

Yoshifumi Kimura

List of Publications by Year in descending order

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93
papers

1,626
citations

279487

23
h-index

377514

34
g-index

96
all docs

96
docs citations

96
times ranked

1118
citing authors

#	ARTICLE	IF	CITATIONS
1	Excited State Intramolecular Proton Transfer Reaction of 4- <i>N,N</i> -Diethylamino-3-hydroxyflavone and Solvation Dynamics in Room Temperature Ionic Liquids Studied by Optical Kerr Gate Fluorescence Measurement. <i>Journal of Physical Chemistry B</i> , 2010, 114, 11847-11858.	1.2	79
2	Comparative Study on the Structural, Optical, and Electrochemical Properties of Bithiophene-Fused Benzo[<i>b</i>]phospholes. <i>Chemistry - A European Journal</i> , 2008, 14, 8102-8115.	1.7	75
3	Raman Spectroscopic Study on Solvation of Diphenylcyclopropanone and Phenol Blue in Room Temperature Ionic Liquids. <i>Journal of Physical Chemistry A</i> , 2006, 110, 6164-6172.	1.1	52
4	Synthesis and Structure-Property Relationships of 2,2-Bis(benzo[<i>b</i>]phosphole) and 2,2-Benzo[<i>b</i>]phosphole-Benzo[<i>b</i>]heterole Hybrid π -Systems. <i>Chemistry - A European Journal</i> , 2012, 18, 15972-15983.	1.7	52
5	Solvent and Solvent Density Effects on the Spectral Shifts and the Bandwidths of the Absorption and the Resonance Raman Spectra of Phenol Blue. <i>Journal of Physical Chemistry A</i> , 1997, 101, 9050-9060.	1.1	49
6	Comparison of 2-Arylnaphtho[2,3- <i>b</i>]phospholes and 2-Arylbenzo[<i>b</i>]phospholes: Effects of 2-Aryl Groups and Fused Arene Moieties on Their Optical and Photophysical Properties. <i>Journal of Organic Chemistry</i> , 2015, 80, 5944-5950.	1.7	46
7	Chemical equilibrium in fluids from the gaseous to liquid states: Solvent density dependence of the dimerization equilibrium of 2-methyl-2-nitrosopropane in carbon dioxide, chlorotrifluoromethane, and trifluoromethane. <i>Journal of Chemical Physics</i> , 1992, 96, 3085-3091.	1.2	45
8	Excitation Wavelength Dependence of Excited State Intramolecular Proton Transfer Reaction of 4- <i>N,N</i> -Diethylamino-3-hydroxyflavone in Room Temperature Ionic Liquids Studied by Optical Kerr Gate Fluorescence Measurement. <i>Journal of Physical Chemistry B</i> , 2013, 117, 12567-12582.	1.2	41
9	Ab Initio Study on an Excited-State Intramolecular Proton-Transfer Reaction in Ionic Liquid. <i>Journal of Physical Chemistry B</i> , 2013, 117, 6759-6767.	1.2	41
10	The role of the attractive and the repulsive interactions in the nonpolar solvation dynamics in simple fluids from the gas-like to the liquid-like densities. <i>Journal of Chemical Physics</i> , 1999, 111, 4169-4185.	1.2	40
11	Chemical reaction in medium density fluid. Solvent density effects on the dimerization equilibrium of 2-methyl-2-nitrosopropane in carbon dioxide. <i>Journal of Chemical Physics</i> , 1989, 90, 5679-5686.	1.2	37
12	Effect of solvent density and species on static and dynamic fluorescence Stokes shifts of coumarin 153. <i>Journal of Chemical Physics</i> , 1999, 111, 5474-5484.	1.2	37
13	Raman Spectroscopic Study on the Solvation of <i>N,N</i> -Dimethyl- <i>p</i> -nitroaniline in Room-Temperature Ionic Liquids. <i>Journal of Physical Chemistry A</i> , 2007, 111, 7081-7089.	1.1	36
14	Solvent Effects on the Local Structure of <i>p</i> -Nitroaniline in Supercritical Water and Supercritical Alcohols. <i>Journal of Physical Chemistry A</i> , 2008, 112, 5515-5526.	1.1	34
15	Preparation of Gold Nanoparticles by the Laser Ablation in Room-temperature Ionic Liquids. <i>Chemistry Letters</i> , 2007, 36, 1130-1131.	0.7	33
16	Chemical equilibrium in simple fluids: Solvent density dependence of the dimerization equilibrium of 2-methyl-2-nitrosopropane in argon and xenon. <i>Journal of Chemical Physics</i> , 1992, 96, 3824-3829.	1.2	32
17	Solvation state selective excitation in resonance Raman spectroscopy. I. Experimental study on the C=N and the C=O stretching modes of phenol blue. <i>Journal of Chemical Physics</i> , 1998, 109, 9075-9083.	1.2	32
18	Synthesis of 2-Alkenyl- and 2-Alkynyl-benzo[<i>b</i>]phospholes by Using Palladium-Catalyzed Cross-Coupling Reactions. <i>Organic Letters</i> , 2013, 15, 4458-4461.	2.4	31

