

# **Curriculum Vitae and Complete Publication List**

**Professor Biman Bagchi**



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## Curriculum Vitae of Prof. Biman Bagchi

**Name:** Biman Bagchi

**Designation:** Amrut Mody Professor

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**Date of birth:** January 1, 1954

### Education:

- **B.Sc.**, 1974, Presidency College, Calcutta University, Chemistry (Hons, First Class).
- **M.Sc.**, 1976, Calcutta University, Physical Chemistry (First Class).
- **Ph.D.**, 1981, Brown University, Providence, R.I., U.S.A. (Adviser: Prof. Julian H. Gibbs).

### Post-doctoral research experience:

- Research Associate (1981-1983) at the James Franck Institute, University of Chicago, U.S.A. (With Prof. David W. Oxtoby).
- Research Associate (1983-1984) at the Institute for Physical Science and Technology, University of Maryland, U.S.A (with Prof. R.W. Zwanzig).

### Work Experience:

- Lecturer at Solid State and Structural Chemistry Unit (SSCU), Indian Institute of Science (August, 1984 - March, 1987).
- Assistant Professor (March 1987 - June 1991).
- Associate Professor (June 1991 - December 1995).

- Professor (1995 - ). And Chairman (2000 - 2004).
- Amrut Mody Chair (2004 - )

#### **Scientific Visits (after returning to India):**

- Visiting Scientist at the University of Maryland (summer, 1986). (Dr. R. Zwanzig)
- Visiting Scientist at the University of Chicago (Summer, 1987 and 1989).
- Visitor at University of Umea, UMEA, Sweden (summer, 1988).
- Visiting Scientist at N.I.H., Bethesda, U.S.A. (1991 -1992). (With R. Zwanzig).
- Visiting Professor at the Univ. Piere et Marie Curie, Paris 5, France (June, 1992).
- Visiting Scientist, IMS, Okazaki, Japan (June - July, 1994). (Host: Prof. K. Yoshihara and Prof. Iwao Ohmine).
- Visiting Professor, University of Wisconsin, Madison, Wisconsin, U.S.A. (1998-1999). (Prof. J.L. Skinner).
- Visiting Professor, University of Texas at Austin, U.S.A. (summer, 1999). (P.F. Barbara and Peter Rossky).
- Invited Professor, Pierre et Marie Curie, Paris 5, France (September, 2001). (Professor Pierre Turq).
- Distinguished Visitor at LMS, Chemistry, California Institute of Technology (Pasadena, USA) (April-May, 2002).
- Invited Professor at Ecole Normal Superior, Paris, France (June, 2002).
- Invited Professor at Nagoya University, Japan (Apr-May, 2004).

#### **Awards & Distinctions:**

- National Merit Certificate (1971)
- First Degree State Scholarship (1971 - 1974).
- Young Associate of the Indian Academy of Science (1985 -1989).
- Young Scientist Medal of the Indian National Science Academy (1986).
- Homi Bhabha Fellow (1989 - 1991).
- A.K. Bose Memorial Award (1990) from Indian National Science Academy for best research paper in physical sciences (published on work done in India).
- **S.S. Bhatnagar Prize in Chemical Sciences (1991).**
- **Elected Fellow of the Indian Academy of Sciences (1991).**
- **Elected Fellow of the Indian National Science Academy, New Delhi, (1994).**
- **Elected Fellow of the National Academy of Sciences, Allahabad, (1995).**
- Jagdish Shankar Memorial Lecture award from Indian National Science Academy (1997).
- **Seventh G.D. Birla Award for Science and Technology, 1997.**
- Platinum Jubilee Lecture of Indian Science Congress, January 1997.
- **Third World Academy of Sciences (TWAS) in Chemistry, 1998.**
- BC Laha Memorial Lecture of the Indian Assoc. Cultivation of Science (Calcutta, December, 2001).
- Alumni Excellence Award in Research, IISc. (2002).
- Silver Medal of Chem. Res. Soc. of India (2002).
- Goyal prize in Chemistry (2003).
- Amrut Mody Chair Professorship (2004-2007).
- Fellowship of the Third World Academy of Sciences, Trieste, Italy (2004).

- J. C. Bose National Fellowship (2006)

#### **Editorial Board:**

- Editorial Advisory Board, **J. Phys. Chem. (ACS)** (2001- ).
- Editorial Advisory Board, **Chemical Physics Letters**. (2002 - 2004).
- Editorial Board, **Journal of Molecular Liquids** (1993 - ) (An Elsevier publication)
- Editorial Advisory Board, **PhysChemComm** (RSC) (2000- )
- International Advisory Board, **Adv. Chem. Phys.** (Wiley) (2000- )
- Editorial Advisory Board, **Theoretical Chemistry Accts.** (2002 - ).
- Editorial Board, **Indian Journal of Chemistry** (1995 -2001).
- Editorial Board, **Proceedings of Indian Academy of Sciences** (1989-1993).

#### **Other Activity:**

- Honorary Professor of the Jawaharlal Nehru Center for Advanced Scientific Research, Bangalore.
- Honorary Professor of S.N. Bose National Center of Basic Sciences, Calcutta.
- Honorary Professor of National Institute of Advanced studies, Bangalore.
- **Member of the Research Council, National Chemical Laboratory (NCL), Pune** (2001-2004).

#### **Named Lectures:**

- A.K. Bose Memorial Lecture (1990) at Indian National Science Academy, New Delhi.
- Jagdish Shankar Memorial Lecture at Indian National Science Academy, New Delhi (1997).
- Platinum Jubilee Lecture of Indian Science Congress, Hyderabad, January 1997.
- B.C. Laha Memorial Lecture of the Indian Assoc. Cultivation of Science (Calcutta, December, 2001).
- Raman-Mizushima Lecture of JSPS & DST (2006).

