion. Combustion and Flame, 2014, 161, 1955-1971.	2.8	28
70	Experimental and kinetic modeling study on methylcyclohexane pyrolysis and combustion. Combustion and Flame, 2014, 161, 84-100.	2.8	126
71	Large Negative Magnetoresistance Induced by Anionic Solid Solutions in Two-Dimensional Spin-Frustrated Transition Metal Chalcogenides. Physical Review Letters, 2014, 113, 157202.	2.9	39
72	Kinetics of Homoallylic/Homobenzylic Rearrangement Reactions under Combustion Conditions. Journal of Physical Chemistry A, 2014, 118, 6741-6748.	1.1	16

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73	Experimental and kinetic modeling study of pyrolysis and oxidation of n-decane. Combustion and Flame, 2014, 161, 1701-1715.	2.8	87
74	Interactions of disulfide-constrained cyclic tetrapeptides with Cu2+. Journal of Biological Inorganic Chemistry, 2013, 18, 277-286.	1.1	4
75	Experimental and kinetic modeling study of tetralin pyrolysis at low pressure. Proceedings of the Combustion Institute, 2013, 34, 1739-1748.	2.4	53
76	An experimental and modeling study of methyl propanoate pyrolysis at low pressure. Combustion and Flame, 2013, 160, 1958-1966.	2.8	50
77	Experimental and kinetic modeling study of 2-butanol pyrolysis and combustion. Combustion and Flame, 2013, 160, 1939-1957.	2.8	58
78	An experimental and theoretical study of pyrrolidine pyrolysis at low pressure. Proceedings of the Combustion Institute, 2013, 34, 641-648.	2.4	9
79	An experimental and kinetic modeling study of premixed nitroethane flames at low pressure. Proceedings of the Combustion Institute, 2013, 34, 617-624.	2.4	29
80	An experimental and kinetic modeling investigation on a rich premixed n-propylbenzene flame at low pressure. Proceedings of the Combustion Institute, 2013, 34, 1785-1793.	2.4	41
81	Experimental and theoretical studies of pyrolysis of chrysophanol and its derivatives. Journal of Analytical and Applied Pyrolysis, 2013, 100, 237-244.	2.6	2
82	Kinetics of Decomposition and Isomerization of Methylcyclohexane: Starting Point for Studying Monoalkylated Cyclohexanes Combustion. Energy & Fuels, 2013, 27, 1679-1687.	2.5	44
83	Product Identification and Mass Spectrometric Analysis of n-Butane and i-Butane Pyrolysis at Low Pressure. Chinese Journal of Chemical Physics, 2013, 26, 151-156.	0.6	9
84	Experimental and Theoretical Study on Pyrolysis of Isopsoralen. Chinese Journal of Chemical Physics, 2012, 25, 249-253.	0.6	1
85	VUV Photoionization and Dissociation of Tyramine and Dopamine: the Joint Experimental and Theoretical Studies. Chinese Journal of Chemical Physics, 2012, 25, 11-18.	0.6	5
86	Thermal Decomposition of 1-Pentanol and Its Isomers: A Theoretical Study. Journal of Physical Chemistry A, 2012, 116, 9238-9244.	1.1	44
87	Experimental and Kinetic Modeling Study of n-Butanol Pyrolysis and Combustion. Energy & Fuels, 2012, 26, 5550-5568.	2.5	123
88	Experimental and kinetic modeling study of tert-butanol combustion at low pressure. Energy, 2012, 43, 94-102.	4.5	29
89	Theoretical Studies on the Unimolecular Decomposition of Propanediols and Glycerol. Journal of Physical Chemistry A, 2012, 116, 4457-4465.	1.1	13
90	Theoretical Studies on the Unimolecular Decomposition of Ethylene Glycol. Journal of Physical Chemistry A, 2012, 116, 55-63.	1.1	30

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91	Synchrotron vacuum ultraviolet (VUV) photo-induced fragmentation of cyclic dipeptides radical cations. Amino Acids, 2012, 43, 279-287.	1.2	2