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19	Translational diffusion of hydrophobic solutes in supercritical water studied by molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 2003, 119, 7328-7334.	1.2	29
20	Universality of Viscosity Dependence of Translational Diffusion Coefficients of Carbon Monoxide, Diphenylacetylene, and Diphenylcyclopropanone in Ionic Liquids under Various Conditions. <i>Journal of Physical Chemistry B</i> , 2015, 119, 8096-8103.	1.2	29
21	Study on the excited state intramolecular proton transfer of 4- <i>N,N</i> -diethylamino-3-hydroxyflavone in imidazolium-based room temperature ionic liquids. <i>Chemical Physics Letters</i> , 2008, 463, 364-368.	1.2	28
22	Solvent Density Dependence of Translational Diffusion of Transient Radicals in the Medium-Density Region of Trifluoromethane and Carbon Dioxide. <i>Journal of Physical Chemistry B</i> , 1997, 101, 4442-4447.	1.2	27
23	Excitation wavelength dependence of photo-induced intramolecular proton transfer reaction of 4- <i>N,N</i> -diethylamino-3-hydroxyflavone in various liquids. <i>Chemical Physics Letters</i> , 2012, 531, 70-74.	1.2	24
24	Solute-solute potential of mean force in simple fluids at infinite dilution. <i>Molecular Physics</i> , 1991, 72, 279-294.	0.8	23
25	Acceptor Number of Room Temperature Ionic Liquid Determined by the Raman Spectrum of Diphenylcyclopropanone. <i>Chemistry Letters</i> , 2005, 34, 338-339.	0.7	23
26	Effect of the solvent density and species on the back-electron transfer rate in the hexamethylbenzene/tetracyanoethylene charge-transfer complex. <i>Journal of Chemical Physics</i> , 1998, 108, 1485-1498.	1.2	22
27	Solute-solvent hydrogen-bonding in room temperature ionic liquids studied by Raman spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 13676.	1.3	22
28	Solvation state selective excitation in resonance Raman spectroscopy. II. Theoretical calculation. <i>Journal of Chemical Physics</i> , 1998, 109, 9084-9095.	1.2	21
29	Raman Spectroscopic Study on the Solvation of <i>p</i> -Aminobenzonitrile in Supercritical Water and Methanol. <i>Journal of Physical Chemistry A</i> , 2009, 113, 3143-3154.	1.1	21
30	Application of the Transient Grating Method to the Measurement of Transport Properties for High Pressure Fluids. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , 1995, 99, 196-203.	0.9	20
31	Polarity and Nonpolarity of Ionic Liquids Viewed from the Rotational Dynamics of Carbon Monoxide. <i>Journal of Physical Chemistry B</i> , 2015, 119, 15493-15501.	1.2	20
32	The excitation energy dependence of the Raman-Stokes shift: The resonance Raman spectra of phenol blue in methanol. <i>Journal of Chemical Physics</i> , 1997, 107, 4436-4438.	1.2	19
33	Vibrational energy relaxation of azulene studied by the transient grating method. I. Supercritical fluids. <i>Journal of Chemical Physics</i> , 2005, 123, 054512.	1.2	19
34	Effects of counter anions, P-substituents, and solvents on optical and photophysical properties of 2-phenylbenzo[b]phospholium salts. <i>Dalton Transactions</i> , 2017, 46, 9517-9527.	1.6	18
35	Synthesis of zero-valent iron nanoparticles via laser ablation in a formate ionic liquid under atmospheric conditions. <i>Chemical Communications</i> , 2018, 54, 7834-7837.	2.2	18
36	Excitation wavelength dependence of the Raman-Stokes shift of <i>N,N</i> -dimethyl- <i>p</i> -nitroaniline. <i>Journal of Chemical Physics</i> , 2006, 124, 184503.	1.2	17