#### **Major Scientific fields of interest:**

- Time dependent statistical mechanics, stochastic processes.
- Dynamics of chemical and biological Reactions.
- Phase transitions, dynamics in glassy liquids.
- Liquid crystals.
- Biological water.
- Protein Folding, biopolymers, DNA intercalation.
- Polar solvation dynamics, electron transfer reactions.
- Vibrational and rotational relaxation in liquids.
- Electrochemistry, transport properties, classical laws, polyelectrolyte.
- Conjugated polymers, polymer reactions, Forster energy transfer.

## **Publications:**

- In reputed journals: **288**
- Presentations in Conferences/Meetings/Seminars: **85**
- Major reviews: **14** (Adv. Chem. Phys.: 4, JPC (Feature): 4, Chem. Rev.: 2, Int. Rev. Phys. Chem.: 2, Annu. Rep. Prog. Chem., Sect. C.: 1, Ann. Rev. Phys. Chem.: 1).
- Edited: Special Issue of JPC: 1, Special Issues of J. Mol. Liq.: 2, A book entitled “Ultrafast Processes in Chemistry” (World Scientific, 1991).
- Appeared in the list of top 1% most cited chemists of the world.

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## Publication List of Professor Biman Bagchi

**Total number of papers: 288 (in journals only)**

**Average citation per year: About 400**

1. *Decoupling phenomena in supercooled liquids: Signatures in the energy landscape*, D. Chakrabarti and **B. Bagchi**, *Phys. Rev. Lett.* **96**, 187801 (2006).
2. *Energy landscape view of phase transitions and slow dynamics in thermotropic liquid crystals*, D. Chakrabarti and **B. Bagchi**, *Proc. Natl. Acad. Sci. USA* **103**, 7217 (2006).
3. *Power law relaxation and glassy dynamics in Lebwohl-Lasher model near the isotropic-nematic phase transition*, S. Chakrabarty, D. Chakrabarti and **B. Bagchi**, *Phys. Rev. E* **73**, 061706 (2006).
4. *Ion dynamics in compacted clays: Derivation of a two-state diffusion-reaction scheme from the lattice Fokker-Planck equation*, B. Rotenberg, J.-F. Dufreche, **B. Bagchi**, E. Giffaut, J.-P. Hansen and P. Turq, *J. Chem. Phys.* **124**, 154701 (2006).
5. *Anomalous orientation dependent effective pair interaction among histidine and other amino acid residues in metalloproteins: Breakdown of the hydrophathy scale index*, A. Mukherjee and **B. Bagchi**, *Biochemistry* **45**, 5129 (2006).
6. *Complete breakdown of the Debye model of rotational relaxation near the isotropic-nematic phase boundary: Effect of intermolecular correlations in orientational dynamics*, P. P. Jose, D. Chakrabarti and **B. Bagchi**, *Phys. Rev. E* **73**, 031705 (2006).
7. *Solvation dynamics in biological systems and organized assemblies*, P. Sen, S. Pal, K. Bhattacharyya and **B. Bagchi**, *J. Chin. Chem. Soc. (Taipei)* **53**, 169 (2006).
8. *Förster energy transfer in thin films of conjugated polymers and in solution*, A. Mukherjee and **B. Bagchi**, *J. Chin. Chem. Soc. (Taipei)* **53**, 153 (2006).
9. *Fluorescence resonance energy transfer (FRET) in chemistry and biology: Non-Förster distance dependence of the FRET rate*, S. Saini, H. Singh and **B. Bagchi**, *J. Chem. Sci.* **118**, 23 (2006).
10. *Anisotropic and sub-diffusive water motion at the surface of a DNA and of an anionic micelle CsPFO*, S. Pal, P. K. Maiti and **B. Bagchi**, *J. Phys. Cond. Matt.* **17**, S4317 (2005).
11. *Universal power law in the orientational relaxation in thermotropic liquid crystals*, D. Chakrabarti, P. P. Jose, S. Chakrabarty and **B. Bagchi**, *Phys. Rev. Lett.* **95**, 197801 (2005).

12. *Secondary structure sensitivity of Hydrogen bond lifetime dynamics in the protein hydration layer*, S. Bandyopadhyay, S. Chakraborty and **B. Bagchi**, *J. Am. Chem. Soc.* **127**, 16660 (2005).
13. *Non-Förster Distance and Orientation Dependence of Energy Transfer and Applications of Fluorescence Resonance Energy Transfer to Polymers and Nanoparticles: How Accurate is the Spectroscopic Ruler with  $1/R^6$  Rule?* H. Singh and **B. Bagchi**, *Curr. Sci.* **89**, 1710 (2005).
14. *Water Dynamics in the Hydration Layer around Proteins and Micelles*, **B. Bagchi**, *Chem. Rev. (Review)* **105**, 3197 (2005).
15. *Bridging the gap between the mode coupling and the random first order transition theories of structural relaxation in liquids*, S. M. Bhattacharyya, **B. Bagchi**, and P. G. Wolynes, *Phys. Rev. E* **72**, 031509 (2005).
16. *Water dynamics at the surface of proteins and micelles: Understanding the fast and the slow components*, S. Pal, S. Chakraborty, S. M. Bhattacharyya, S. Bandyopadhyay, S. Balasubramanian and **B. Bagchi**, Proceedings of IXth International Symposium On the Properties Of Water, CRC Press, 2005.
17. *Rotational friction on globular proteins combining dielectric and hydrodynamic effects*, A. Mukherjee and **B. Bagchi**, *Chem. Phys. Lett.* **404**, 409 (2005).
18. *Orientation-dependent potential of mean force for protein folding*, A. Mukherjee, P. Bhimalapuram and **B. Bagchi**, *J. Chem. Phys.* **123**, 014901 (2005).
19. *Hydration Layer of a Cationic Micelle, C<sub>10</sub>TAB: Structure, Rigidity, Slow Reorientation, Hydrogen Bond Lifetime, and Solvation Dynamics*, S. Pal, **B. Bagchi** and S. Balasubramanian, *J. Phys. Chem. B*, **109**, 12879 (2005).
20. *Non-monotonic composition dependence of vibrational phase relaxation rate in binary mixtures*, S. Roychowdhury and **B. Bagchi**, *J. Chem. Phys.* **122**, 144507 (2005).
21. *Anomalous glassy relaxation near the isotropic-nematic phase transition*, P. P. Jose, D. Chakrabarti and **B. Bagchi**, *Phys. Rev. E*, **71**, 030701(R) (2005).
22. *Sensitivity of polar solvation dynamics to the secondary structures of aqueous proteins and the role of surface exposure of the probe*, S. Bandyopadhyay, S. Chakraborty, S. Balasubramanian and **B. Bagchi**, *J. Am. Chem. Soc.* **127**, 4071 (2005).
23. *Frequency dependent heat capacity within a kinetic model of glassy dynamics*, D. Chakrabarti and **B. Bagchi**, *J. Chem. Phys.* **122**, 14501 (2005).
24. *Atomistic simulation study of the coupled motion of amino acid residues and water molecules around protein HP-36: Fluctuations at and around the active*

