

Yasuhiro Umebayashi

List of Publications by Year in descending order

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84
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4,977
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70
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docs citations

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times ranked

3568
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | Ionic conduction within non-stoichiometric N-Methylimidazole-Acetic Acid Pseudo-Protic ionic liquid mixtures. <i>Journal of Molecular Liquids</i> , 2022, 352, 118705. | 4.9 | 5 |
| 2 | Tools for studying ion solvation and ion pair formation in ionic liquids: isotopic substitution Raman spectroscopy. <i>Analytical Sciences</i> , 2022, 38, 1025-1031. | 1.6 | 1 |
| 3 | Thermodynamic and Structural Aspects of Solvate Ionic Liquid Formation. , 2021, , 287-300. | | 0 |
| 4 | Local Structure of Li ⁺ in Superconcentrated Aqueous LiTfSA Solutions. <i>Journal of Physical Chemistry B</i> , 2021, 125, 7477-7484. | 2.6 | 9 |
| 5 | Effect of Brønsted Acidity on Ion Conduction in Fluorinated Acetic Acid and <i>N</i> -Methylimidazole Equimolar Mixtures as Pseudo-protic Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2020, 124, 11157-11164. | 2.6 | 13 |
| 6 | Transport Properties of Ionic Liquid and Sodium Salt Mixtures for Sodium-Ion Battery Electrolytes from Molecular Dynamics Simulation with a Self-Consistent Atomic Charge Determination. <i>Journal of Physical Chemistry B</i> , 2020, 124, 7291-7305. | 2.6 | 22 |
| 7 | Speciation Analysis and Thermodynamic Criteria of Solvated Ionic Liquids: Ionic Liquids or Superconcentrated Solutions?. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 4517-4523. | 4.6 | 16 |
| 8 | Solvation Structure of Li ⁺ in Concentrated Acetonitrile and <i>N,N</i> -Dimethylformamide Solutions Studied by Neutron Diffraction with ⁶ Li/ ⁷ Li Isotopic Substitution Methods. <i>Journal of Physical Chemistry B</i> , 2020, 124, 10456-10464. | 2.6 | 9 |
| 9 | Possible Proton Conduction Mechanism in Pseudo-Protic Ionic Liquids: A Concept of Specific Proton Conduction. <i>Journal of Physical Chemistry B</i> , 2019, 123, 6244-6252. | 2.6 | 43 |
| 10 | Mixture of monoglyme-based solvent and lithium Bis(trifluoromethanesulfonyl)amide as electrolyte for lithium ion battery using silicon electrode. <i>Materials Chemistry and Physics</i> , 2019, 225, 105-110. | 4.0 | 8 |
| 11 | Solvation Structure of Li ⁺ in Methanol and 2-Propanol Solutions Studied by ATR-IR and Neutron Diffraction with ⁶ Li/ ⁷ Li Isotopic Substitution Methods. <i>Journal of Physical Chemistry B</i> , 2019, 123, 4967-4975. | 2.6 | 2 |
| 12 | Dynamic Chelate Effect on the Li ⁺ -Ion Conduction in Solvate Ionic Liquids. <i>Journal of Physical Chemistry C</i> , 2019, 123, 30228-30233. | 3.1 | 10 |
| 13 | Anion Coordination Characteristics of Ion-pair Complexes in Highly Concentrated Aqueous Lithium Bis(trifluoromethane- sulfonyl)amide Electrolytes. <i>Analytical Sciences</i> , 2019, 35, 289-294. | 1.6 | 15 |
| 14 | Direct Evidence for Li Ion Hopping Conduction in Highly Concentrated Sulfolane-Based Liquid Electrolytes. <i>Journal of Physical Chemistry B</i> , 2018, 122, 10736-10745. | 2.6 | 165 |
| 15 | Neutron Diffraction Study on Partial Pair Correlation Functions of Water at Ambient Temperature. <i>Bulletin of the Chemical Society of Japan</i> , 2018, 91, 1586-1595. | 3.2 | 9 |
| 16 | Enhanced Electrochemical Stability of Molten Li Salt Hydrate Electrolytes by the Addition of Divalent Cations. <i>Journal of Physical Chemistry C</i> , 2018, 122, 20167-20175. | 3.1 | 23 |
| 17 | Local Structure of Li ⁺ in Concentrated Ethylene Carbonate Solutions Studied by Low-Frequency Raman Scattering and Neutron Diffraction with ⁶ Li/ ⁷ Li Isotopic Substitution Methods. <i>Journal of Physical Chemistry B</i> , 2017, 121, 10979-10987. | 2.6 | 23 |
| 18 | Li ⁺ Local Structure in Li ⁺ -Tetraglyme Solvate Ionic Liquid Revealed by Neutron Total Scattering Experiments with the ⁶ Li/ ⁷ Li Isotopic Substitution Technique. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2832-2837. | 4.6 | 44 |