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37	Origin of low melting point of ionic liquids: dominant role of entropy. <i>Chemical Science</i> , 2022, 13, 7560-7565.	3.7	16
38	Volume Profile of the Reversible Dimerization Reaction of 2-Methyl-2-Nitrosopropane: Validity of the Hard Sphere Fluid Model and the Transition State Theory. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , 1988, 92, 1095-1103.	0.9	15
39	Charge Effect on the Diffusion Coefficient and the Bimolecular Reaction Rate of Diiodide Anion Radical in Room Temperature Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2009, 113, 5188-5193.	1.2	15
40	Effects of boryl, phosphino, and phosphonio substituents on optical, electrochemical, and photophysical properties of 2,5-dithienylphospholes and 2-phenyl-5-thienylphospholes. <i>Dalton Transactions</i> , 2016, 45, 2190-2200.	1.6	15
41	Vibrational energy relaxation of azulene in the S2 state. I. Solvent species dependence. <i>Journal of Chemical Physics</i> , 2000, 113, 2772-2783.	1.2	14
42	Excitation Wavelength Dependence of the Solvation Dynamics of 4-N,N-Diethylamino-3-methoxyflavon in Ionic Liquids. <i>Bulletin of the Chemical Society of Japan</i> , 2015, 88, 939-945.	2.0	14
43	Effect of Temperature and Water Concentration on CO2 Absorption by Tetrabutylphosphonium Formate Ionic Liquid. <i>Journal of Chemical & Engineering Data</i> , 2016, 61, 837-845.	1.0	14
44	Photodissociation Quantum Yield of Iodine in the Low-, Medium-, and High-Density Fluids Studied by the Transient Grating Method. <i>Journal of Physical Chemistry A</i> , 1999, 103, 7730-7741.	1.1	13
45	Solvophobic interaction in a simple associated fluid. <i>Molecular Physics</i> , 1992, 76, 737-755.	0.8	12
46	Study on the Chemical Reaction of Spiropyran in Medium- and High-Density Fluids. <i>The Journal of Physical Chemistry</i> , 1996, 100, 11009-11013.	2.9	12
47	Diffusion of Transient Radicals in Alcohols and Cyclohexane from Ambient to Supercritical Conditions Studied by the Transient Grating Method. <i>Journal of Physical Chemistry B</i> , 2003, 107, 5958-5966.	1.2	12
48	Transport Properties and Solvation Structure of Mixtures of Carbon Dioxide and Room-Temperature Ionic Liquids. <i>Bulletin of the Chemical Society of Japan</i> , 2011, 84, 70-78.	2.0	12
49	Study of the Excited-State Proton-Transfer Reaction of 5-Cyano-2-naphthol in Sub- and Supercritical Water. <i>Journal of Physical Chemistry B</i> , 2012, 116, 1043-1052.	1.2	12
50	Electron Transfer Reaction Dynamics of p-Nitroaniline in Water from Liquid to Supercritical Conditions. <i>Journal of Physical Chemistry B</i> , 2012, 116, 11508-11516.	1.2	12
51	Anomalous ground-state proton transfer of 4-N,N-diethylamino-3-hydroxyflavone in ionic liquids of imidazolium-based cations with tetrafluoroborate. <i>Chemical Communications</i> , 2013, 49, 3976.	2.2	12
52	Synthesis of 3,5-Disubstituted BODIPYs Bearing N-Containing Five-Membered Heteroaryl Groups via Nucleophilic C-N Bond Formation. <i>Journal of Organic Chemistry</i> , 2018, 83, 5274-5281.	1.7	11
53	Excited-State Proton Transfer of 5,8-Dicyano-2-naphthol in High-Temperature and High-Pressure Methanol: Effect of Solvent Polarity and Hydrogen Bonding Ability. <i>Journal of Physical Chemistry B</i> , 2018, 122, 12363-12374.	1.2	11
54	Understanding Structural Changes through Excited-State Intramolecular Proton Transfer in 4-N,N-Diethylamino-3-hydroxyflavone (DEAHF) in Solution Based on Quantum Chemical Calculations. <i>Journal of Physical Chemistry B</i> , 2019, 123, 9872-9881.	1.2	11