- sites, S. Bandyopadhyay, S. Chakraborty, S. Balasubramanian, S. Pal, and **B. Bagchi**, *J. Phys. Chem. B.* **108**, 12608 (2004).
25. *Self-consistent mode-coupling theory for the viscosity of rodlike polyelectrolyte solutions*, K. Miyazaki, **B. Bagchi**, and A. Yethiraj, *J. Chem. Phys.* **121**, 8120 (2004).
  26. *Anomalous viscoelasticity near the isotropic-nematic phase transition in liquid crystals*, P. P. Jose and **B. Bagchi**, *J. Chem. Phys.* **121**, 6978 (2004).
  27. *Distance and orientation dependence of excitation transfer rates in conjugated systems: Beyond the Forster theory*, K. F. Wong, **B. Bagchi**, and P. J. Rossky, *J. Phys. Chem. A*, **108**, 5752 (2004).
  28. *Nonmonotonic temperature dependence of heat capacity through the glass transition within a kinetic model*, D. Chakrabarti and **B. Bagchi**, *J. Chem. Phys.* **120**, 11678 (2004).
  29. *In search of temporal power laws in the orientational relaxation near isotropic-nematic phase transition in model nematogens*, P. P. Jose and **B. Bagchi**, *J. Chem. Phys.* **120**, 11256 (2004).
  30. *Density and energy relaxation in open one-dimensional system*, Prasanth P. Jose and **B. Bagchi**, *J. Chem. Phys.* **120**, 8327 (2004).
  31. *Dynamics of water at surface of complex systems: Study of aqueous micelles and proteins*, S. Pal, S. Bandyopadhyay, S. Balasubramanian, and **B. Bagchi**, *Femtochemistry and Femtobiology (Proceedings of the VIth International Conference on Femtochemistry (2003))*, 213 (2004).
  32. *Anomalous dielectric relaxation of water molecules at the surface of an aqueous micelle*, S. Pal, S. Balasubramanian, and **B. Bagchi**, *J. Chem. Phys.* **120**, 1912 (2004).
  33. *Contact pair dynamics during folding of two small proteins: chicken villin head piece and the Alzheimer protein  $\beta$ -amyloid*, A. Mukherjee and **B. Bagchi**, *J. Chem. Phys.* **120**, 1602 (2004).
  34. *Dynamics of water at the interface of a small protein, enterotoxin*, S. Balasubramanian, S. Bandyopadhyay, S. Pal, and **B. Bagchi**, *Curr. Sci.* **85**, 1571 (2003).
  35. *Probing folding free energy landscape of small proteins through minimalistic models: Folding of HP-36 and  $\beta$ -Amyloid*, A. Mukherjee and **B. Bagchi**, *Proc. Indian Acad. Sci. (Chem. Sci.)* **115**, 621 (2003).
  36. *Identity, energy, environment, and dynamics of interfacial water molecules in a micellar solution*, S. Balasubramanian, S. Pal, and **B. Bagchi**, *J. Indian Inst. Sci.* **83**, 27 (2003, Jan.-Apr.).