92	An experimental and kinetic modeling study of three butene isomers pyrolysis at low pressure. Combustion and Flame, 2012, 159, 905-917.	2.8	141
93	An experimental and kinetic modeling study of cyclohexane pyrolysis at low pressure. Combustion and Flame, 2012, 159, 2243-2253.	2.8	110
94	Pyrolysis ofn-Heptane: Experimental and Theoretical Study. Journal of Physical Chemistry A, 2011, 115, 1593-1601.	1.1	50
95	Unimolecular Decomposition of Ethyl Hydroperoxide: Ab Initio/Riceâ"Ramspergerâ"Kasselâ"Marcus Theoretical Prediction of Rate Constants. Journal of Physical Chemistry A, 2011, 115, 602-611.	1.1	19
96	Experimental and theoretical studies on decomposition of pyrrolidine. Proceedings of the Combustion Institute, 2011, 33, 415-423.	2.4	16
97	Experimental and modeling investigation on premixed ethylbenzene flames at low pressure. Proceedings of the Combustion Institute, 2011, 33, 617-624.	2.4	56
98	Investigation on chemical structures of premixed toluene flames at low pressure. Proceedings of the Combustion Institute, 2011, 33, 593-600.	2.4	113
99	VUV photonâ€induced ionization/dissociation of antipyrine and propyphenazone: mass spectrometric and theoretical insights. Journal of Mass Spectrometry, 2010, 45, 734-739.	0.7	9
100	Investigation on fuel-rich premixed flames of monocyclic aromatic hydrocarbons: Part I. Intermediate identification and mass spectrometric analysis. Combustion and Flame, 2010, 157, 143-154.	2.8	83
101	Kinetic modeling study of toluene pyrolysis at low pressure. Combustion and Flame, 2010, 157, 1686-1697.	2.8	111
102	Determination of absolute photoionization crossâ€sections of alkanes and cycloâ€alkanes. Rapid Communications in Mass Spectrometry, 2010, 24, 1335-1342.	0.7	53
103	Photoionisation and photodissociation studies of nonvolatile organic molecules by synchrotron VUV photoionisation mass spectrometry and theoretical calculations. International Reviews in Physical Chemistry, 2010, 29, 369-401.	0.9	20
104	Vacuum Ultraviolet Photofragmentation of Sarcosine: Photoionization Mass Spectrometric and Theoretical Insights. Journal of Physical Chemistry A, 2010, 114, 3411-3417.	1.1	14
105	THEORETICAL STUDIES ON PHOTOIONIZATION OF GUANINE TAUTOMERS AND INTERCONVERSION OF CATION RADICALS. Journal of Theoretical and Computational Chemistry, 2009, 08, 1103-1115.	1.8	6
106	Photoionization and Dissociative Photoionization Study of Cholesterol by IR Laser Desorption/Tunable Synchrotron VUV Photoionization Mass Spectrometry. Chinese Journal of Chemical Physics, 2009, 22, 129-133.	0.6	3
107	Identification of Intermediates in Pyridine Pyrolysis with Molecular-beam Mass Spectrometry and Tunable Synchrotron VUV Photoionization. Chinese Journal of Chemical Physics, 2009, 22, 204-209.	0.6	12
108	An experimental study of the rich premixed ethylbenzene flame at low pressure. Proceedings of the Combustion Institute, 2009, 32, 647-655.	2.4	51

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109	An experimental and kinetic modeling study of a premixed nitromethane flame at low pressure. Proceedings of the Combustion Institute, 2009, 32, 311-318.	2.4	70
110	Investigation of the rich premixed laminar acetylene/oxygen/argon flame: Comprehensive flame structure and special concerns of polyynes. Proceedings of the Combustion Institute, 2009, 32, 1293-1300.	2.4	66