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55	Photo-dissociation dynamics of bis(p-dimethylaminophenyl) disulfide in ionic liquids studied by ultrafast transient absorption spectroscopy. <i>Chemical Physics Letters</i> , 2013, 564, 21-25.	1.2	10
56	Systematic estimation and interpretation of fractional free volume in 1-alkyl-3-methylimidazolium-based ionic liquids. <i>Fluid Phase Equilibria</i> , 2019, 498, 144-150.	1.4	10
57	Vibrational energy relaxation of azulene in the S2 state. II. Solvent density dependence. <i>Journal of Chemical Physics</i> , 2000, 113, 4340-4348.	1.2	9
58	Effects of solute-solvent and solvent-solvent attractive interactions on solute diffusion. <i>Molecular Physics</i> , 2000, 98, 1553-1563.	0.8	9
59	Non-Gaussian dynamics of a dilute hard-sphere gas. <i>Journal of Chemical Physics</i> , 2001, 114, 3029-3034.	1.2	9
60	Chemisorption of Carbon Dioxide in Carboxylate-Functionalized Ionic Liquids: A Mechanistic Study. <i>Chemistry Letters</i> , 2014, 43, 626-628.	0.7	9
61	Non-flammable and Highly Concentrated Carbonate Ester-free Electrolyte Solutions for 5 V Class Positive Electrodes in Lithium-ion Batteries. <i>ChemSusChem</i> , 2021, 14, 2445-2451.	3.6	9
62	Photo-excitation dynamics of Phenol Blue. <i>Physical Chemistry Chemical Physics</i> , 2000, 2, 1415-1420.	1.3	8
63	Excited-State Proton Transfer of Cyanonaphthols in Protic Ionic Liquids: Appearance of a New Fluorescent Species. <i>Journal of Physical Chemistry B</i> , 2017, 121, 6042-6049.	1.2	8
64	Extension of Anodic Potential Window of Ester-Based Electrolyte Solutions for High-Voltage Lithium Ion Batteries. <i>ACS Applied Energy Materials</i> , 2019, 2, 7728-7732.	2.5	8
65	Photo-excitation dynamics of N, N-dimethyl-p-nitroaniline in ionic liquids: Effect of cation alkyl-chain length. <i>Journal of Molecular Liquids</i> , 2019, 289, 111128.	2.3	8
66	Heterogeneous Structures of Ionic Liquids as Probed by CO Rotation with Nuclear Magnetic Resonance Relaxation Analysis and Molecular Dynamics Simulations. <i>Journal of Physical Chemistry B</i> , 2020, 124, 10465-10476.	1.2	8
67	Non-linear Laser Spectroscopy in Supercritical Fluids. <i>Review of High Pressure Science and Technology/Koatsuryoku No Kagaku To Gijutsu</i> , 2006, 16, 87-94.	0.1	7
68	Regioselective functionalization at the 7-position of 1,2,3-triphenylbenzo[b]phosphole oxide via Pd-directed lithiation. <i>Dalton Transactions</i> , 2018, 47, 7123-7127.	1.6	7
69	Experimental observation of the unique solvation process along multiple solvation coordinates of photodissociated products. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 4569-4579.	1.3	7
70	Translational Diffusion of a Transient Charge-separated Species in Carbon Dioxide and Trifluoromethane Studied by the Transient Grating Method. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , 1996, 100, 656-660.	0.9	6
71	SO ₂ capture by ionic liquid and spectroscopic speciation of sulfur(IV) therein. <i>RSC Advances</i> , 2017, 7, 6538-6547.	1.7	6
72	Photoexcitation dynamics of p-nitroaniline and N,N-dimethyl-p-nitroaniline in 1-alkyl-3-methylimidazolium-cation based ionic liquids with different alkyl-chain lengths. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 22161-22168.	1.3	6