37. *Water solvation dynamics in the bulk and in the hydration layer of proteins and self-assemblies*, **B. Bagchi**, *Annu. Rep. Prog. Chem., Sect. C.* **99**, 127 (2003).
38. *Identity, energy, and environment of interfacial water molecules in a micellar solution*, S. Pal, S. Balasubramanian, and **B. Bagchi**, *J. Phys. Chem. B* **107**, 5194 (2003).
39. *Dynamics of bound and free water in an aqueous micellar solution: Analysis of the lifetime and vibrational frequencies of hydrogen bonds at a complex interface*, S. Pal, S. Balasubramanian, and **B. Bagchi**, *Phys. Rev. E* **67**, 61502 (2003).
40. *Study of pair contact formation among hydrophobic residues in a model HP-36 protein: Relationship between contact order parameter and rate of folding and collapse*, G. Srinivas and **B. Bagchi**, *J. Phys. Chem. A*, **107**, 11768 (2003).
41. *Simulation and theory of vibrational phase relaxation in the critical and supercritical nitrogen: Origin of observed anomalies*, S. Roychowdhury and **B. Bagchi**, *J. Chem. Phys.* **119**, 3278 (2003).
42. *Fluorescence resonance energy transfer dynamics during protein folding: Evidence of multistage folding kinetics*, A. Mukherjee and **B. Bagchi**, *Curr. Sci.* **85**, 68 (2003).
43. *Pair dynamics in a glass-forming binary mixture: Simulations and theory*, R. K. Murarka and **B. Bagchi**, *Phys. Rev. E* **67**, 41501 (2003).
44. *Diffusion and viscosity in a supercooled polydisperse system*, R. K. Murarka and **B. Bagchi**, *Phys. Rev. E* **67**, 51504 (2003).
45. *Waiting time distribution and nonexponential relaxation in single molecule spectroscopic studies: Realization of entropic bottleneck in a simple model*, D. Chakrabarti and **B. Bagchi**, *J. Chem. Phys.* **118**, 7965 (2003).
46. *Origin of the sub-diffusive behavior and crossover from sub-diffusive to super-diffusive dynamics near a biological surface*, A. Mukherjee and **B. Bagchi**, *PhysChemComm.* **6**, 28 (2003).
47. *Correlation between rate of folding, energy landscape and topology in the folding of a model protein HP-36*, A. Mukherjee and **B. Bagchi**, *J. Chem. Phys.* **118**, 4733 (2003).
48. *Vibrational phase relaxation along the critical isochore of nitrogen: The role of local density fluctuations in the rate enhancement*, S. Roychowdhury and **B. Bagchi**, *Phys. Rev. Lett.* **90**, 75701 (2003).
49. *Evidence for bound and free water species in the hydration shell of an aqueous micelle*, S. Balasubramanian, S. Pal, and **B. Bagchi**, *Curr. Sci.* **84**, 428 (2003).

50. *Study of the dynamics of protein folding through minimalistic models*, G. Srinivas and **B. Bagchi**, *Theo. Chem. Acc.* **109**, 8 (2003).
51. *Dynamics of water in the hydration layer of a self-assembled system and in biomolecules*, **B. Bagchi**, *Proc. Ind. Natl. Sci. Acad. (Phys. Sci.)* **69A**, 15 (2003).
52. *Biological water: Femtosecond dynamics of Macromolecular hydration*, S. K. Pal, J. Peon, **B. Bagchi**, and A.H. Zewail, *J. Phys. Chem. B* **106**, 12376 (2002).
53. *Diffusion of small light particles in a solvent of large massive molecules*, R. K. Murarka, S. Bhattacharyya, and **B. Bagchi**, *J. Chem. Phys.* **117**, 10730 (2002).
54. *Anisotropic local stress and particle hopping in a deeply supercooled liquid*, S. Bhattacharyya and **B. Bagchi**, *Phys. Rev. Lett.* **89**, 25504 (2002).
55. *Hydrogen-bond dynamics near a micellar surface: Origin of the universal slow relaxation at complex aqueous interfaces*, S. Balasubramanian, S. Pal and **B. Bagchi**, *Phys. Rev. Lett.* **89**, 115505 (2002).
56. *Ionic Self-Diffusion in Concentrated Aqueous Electrolyte Solutions*, J. F. Dufreche, O. Bernard, P. Turq, A. Mukherjee, and **B. Bagchi**, *Phys. Rev. Lett.* **88**, 95902 (2002).
57. *Temperature dependence of water dynamics at an aqueous micellar surface: Atomistic molecular dynamics simulation studies of a complex system*, S. Pal, S. Balasubramanian, and **B. Bagchi**, *J. Chem. Phys.* **117**, 2852 (2002).
58. *Local composition fluctuations in strongly Nonideal binary mixtures*, R. K. Murarka and **B. Bagchi**, *J. Chem. Phys.* **117**, 1155 (2002).
59. *Foldability and the funnel of HP-36 protein sequence: Use of hydrophathy scale in protein folding*, G. Srinivas and **B. Bagchi**, *J. Chem. Phys.* **116**, 8579 (2002).
60. *Time-dependent survival probability in diffusion-controlled reactions in a polymer chain: Beyond the Wilemski-Fixman theory*, G. Srinivas, K. L. Sebastian, and **B. Bagchi**, *J. Chem. Phys.* **116**, 7276 (2002).
61. *Comparison of the ultrafast to slow time scale dynamics of three liquid crystals in the isotropic phase*, S. D. Gottke, H. Cang, **B. Bagchi**, and M. D. Fayer, *J. Chem. Phys.* **116**, 6339 (2002).
62. *Liquid crystal dynamics in the isotropic phase*, S. D. Gottke, D. D. Brace, H. Cang, **B. Bagchi**, and M. D. Fayer, *J. Chem. Phys.* **116**, 360 (2002).
63. *Dynamics of water molecules at the surface of an aqueous micelle: Atomistic molecular dynamics simulation study of a complex system*, S. Balasubramanian, S. Pal, and **B. Bagchi**, *Curr. Sci.* **82**, 845 (2002).