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|----|--|------|-----------|
| 19 | A pH Scale for the Protic Ionic Liquid Ethylammonium Nitrate. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 6266-6269. | 13.8 | 34 |
| 20 | Raman Spectroscopic Speciation Analyses and Liquid Structures by High-Energy X-ray Total Scattering and Molecular Dynamics Simulations for <i>N</i> -methylimidazolium-Based Protic Ionic Liquids. <i>Bulletin of the Chemical Society of Japan</i> , 2016, 89, 965-972. | 3.2 | 5 |
| 21 | Local structure of Li ⁺ in concentrated LiPF ₆ -dimethyl carbonate solutions. <i>Journal of Molecular Liquids</i> , 2016, 217, 17-22. | 4.9 | 24 |
| 22 | A pH Scale for the Protic Ionic Liquid Ethylammonium Nitrate. <i>Angewandte Chemie</i> , 2016, 128, 6374-6377. | 2.0 | 22 |
| 23 | Effects of non-equimolar lithium salt glyme solvate ionic liquid on the control of interfacial degradation in lithium secondary batteries. <i>RSC Advances</i> , 2016, 6, 33043-33047. | 3.6 | 18 |
| 24 | Li ⁺ Local Structure in Hydrofluoroether Diluted Li-Glyme Solvate Ionic Liquid. <i>Journal of Physical Chemistry B</i> , 2016, 120, 3378-3387. | 2.6 | 81 |
| 25 | Hydrogen bond in imidazolium based protic and aprotic ionic liquids. <i>Journal of Molecular Liquids</i> , 2016, 217, 35-42. | 4.9 | 45 |
| 26 | Li ⁺ solvation in glyme-Li salt solvate ionic liquids. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 8248-8257. | 2.8 | 222 |
| 27 | Structural and aggregate analyses of (Li salt + glyme) mixtures: the complex nature of solvate ionic liquids. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 22321-22335. | 2.8 | 78 |
| 28 | Microscopic Solvation Structure of Glucose in 1-Ethyl-3-methylimidazolium Methylphosphonate Ionic Liquid. <i>Journal of Physical Chemistry B</i> , 2015, 119, 6262-6270. | 2.6 | 9 |
| 29 | Structural modification by adding Li cations into Mg/Cs-TFSA molten salt facilitating Mg electrodeposition. <i>RSC Advances</i> , 2015, 5, 3063-3069. | 3.6 | 3 |
| 30 | Structures of [Li(glyme)] ⁺ complexes and their interactions with anions in equimolar mixtures of glymes and Li[TFSA]: analysis by molecular dynamics simulations. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 126-129. | 2.8 | 87 |
| 31 | High-Energy X-ray Diffraction and MD Simulation Study on the Ion-Ion Interactions in 1-Ethyl-3-methylimidazolium Bis(fluorosulfonyl)amide. <i>Journal of Solution Chemistry</i> , 2014, 43, 1655-1668. | 1.2 | 11 |
| 32 | Chelate Effects in Glyme/Lithium Bis(trifluoromethanesulfonyl)amide Solvate Ionic Liquids. I. Stability of Solvate Cations and Correlation with Electrolyte Properties. <i>Journal of Physical Chemistry B</i> , 2014, 118, 5144-5153. | 2.6 | 194 |
| 33 | Solvation Structure of Poly(ethylene glycol) in Ionic Liquids Studied by High-energy X-ray Diffraction and Molecular Dynamics Simulations. <i>Macromolecules</i> , 2013, 46, 2369-2375. | 4.8 | 33 |
| 34 | Unusual Li ⁺ Ion Solvation Structure in Bis(fluorosulfonyl)amide Based Ionic Liquid. <i>Journal of Physical Chemistry C</i> , 2013, 117, 19314-19324. | 3.1 | 133 |
| 35 | A New Proton Conductive Liquid with No Ions: Pseudo-Protic Ionic Liquids. <i>Chemistry - A European Journal</i> , 2013, 19, 11522-11526. | 3.3 | 60 |
| 36 | Communication: Collective dynamics of room-temperature ionic liquids and their Li ion solutions studied by high-resolution inelastic X-ray scattering. <i>Journal of Chemical Physics</i> , 2013, 138, 151101. | 3.0 | 15 |