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73	Application of the time-resolved spectroscopy on the photo-dissociation dynamics of disulfide compounds in supercritical fluids. <i>Journal of Molecular Liquids</i> , 2005, 119, 113-117.	2.3	5
74	Excited-State Intramolecular Proton Transfer Reaction and Ground-State Hole Dynamics of 4-(<i>N,N</i> -Dialkylamino-3-hydroxyflavone in Ionic Liquids Studied by Transient Absorption Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2021, 125, 5373-5386.	1.2	5
75	Solvation heterogeneity in ionic liquids as demonstrated by photo-chemical reactions. <i>Pure and Applied Chemistry</i> , 2020, 92, 1695-1708.	0.9	5
76	Chemical equilibrium in argon: Anomaly of density dependence of dimerization reaction in the medium density region. <i>Journal of Chemical Physics</i> , 1989, 91, 2758-2759.	1.2	4
77	Solvent Density Dependence of the Dissociation Rates of <i>N,N</i> -Methyl- <i>N</i> -Nitrosopropane Dimer in Simple and Molecular Fluids. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , 1993, 97, 29-32.	0.9	4
78	Rotational dynamics of carbon dioxide in ionic liquids. <i>Journal of Molecular Liquids</i> , 2017, 226, 43-47.	2.3	4
79	Self-Assembly and Complexation of Cellulose/Ionic Liquid at High Cellulose Concentration: Anion Dependence. <i>Crystal Growth and Design</i> , 2020, 20, 6267-6271.	1.4	4
80	Study on the Photolysis Quantum Yield of Diphenyl Disulfide by the Transient Grating Method. <i>Bulletin of the Chemical Society of Japan</i> , 1997, 70, 2657-2664.	2.0	3
81	Raman spectroscopic study on the acceptor number of supercritical methanol and ethanol. <i>Journal of Molecular Liquids</i> , 2017, 245, 11-16.	2.3	3
82	Formate Ionic Liquids Playing the Roles of Reducer and Stabilizer for the Synthesis of Noble Metal Nanoparticles. <i>Chemistry Letters</i> , 2017, 46, 1344-1346.	0.7	3
83	Role of Hydrogen Bond Interactions in CO ₂ Capture by Wet Phosphonium Formate Ionic Liquid: A Raman Spectroscopic Study. <i>ChemPhysChem</i> , 2018, 19, 1674-1682.	1.0	3
84	Structure-Property Relationship for 1-Isopropyl-3-methylimidazolium- and 1- <i>tert</i> -Butyl-3-methylimidazolium-Based Ionic Liquids: Thermal Properties, Densities, Viscosities, and Quantum Chemical Calculations. <i>Journal of Chemical & Engineering Data</i> , 2019, 64, 5857-5868.	1.0	3
85	Experimental and theoretical study on <i>p</i> -aminophenylthyl radical geminate recombination in ionic liquids; analysis using the Smoluchowski-Collins-Kimball equation. <i>Journal of Chemical Physics</i> , 2021, 154, 154504.	1.2	2
86	Factors Affecting VUV Emission Spectrum near Lyman- α from a Hydrogen Plasma Source. <i>AIP Conference Proceedings</i> , 2011, , .	0.3	1
87	Synthesis and Optical Properties of 1,2,5,10-Tetraphenylanthra[2,3- <i>b</i>]phosphole Derivatives. <i>Journal of Organic Chemistry</i> , 2022, 87, 10493-10500.	1.7	1
88	3P176 Pressure-induced reversal in the rotational direction of the bacterial flagellar motor(Molecular motors,Oral Presentations). <i>Seibutsu Butsuri</i> , 2007, 47, S247.	0.0	0
89	1P-170 Pressure-induced effects of the motility of <i>Vibrio alginolyticus</i> (The 46th Annual Meeting of the) Tj ETQq1 1 0,784314 rgBT /Over	0.0	0
90	3P087 Determination of the photoreaction product of blue light sensor BLUF protein TePixD(Protein:) Tj ETQq0 0 0,rgBT /Overlock 10 T	0.0	0

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91	Effect of Hydrated Ionic Liquid on Photocycle and Dynamics of Photoactive Yellow Protein. <i>Molecules</i> , 2021, 26, 4554.	1.7	0
92	Solvent Density Dependence of the Photolysis Quantum Yield in Supercritical Fluids.. Review of High Pressure Science and Technology/ <i>Koatsuryoku No Kagaku To Gijutsu</i> , 1998, 7, 1259-1261.	0.1	0
93	Electron-Transfer Rate in the Charge Transfer Complex in the Supercritical Fluids.. Review of High Pressure Science and Technology/ <i>Koatsuryoku No Kagaku To Gijutsu</i> , 1998, 7, 1230-1232.	0.1	0