64. *Pressure and temperature dependence of viscosity and diffusion coefficients of a glassy binary mixture*, S. Bhattacharyya, A. Mukherjee, and **B. Bagchi**, *J. Chem. Phys.*, **116**, 4577 (2002).
65. *Intermittency, current flows, and short time diffusion in interacting finite sized one-dimensional fluids*, S. Pal, G. Srinivas, S. Bhattacharyya, and **B. Bagchi**, *J. Chem. Phys.* **116**, 5941 (2002).
66. *Formation of nanoclusters under radiation pressure in solution: A Brownian dynamics simulation study*, P. P. Jose and **B. Bagchi**, *J. Chem. Phys.* **116**, 2556 (2002).
67. *Correlated orientational and translational motions in supercooled liquids*, S. Bhattacharyya, A. Mukherjee, and **Biman Bagchi**, *J. Chem. Phys.* **117**, 2741 (2002).
68. *Anisotropic diffusion of spheroids in liquids: Slow orientational relaxation of the oblates*, R. Vasanthi, S. Bhattacharyya, and **B. Bagchi**, *J. Chem. Phys.* **116**, 1092 (2002).
69. *Detection of collapsed and ordered polymer structures by fluorescence resonance energy transfer in stiff homopolymers: Bimodality in the reaction efficiency distribution*, G. Srinivas and **B. Bagchi**, *J. Chem. Phys.*, **116**, 837 (2002).
70. *Folding and unfolding of chicken villin headpiece: Energy landscape of a single-domain model protein*, G. Srinivas and **B. Bagchi**, *Curr. Sci.*, **82** 179 (2002).
71. *Slow orientational dynamics of water at a micellar surface*, S. Balasubramanian and **B. Bagchi**, *J. Phys. Chem. B* **106**, 3668 (2002).
72. *Energy transfer efficiency distributions in polymers in solution during folding and unfolding*, G. Srinivas and **B. Bagchi**, *PhysChemComm* **8**, 59 (2002).
73. *Relaxation in binary mixtures: Non-Ideality, heterogeneity and re-entrance*, A. Mukherjee, G. Srinivas, S. Bhattacharyya, and **B. Bagchi**, *Proc. Ind. Acad. Sci. (Chem. Sci.)* **113**, 393, (2001).
74. *Anisotropic diffusion of tagged spheres near the isotropic- nematic phase transition*, R. Vasanthi, S. Ravichandran, and **B. Bagchi**, *J. Chem. Phys.* **115**, 10022 (2001).
75. *Anomalous orientational relaxation of solute probes in binary mixtures*, S. Bhattacharyya and **B. Bagchi**, *J. Chem. Phys.* **115**, 9061 (2001).
76. *Chemical reaction dynamics and relaxation phenomena in one component liquids, binary mixtures and electrolyte solutions*, **B. Bagchi**, *Curr. Sci.* **81**, 1054 (2001).
77. *Nonideality in binary mixtures: Correlations between excess volume, excess viscosity, and diffusion coefficients*, A. Mukherjee and **B. Bagchi**, *J. Phys. Chem. B* **105**, 9581 (2001).

78. *Effects of vibrational energy relaxation and reverse reaction on electron transfer kinetics and fluorescence line shapes in solution*, R. A. Denny, **B. Bagchi**, and P. F. Barbara, *J. Chem. Phys.* **115**, 6058 (2001).
79. *Effect of orientational motion of mobile chromophores on the dynamics of Forster energy transfer in polymers*, G. Srinivas and **B. Bagchi**, *J. Phys. Chem. B* **105**, 9370 (2001).
80. *Slow solvation dynamics near an aqueous micellar surface*, S. Balasubramanian and **B. Bagchi**, *J. Phys. Chem. B* **105**, 12529 (2001).
81. *Heterogeneous relaxation in supercooled liquids: A density functional theory analysis*, R. K. Murarka and **B. Bagchi**, *J. Chem. Phys.* **115**, 5513 (2001).
82. *The Enskog theory for classical vibrational energy relaxation in fluids with continuous potentials*, **B. Bagchi**, G. Srinivas, and K. Miyazaki, *J. Chem. Phys.* **115**, 4195 (2001).
83. *The extended Enskog operator for simple fluids with continuous potentials: single particle and collective properties*, K. Miyazaki, I. M.de Schepper, and B. Bagchi, *Physica A*, **298**, 101 (2001).
84. *Relation between orientational correlation time and the self-diffusion coefficient of tagged probes in viscous liquids: A density functional theory analysis*, **B. Bagchi**, *J. Chem. Phys.* **115**, 2207 (2001).
85. *Structural and electronic characterization of chemical and conformational defects in conjugated polymers*, K. F. Wong, M. S. Skaf, C. -Y. Yang, P. J. Rossky, **B. Bagchi**, D. Hu, J. Yu, and P.F. Barbara, *J. Phys. Chem. B* **105**, 6103 (2001).
86. *Reentrant behavior of relaxation time with viscosity at varying composition in binary mixtures*, A. Mukherjee, G. Srinivas, and **B. Bagchi**, *Phys. Rev. Lett.* **86**, 5926 (2001).
87. *Nonexponentiality of time dependent survival probability and the fractional viscosity dependence of the rate in diffusion controlled reactions in a polymer chain*, G. Srinivas, A. Yethiraj, and **B. Bagchi**, *J. Chem. Phys.* **114**, 9170 (2001).
88. *Needlelike motion of prolate ellipsoids in the sea of spheres*, R. Vasanthi, S. Ravichandran, and **B. Bagchi**, *J. Chem. Phys.* **114**, 7989 (2001).
89. *FRET by FET and dynamics of polymer folding*, G. Srinivas, A. Yethiraj, and **B. Bagchi**, *J. Phys. Chem. B* **105**, 2475 (2001).
90. *The Enskog theory for transport coefficients of simple fluids with continuous potentials*, K. Miyazaki, G. Srinivas, and **B. Bagchi**, *J. Chem. Phys.* **114**, 6276 (2001).