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|----|---|-----|-----------|
| 37 | Brønsted Basicity of Solute Butylamine in an Aprotic Ionic Liquid Investigated by Potentiometric Titration. <i>Chemistry Letters</i> , 2013, 42, 1250-1251. | 1.3 | 16 |
| 38 | Intermolecular Interactions in Li ⁺ Glyme and Li ⁺ Glyme-TFSA ⁻ Complexes: Relationship with Physicochemical Properties of [Li(glyme)][TFSA] Ionic Liquids. <i>ChemPhysChem</i> , 2013, 14, 1993-2001. | 2.1 | 79 |
| 39 | Specific Solvation of Benzyl Methacrylate in 1-Ethyl-3-methylimidazolium Bis(trifluoromethanesulfonyl)amide Ionic Liquid. <i>Analytical Sciences</i> , 2013, 29, 311-314. | 1.6 | 27 |
| 40 | Relationship between mesoscale dynamics and shear relaxation of ionic liquids with long alkyl chain. <i>Journal of Chemical Physics</i> , 2012, 137, 104511. | 3.0 | 35 |
| 41 | Physicochemical and Acid-base Properties of a Series of 2-Hydroxyethylammonium-based Protic Ionic Liquids. <i>Analytical Sciences</i> , 2012, 28, 469-474. | 1.6 | 30 |
| 42 | Structural Heterogeneity and Unique Distorted Hydrogen Bonding in Primary Ammonium Nitrate Ionic Liquids Studied by High-Energy X-ray Diffraction Experiments and MD Simulations. <i>Journal of Physical Chemistry B</i> , 2012, 116, 2801-2813. | 2.6 | 116 |
| 43 | Comprehensive Refractive Index Property for Room-Temperature Ionic Liquids. <i>Journal of Chemical & Engineering Data</i> , 2012, 57, 2211-2216. | 1.9 | 191 |
| 44 | Acid-Base Property of N-Methylimidazolium-Based Protic Ionic Liquids Depending on Anion. <i>Journal of Physical Chemistry B</i> , 2012, 116, 14146-14152. | 2.6 | 57 |
| 45 | Liquid Structure of and Li ⁺ Ion Solvation in Bis(trifluoromethanesulfonyl)amide Based Ionic Liquids Composed of 1-Ethyl-3-methylimidazolium and N-Methyl-N-propylpyrrolidinium Cations. <i>Journal of Physical Chemistry B</i> , 2011, 115, 12179-12191. | 2.6 | 102 |
| 46 | Experimental evidences for molecular origin of low-Q peak in neutron/x-ray scattering of 1-alkyl-3-methylimidazolium bis(trifluoromethanesulfonyl)amide ionic liquids. <i>Journal of Chemical Physics</i> , 2011, 135, 244502. | 3.0 | 140 |
| 47 | Thermodynamic Study of the Solvation States of Acid and Base in a Protic Ionic Liquid, Ethylammonium Nitrate, and Its Aqueous Mixtures. <i>Chemistry Letters</i> , 2010, 39, 578-579. | 1.3 | 27 |
| 48 | Structure, solvation, and acid-base property in ionic liquids. <i>Pure and Applied Chemistry</i> , 2010, 82, 1927-1941. | 1.9 | 14 |
| 49 | Studies on the translational and rotational motions of ionic liquids composed of N-methyl-N-propyl-pyrrolidinium (P13) cation and bis(trifluoromethanesulfonyl)amide and bis(fluorosulfonyl)amide anions and their binary systems including lithium salts. <i>Journal of Chemical Physics</i> , 2010, 133, 194505. | 3.0 | 129 |
| 50 | Raman Spectroscopic Studies and Ab Initio Calculations on Conformational Isomerism of 1-Butyl-3-methylimidazolium Bis-(trifluoromethanesulfonyl)amide Solvated to a Lithium Ion in Ionic Liquids: Effects of the Second Solvation Sphere of the Lithium Ion. <i>Journal of Physical Chemistry B</i> , 2010, 114, 6513-6521. | 2.6 | 107 |
| 51 | Dependence of the Conformational Isomerism in 1-N-Butyl-3-methylimidazolium Ionic Liquids on the Nature of the Halide Anion. <i>Journal of Physical Chemistry B</i> , 2010, 114, 11715-11724. | 2.6 | 66 |
| 52 | Solvation and microscopic properties of ionic liquid/acetonitrile mixtures probed by high-pressure infrared spectroscopy. <i>Journal of Chemical Physics</i> , 2009, 131, 234502. | 3.0 | 29 |
| 53 | Structural change of ionic association in ionic liquid/water mixtures: A high-pressure infrared spectroscopic study. <i>Journal of Chemical Physics</i> , 2009, 130, 124503. | 3.0 | 43 |
| 54 | Ion-ion interaction in room temperature ionic liquid 1-ethyl-3-methylimidazolium tetrafluoroborate studied by large angle X-ray scattering experiment and molecular dynamics simulations. <i>Journal of Molecular Liquids</i> , 2009, 147, 77-82. | 4.9 | 53 |