111	An experimental and kinetic modeling study of premixed NH3/CH4/O2/Ar flames at low pressure. Combustion and Flame, 2009, 156, 1413-1426.	2.8	359
112	An experimental and theoretical study of toluene pyrolysis with tunable synchrotron VUV photoionization and molecular-beam mass spectrometry. Combustion and Flame, 2009, 156, 2071-2083.	2.8	111
113	Conformation-Specific Pathways of Î²-Alanine: A Vacuum Ultraviolet Photoionization and Theoretical Study. Journal of Physical Chemistry A, 2009, 113, 5838-5845.	1.1	22
114	An Experimental and Theoretical Study of Pyrrole Pyrolysis with Tunable Synchrotron VUV Photoionization and Molecular-Beam Mass Spectrometry. Journal of Physical Chemistry A, 2009, 113, 5397-5405.	1.1	35
115	Experimental Study of a Fuel-Rich Premixed Toluene Flame at Low Pressure. Energy & Fuels, 2009, 23, 1473-1485.	2.5	184
116	Intramolecular hydrogen transfer in the ionization process of Î±-alanine. Physical Chemistry Chemical Physics, 2009, 11, 1189.	1.3	22
117	Photoionization studies on various quinones by an infrared laser desorption/tunable VUV photoionization TOF mass spectrometry. Journal of Mass Spectrometry, 2008, 43, 1701-1710.	0.7	25
118	Dissociative Photoionization Mechanism of 1,8-Dihydroxyanthraquinone: An Experimental and Theoretical Study. Journal of Physical Chemistry A, 2008, 112, 10977-10984.	1.1	7
119	Pyrolysis of Methyl tert-Butyl Ether (MTBE). 2. Theoretical Study of Decomposition Pathways. Journal of Physical Chemistry A, 2008, 112, 10495-10501.	1.1	19
120	Pyrolysis of Methyl tert-Butyl Ether (MTBE). 1. Experimental Study with Molecular-Beam Mass Spectrometry and Tunable Synchrotron VUV Photoionization. Journal of Physical Chemistry A, 2008, 112, 10487-10494.	1.1	74
121	Theoretical Study of the AlEt3-Promoted Tandem Reductive Rearrangement of Epoxides. Chinese Journal of Chemical Physics, 2008, 21, 547-554.	0.6	3
122	Molecular Dynamics and Density Functional Studies of Substrate Binding and Catalysis of Arginine Deiminase. Journal of Physical Chemistry B, 2007, 111, 3267-3273.	1.2	10
123	Supermolecule density functional calculations suggest a key role for solvent in alkaline hydrolysis of p-nitrophenyl phosphate. Chemical Communications, 2007, , 1638.	2.2	42
124	Electronic Excitations of Green Fluorescent Proteins:â€‰ Modeling Solvatochromatic Shifts of Red Fluorescent Protein Chromophore Model Compound in Aqueous Solutions. Journal of Physical Chemistry B, 2007, 111, 14055-14063.	1.2	21
125	A computational investigation on the sequential rearrangement mechanism of 2-allyl-2,4,5-hexatrienaldehyde involving [1,5]-hydrogen migration and 8Ï€-electrocyclization. Journal of Computational Chemistry, 2007, 28, 2164-2169.	1.5	5
126	ELECTRONIC EXCITATIONS OF GREEN FLUORESCENT PROTEINS: MODELING SOLVATOCHROMATIC SHIFTS OF CHROMOPHORE MODEL COMPOUNDS IN SOLUTIONS. Journal of Theoretical and Computational Chemistry, 2006, 05, 375-390.	1.8	7

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127	Density functional theory study of proton transfer in carbonic anhydrase. Science Bulletin, 2005, 50, 2557.	1.7	0
128	Theoretical Study of General Base-Catalyzed Hydrolysis of Aryl Esters and Implications for Enzymatic Reactions. Journal of Physical Chemistry B, 2005, 109, 5259-5266.	1.2	11
129	Reactivity of Metaphosphate and Thiometaphosphate in Water:â€‰ A DFT Study. Journal of Physical Chemistry A, 2005, 109, 11295-11303.	1.1	14