91. *Nonideality in the composition dependence of viscosity in binary mixtures*, G. Srinivas, A. Mukherjee, and **B. Bagchi**, *J. Chem. Phys.* **114**, 6220 (2001).
92. *Mode coupling theory approach to liquid-state dynamics*, **B. Bagchi** and S. Bhattacharyya, *Adv. Chem. Phys.* **116**, 67 (2001).
93. *Slow dynamics of constrained water in complex geometries*, K. Bhattacharyya and **B. Bagchi**, *J. Phys. Chem. A* **104**, 10603 (2000).
94. *Beyond the classical transport laws of electrochemistry: New microscopic approach to ionic conductance and viscosity*, A. Chandra and **B. Bagchi**, *J. Phys. Chem. B* **104**, 9067 (2000).
95. *Distribution of reaction times in diffusion controlled reactions in polymers*, G. Srinivas and **B. Bagchi**, *Chem. Phys. Lett.* **328**, 420 (2000).
96. *Ionic contribution to the viscosity of dilute electrolyte solutions: Towards a microscopic theory*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **113**, 3236 (2000).
97. *Dielectric relaxation and solvation dynamics of water in complex chemical and biological systems*, N. Nandi, K. Bhattacharyya, and **B. Bagchi**, *Chem. Rev.* **100**, 2013 (2000).
98. *Power law mass dependence of diffusion: A mode coupling theory analysis*, S. Bhattacharyya and **B. Bagchi**, *Phys. Rev. E* **61**, 3850 (2000).
99. *Collapse of stiff conjugated polymers with chemical defects into ordered, cylindrical conformations*, De Hong, Ji Hu, **B. Bagchi**, P. Rossky and P. Barbara, *Nature* **405**, 1030 (2000).
100. *Cage dynamics in the third-order off-resonant response of liquid molecules: A theoretical realization*, K. Okumura, **B. Bagchi**, and Y. Tanimura, *Bull. Chem. Soc. Jpn.* **73**, 873 (2000).
101. *Understanding the anomalous  $1/t^3$  time dependence of velocity correlation function in one dimensional Lennard-Jones systems*, G. Srinivas and **B. Bagchi**, *J. Chem. Phys.* **112**, 7557 (2000).
102. *Computer simulation and mode-coupling theory analysis of time-dependent diffusion in two dimensional Lennard-Jones fluids*, S. Bhattacharyya, G. Srinivas and **B. Bagchi**, *Phys. Lett. A*, **266**, 394 (2000).
103. *Frequency dependence of ionic conductivity of electrolyte solutions*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **112**, 1876 (2000).
104. *Computer simulation study of the density and temperature dependence of fundamental and overtone vibrational dephasing in nitrogen: Interplay between different mechanisms of dephasing*, N. Gayathri and **B. Bagchi**, *J. Phys. Chem. A* **103**, 9579 (1999).

105. *Barrierless isomerization dynamics in viscous liquids: Decoupling of the reaction rate from the slow frictional forces*, R.A. Denny and **B. Bagchi**, *J. Phys. Chem. A* **103**, 9061 (1999).
106. *Free energy gap dependence of the electron-transfer rate from the inverted to the normal region*, N. Gayathri and **B. Bagchi**, *J. Phys. Chem. A* **103** 8596(1999).
107. *Anisotropic diffusion of nonspherical molecules in dense liquids: A molecular dynamics simulation of isolated ellipsoids in the sea of spheres*, S. Ravichandran and **B. Bagchi**, *J. Chem. Phys.* **111**, 7505 (1999).
108. *Anomalous solubility of organic solutes in supercritical water: A molecular explanation*, R. Biswas and **B. Bagchi**, *Proc. Indian Acad. Chem. Sc.* **111**, 387 (1999).
109. *Subquadratic quantum number dependence and other anomalies of vibrational dephasing in liquid nitrogen: Molecular dynamics simulation study from the triple point to the critical point and beyond*, N. Gayathri and **B. Bagchi**, *Phys. Rev. Lett.* **82**, 4851 (1999).
110. *Ion conductance in electrolyte solutions*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **110**, 10024 (1999).
111. *Molecular origin of the Debye-Huckel-Onsager limiting law of ion conductance and its extension to high concentrations: Mode coupling theory approach to electrolyte friction*, A. Chandra, R. Biswas and **B. Bagchi**, *J. Am. Chem. Soc.* **121**, 4082 (1999).
112. *Solvation dynamics in nonassociated polar solvents*, R. Biswas and **B. Bagchi**, *J. Phys. Chem. A* **103**, 2495 (1999).
113. *Isomerization dynamics in viscous liquids: Microscopic investigation of the coupling and decoupling of the rate to and from solvent viscosity and dependence on the intermolecular potential*, R.K. Murarka, S. Bhattacharyya, R. Biswas and **B. Bagchi**, *J. Chem. Phys.***110**, 7365 (1999).
114. *Polar and nonpolar solvation dynamics, ion diffusion, and vibrational relaxation: Role of biphasic solvent response in chemical dynamics*, **B. Bagchi** and R. Biswas, *Adv. Chem. Phys.* **109**, 207 (1999).
115. *Time dependent diffusion coefficient and the transient dynamics of diffusion controlled bimolecular reactions in liquids: A mode coupling theory analysis*, A. Morita and **B. Bagchi**, *J. Chem. Phys.*, **110**, 8643 (1999).
116. *Limiting ionic conductance of symmetrical rigid ions in aqueous solution*, R. Biswas and **B. Bagchi**, *Ind. J. Chem.* (1999).
117. *Computer simulation and mode coupling theory study of the effects of specific solute-solvent interactions on diffusion: Crossover from a sub-slip to a super-stick*