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| 55 | Raman Spectroscopic Study, DFT Calculations and MD Simulations on the Conformational Isomerism of <i>N</i> -Alkyl- <i>N</i> -methylpyrrolidinium Bis-(trifluoromethanesulfonyl) Amide Ionic Liquids. <i>Journal of Physical Chemistry B</i> , 2009, 113, 4338-4346. | 2.6 | 56 |
| 56 | Relationships between center atom species (N, P) and ionic conductivity, viscosity, density, self-diffusion coefficient of quaternary cation room-temperature ionic liquids. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 3509. | 2.8 | 80 |
| 57 | Effect of Methylation at the C2 Position of Imidazolium on the Structure of Ionic Liquids Revealed by Large Angle X-ray Scattering Experiments and MD Simulations. <i>Chemistry Letters</i> , 2009, 38, 340-341. | 1.3 | 42 |
| 58 | Liquid structure of <i>N</i> -butyl- <i>N</i> -methylpyrrolidinium bis-(trifluoromethanesulfonyl) amide ionic liquid studied by large angle X-ray scattering and molecular dynamics simulations. <i>Journal of Molecular Liquids</i> , 2008, 143, 2-7. | 4.9 | 54 |
| 59 | Liquid structure and conformation of a low-viscosity ionic liquid, <i>N</i> -methyl- <i>N</i> -propyl-pyrrolidinium bis(fluorosulfonyl) imide studied by high-energy X-ray scattering. <i>Journal of Molecular Liquids</i> , 2008, 143, 64-69. | 4.9 | 75 |
| 60 | Liquid Structure of Room-Temperature Ionic Liquid, 1-Ethyl-3-methylimidazolium Bis-(trifluoromethanesulfonyl) Imide. <i>Journal of Physical Chemistry B</i> , 2008, 112, 4329-4336. | 2.6 | 159 |
| 61 | Potential Energy Landscape of Bis(fluorosulfonyl)amide. <i>Journal of Physical Chemistry B</i> , 2008, 112, 9449-9455. | 2.6 | 81 |
| 62 | Raman Spectroscopic Study on Alkaline Metal Ion Solvation in 1-Butyl-3-methylimidazolium Bis(trifluoromethanesulfonyl)amide Ionic Liquid. <i>Analytical Sciences</i> , 2008, 24, 1297-1304. | 1.6 | 38 |
| 63 | Acidity and Basicity of Aqueous Mixtures of a Protic Ionic Liquid, Ethylammonium Nitrate. <i>Analytical Sciences</i> , 2008, 24, 1347-1349. | 1.6 | 54 |
| 64 | Solvation Structures of Some Transition Metal(II) Ions in a Room-Temperature Ionic Liquid, 1-Ethyl-3-methylimidazolium Bis(trifluoromethanesulfonyl)amide. <i>Analytical Sciences</i> , 2008, 24, 1377-1380. | 1.6 | 76 |
| 65 | Liquid Structure and the Ion-Ion Interactions of Ethylammonium Nitrate Ionic Liquid Studied by Large Angle X-Ray Scattering and Molecular Dynamics Simulations. <i>Journal of Computer Chemistry Japan</i> , 2008, 7, 125-134. | 0.1 | 97 |
| 66 | Acid-Base Property of Ethylammonium Nitrate Ionic Liquid Directly Obtained Using Ion-selective Field Effect Transistor Electrode. <i>Chemistry Letters</i> , 2007, 36, 684-685. | 1.3 | 61 |
| 67 | Anion Conformation of Low-Viscosity Room-Temperature Ionic Liquid 1-Ethyl-3-methylimidazolium Bis(fluorosulfonyl) Imide. <i>Journal of Physical Chemistry B</i> , 2007, 111, 12829-12833. | 2.6 | 127 |
| 68 | Lithium Ion Solvation in Room-Temperature Ionic Liquids Involving Bis(trifluoromethanesulfonyl) Imide Anion Studied by Raman Spectroscopy and DFT Calculations. <i>Journal of Physical Chemistry B</i> , 2007, 111, 13028-13032. | 2.6 | 321 |
| 69 | Solvation Structure of Li ⁺ in Concentrated LiPF ₆ /Propylene Carbonate Solutions. <i>Journal of Physical Chemistry B</i> , 2007, 111, 6104-6109. | 2.6 | 131 |
| 70 | Solvation structure of magnesium, zinc, and alkaline earth metal ions in <i>N,N</i> -dimethylformamide, <i>N,N</i> -dimethylacetamide, and their mixtures studied by means of Raman spectroscopy and DFT calculations ³⁹ Ionic size and electronic effects on steric congestion. <i>Journal of Raman Spectroscopy</i> , 2007, 38, 417-426. | 2.5 | 33 |
| 71 | Conformational Equilibrium of Bis(trifluoromethanesulfonyl) Imide Anion of a Room-Temperature Ionic Liquid: A Raman Spectroscopic Study and DFT Calculations. <i>Journal of Physical Chemistry B</i> , 2006, 110, 8179-8183. | 2.6 | 333 |
| 72 | Solvent conformation and ion solvation: From molecular to ionic liquids. <i>Pure and Applied Chemistry</i> , 2006, 78, 1595-1609. | 1.9 | 13 |