- limit of diffusion*, G. Srinivas, S. Bhattacharyya and **B. Bagchi**, *J. Chem. Phys.* **110**, 4477 (1999).
118. *Solvation dynamics of a quadrupolar solute in dipolar liquids*, A. Sethia and **B. Bagchi**, *J. Phys. Soc. Jap.* **68**, 303 (1999).
  119. *Interplay between ultrafast polar solvation and vibrational dynamics in electron transfer reactions: Role of high-frequency vibrational modes*, **B. Bagchi** and N. Gayathri, *Adv. Chem. Phys.* **107**, 1 (1999).
  120. *Comment on "Dynamics of solvated ion in polar liquids: An interaction-site-model description"*, R. Biswas and **B. Bagchi**, *J. Chem. Phys.* **110**, 1833 (1999).
  121. *Computer simulation study of the subquadratic quantum number dependence of vibrational overtone dephasing: Comparison with the mode-coupling theory predictions*, N. Gayathri and **B. Bagchi**, *J. Chem. Phys.* **110**, 539 (1999).
  122. *Bimodality of the viscoelastic response of a dense liquid and comparison with the frictional responses at short times*, S. Bhattacharyya and **B. Bagchi**, *J. Chem. Phys.* **109**, 7885 (1998).
  123. *Ultra-fast chemical phenomena: A frontier of Physical Chemistry*, **B. Bagchi**, *Proc. Ind. Nat. Sci. Acad.* **64A**, 651 (1998).
  124. *Universality in the fast orientational relaxation near isotropic-nematic transition*, S. Ravichandran, A. Perera, and **B. Bagchi**, *J. Chem. Phys.* **109**, 7349 (1998).
  125. *Anomalous dielectric relaxation of aqueous protein solutions*, N. Nandi and **B. Bagchi**, *J. Phys. Chem. A* **102**, 8217 (1998).
  126. *A new model for promoting protein crystallization in solution*, **B. Bagchi**, *Curr. Sci.* **75**, 644 (1998).
  127. *Microscopic derivation of the Hubbard-Onsager-Zwanzig expression of limiting ionic conductivity*, **B. Bagchi**, *J. Chem. Phys.* **109**, 3989 (1998).
  128. *Self-consistent molecular theory of orientational relaxation and dielectric friction in a dense dipolar liquid*, **B. Bagchi**, *J. Mol. Liq.* **77**, 177 (1998).
  129. *Ion solvation dynamics in supercritical water*, R. Biswas and **B. Bagchi**, *Chem. Phys. Lett.* **290**, 223 (1998).
  130. *Molecular theory for the effects of specific solute-solvent interaction on the diffusion of a solute particle in a molecular liquid*, R. Biswas, S. Bhattacharyya, and **B. Bagchi**, *J. Phys. Chem. B* **102**, 3252 (1998).
  131. *Ionic mobility and ultrafast solvation: Control of a slow phenomenon by fast dynamics*, **B. Bagchi** and R. Biswas, *Acc. Chem. Res.* **31**, 181 (1998).

132. *Vibrational energy relaxation, nonpolar solvation dynamics and instantaneous normal modes: Role of binary interaction in the ultrafast response of a dense liquid*, R. Biswas, S. Bhattacharyya, and **B. Bagchi**, *J. Chem. Phys.* **108**, 4963, (1998).
133. *Dynamic solvent effects on the vibrational overtone dephasing in molecular liquids: Subquadratic quantum number dependence*, N. Gayathri, S. Bhattacharyya, and **B. Bagchi**, *J. Chem. Phys.* **107**, 10381 (1997).
134. *Dielectric relaxation of biological water*, N. Nandi and **B. Bagchi**, *J. Phys. Chem. B* **101**, 10954 (1997).
135. *Translational and rotational motion in molecular liquids: A computer simulation study of Lennard-Jones ellipsoids*, S. Ravichandran, A. Perera, and **B. Bagchi**, *J. Chem. Phys.* **107**, 8469 (1997).
136. *Solvation dynamics of a charge bubble in water*, R. Biswas and **B. Bagchi**, *Proc. Ind. Acad. Sci. Chem Sci.*, **109**, 347 (1997).
137. *Decoupling of tracer diffusion from viscosity in a supercooled liquid near the glass transition*, S. Bhattacharyya and **B. Bagchi**, *J. Chem. Phys.* **107**, 5852 (1997).
138. *From dielectric relaxation to polar solvation dynamics in alcohols and amides*, R. Biswas and **B. Bagchi**, *Ind. J. Chem.* **36A**, 635 (1997).
139. *Limiting ionic conductance of symmetrical, rigid ions in aqueous solutions: Temperature dependence and solvent isotope effects*, R. Biswas, and **B. Bagchi**, *J. Am. Chem. Soc.* **119**, 5946 (1997).
140. *Bimodality in the dynamic response of a supercooled liquid*, S. Bhattacharyya and **B. Bagchi**, *J. Chem. Phys.* **106**, 7262 (1997).
141. *Slow Polar Solvation Dynamics in a Finite Complex System: Coumarin 480 in Reverse Micelle*, N. Nandi and **B. Bagchi**, *J. Phys. Chem.*
142. *Solvation dynamics in monohydroxy alcohols: Agreement between theory and different experiments*, R. Biswas, N. Nandi, and **B. Bagchi**, *J. Phys. Chem. B* **101**, 2968 (1997).
143. *Ionic mobility in alcohols: From dielectric friction to the solvent-berg model*, R. Biswas and **B. Bagchi**, *J. Chem. Phys.* **106**, 5587 (1997).
144. *Prediction of the senses of helical amphiphilic assemblies from effective intermolecular pair potential: Studies on chiral monolayers and bilayers*, N. Nandi and **B. Bagchi**, *J. Phys. Chem.* **101**, 1343 (1997).
145. *Anomalous diffusion of small particles in dense liquids*, S. Bhattacharyya and **B. Bagchi**, *J. Chem. Phys.* **106**, 1757 (1997).