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|----|---|-----|-----------|
| 73 | Evidence of Conformational Equilibrium of 1-Ethyl-3-methylimidazolium in Its Ionic Liquid Salts: A Raman Spectroscopic Study and Quantum Chemical Calculations. <i>Journal of Physical Chemistry A</i> , 2005, 109, 8976-8982. | 2.5 | 199 |
| 74 | Thermodynamic Aspects of Metal Ion Complexation in the Structured Solvent, N-Methylformamide. <i>Journal of Solution Chemistry</i> , 2005, 34, 739-753. | 1.2 | 14 |
| 75 | Solvation Structure and Complexation of the Manganese(II) Ion in N,N-Dimethylpropionamide and N,N,N',N'-Tetramethylurea Studied by Means of Titration Calorimetry and Raman Spectroscopy. <i>Journal of Solution Chemistry</i> , 2005, 34, 1429-1443. | 1.2 | 7 |
| 76 | Conformation of Solvent N,N-Dimethylpropionamide in the Coordination Sphere of the Zinc(II) Ion Studied by Raman Spectroscopy and DFT Calculations. <i>Journal of Physical Chemistry A</i> , 2005, 109, 4862-4868. | 2.5 | 18 |
| 77 | Thermodynamics and Fluorescence Spectra of 1,10-Phenanthroline in Micelles of Poly (Ethylene Terephthalate) / Overlooked | 1.2 | 4 |
| 78 | Conformational equilibria of solvent N,N-dimethylpropionamide in the bulk and in the coordination sphere of the manganese(ii) ion Electronic supplementary information (ESI) available: non-planar staggered and planar cis Gaussian results. See http://www.rsc.org/suppdata/cp/b3/b302143b/ . <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 2552. | 2.8 | 24 |
| 79 | Solvation structure of lanthanide(III) ions in solvent mixtures of N,N-dimethylformamide and N,N-dimethylacetamide studied by titration Raman spectroscopy Electronic supplementary information (ESI) available: Crystallographic data (single crystal, [Gd(DMF)4(DMA)4](ClO4)3), (CCDC reference) http://www.rsc.org/suppdata/cp/b3/b302143b/ . <i>Physical Chemistry Chemical Physics</i> , 2002, 4, 5599-5605. | 1.2 | 4 |
| 80 | Title is missing!. <i>Journal of Solution Chemistry</i> , 2002, 31, 931-946. | 1.2 | 5 |
| 81 | Individual solvation number of first-row transition metal(II) ions in solvent mixtures of N,N-dimethylformamide and N,N-dimethylacetamide Solvation steric effect. <i>Physical Chemistry Chemical Physics</i> , 2001, 3, 5475-5481. | 2.8 | 38 |
| 82 | Individual Solvation Numbers around the Nickel (II) Ion in an N,N-Dimethylformamide and N,N-Dimethylacetamide Mixture Determined by Raman Spectrophotometry. <i>Analytical Sciences</i> , 2001, 17, 323-326. | 1.6 | 31 |
| 83 | Thermodynamics of [Co(NCS)4]2- at Poly(ethylene Oxide) and Octylphenyl Moieties in Micelles of Nonionic Surfactants. <i>Journal of Colloid and Interface Science</i> , 2001, 237, 167-173. | 9.4 | 14 |
| 84 | Spectrophotometric study of thiocyanato complexation of cobalt(II) and nickel(II) ions in micellar solutions of a nonionic surfactant triton X-100. <i>Journal of Solution Chemistry</i> , 1996, 25, 731-746. | 1.2 | 12 |