146. *Single particle and collective orientational relaxation in an anisotropic liquid near the isotropic-nematic transition*, A. Perera, S. Ravichandran, M. Moreau, and **B. Bagchi**, *J. Chem. Phys.* **106**, 1280 (1997).
147. *Molecular origin of the intrinsic bending force for helical morphology observed in chiral amphiphilic assemblies: Concentration and size dependence*, N. Nandi and **B. Bagchi**, *J. Am. Chem. Soc.* **118**, 11208 (1996).
148. *Activated barrier crossing dynamics in slow, viscous liquids*, R. Biswas and **B. Bagchi**, *J. Chem. Phys.* **105**, 7543 (1996).
149. *Solvent dynamic effects in electron transfer reactions in slow liquids: Interplay between ultra-fast solvation and vibronic coupling in Betaines*, N. Gayathri and **B. Bagchi**, *J. de Chimie Physique* **93**, 1652 (1996).
150. *Orientalional relaxation in a random dipolar lattice: Wave-number and frequency dependence*, S. Ravichandran and **B. Bagchi**, *Phys. Rev. E* **54**, 3693 (1996).
151. *Ultrafast solvation dynamics of an ion in the gamma-cyclodextrin cavity: The role of restricted environment*, N. Nandi and **B. Bagchi**, *J. Phys. Chem.* **100**, 13914 (1996).
152. *Microscopic origin of the chirality driven morphologies of the amphiphilic monolayers and bilayers*, N. Nandi and **B. Bagchi**, *Ind. J. Chem.* **35A**, 536 (1996).
153. *Self-consistent microscopic treatment of the effects of self-motion of the probe on ionic and dipolar solvation dynamics*, R. Biswas and **B. Bagchi**, *J. Phys. Chem.* **100**, 4261 (1996).
154. *Novel inter effects of High Frequency Modes in the Dynamics of Electron Transfer Reaction in the Marcus Inverted Regime*, N. Gayathri and **B. Bagchi**, *Fast Elementary Processes in Chemical and Biological Systems*, Ed. Andre Tramer, American Institute of Physics, AIP Conference Proceedings **364**, 340 (1996).
155. *Quantum and non-Markovian effects in the electron transfer reaction dynamics in the Marcus inverted region*, N. Gayathri and **B. Bagchi**, *J. Phys. Chem.* **100**, 3056 (1996).
156. *Non-Marcus energy gap dependence of electron transfer rate in contact ion pairs. Novel interplay between relaxation and reaction in solution*, N. Gayathri and **B. Bagchi**, *J. Mol. Struc. (Theo. Chem.)* **361**, 117 (1996).
157. *Solvation dynamics in slow, viscous liquids: Application to amides*, R. Biswas and **B. Bagchi**, *J. Phys. Chem.* **100**, 1238 (1996).
158. *Ultrafast salvation dynamics of an ion in a restricted environment*, N. Nandi and **B. Bagchi**, *Ind. J. Chem.* **34A**, 845 (1995).

159. *Orientational relaxation in dipolar systems: How much do we understand the role of correlations?* S. Ravichandran and **B. Bagchi**, *Int. Rev. Phys. Chem.* **14**, 271 (1995).
160. *Orientational relaxation in a random dipolar lattice: Role of spatial density fluctuations in supercooled liquids*, S. Ravichandran and **B. Bagchi**, *Phys. Rev. Lett.* **76**, 644 (1996).
161. *Anomalous ion diffusion in dense dipolar liquids*, R. Biswas, S. Roy, and **B. Bagchi**, *Phys. Rev. Lett.* **75**, 1098 (1995).
162. *How fast is ultrafast chemistry?* **B. Bagchi**, *Curr. Sci.* **69**, 129 (1995).
163. *Effects of solvent polarization relaxation on nonadiabatic outersphere electron transfer reactions in ultrafast dipolar solvents*, S. Roy and **B. Bagchi**, *J. Chem. Phys.* **102**, 7937 (1995).
164. *Adiabatic and nonadiabatic outersphere electron transfer reactions in methanol: Effects of ultrafast solvent polarization modes*, S. Roy and **B. Bagchi**, *J. Chem. Phys.* **102**, 6719 (1995).
165. *Collective effects on single-particle orientational relaxation in slow dipolar liquids*, S. Ravichandran, S. Roy, and **B. Bagchi**, *J. Phys. Chem.* **99**, 2489 (1995).
166. *Ultrafast solvation dynamics in water: Isotope effects and comparison with experimental results*, N. Nandi, S. Roy and **B. Bagchi**, *J. Chem. Phys.* **102**, 1390 (1995).
167. *A molecular explanation of the transition from viscous to hopping mechanism of mass transport in the supercooled liquid near the glass transition*, **B. Bagchi**, *J. Chem. Phys.* **101**, 9946 (1994).
168. *Non-exponential orientational relaxation in dipolar solids: The role of dipolar interactions and dielectric frictions*, S. Ravichandran and **B. Bagchi**, *J. Mol. Struct.* **327**, 247 (1994).
169. *Molecular dynamics simulations of orientational relaxation in dipolar lattice: Lack of diffusive decay for second and higher rank correlation functions*, S. Ravichandran and **B. Bagchi**, *J. Phys. Chem.* **98**, 11242 (1994).
170. *Ionic and dipolar solvation dynamics in liquid water*, N. Nandi, S. Roy, and **B. Bagchi**, *Proc. Ind. Acad. Sci. (Chem. Sci.)* **106**, 1297 (1994).
171. *Effects of ultrafast solvation on the rate of adiabatic outersphere electron transfer reactions*, S. Roy and **B. Bagchi**, *J. Phys. Chem.* **98**, 9207 (1994).
172. *Microscopic theory of ion solvation dynamics in liquid methanol*, S. Roy and **B. Bagchi**, *J. Chem. Phys.* **101**, 4150 (1994).

173. *Time dependent solution of generalized Zusman model of outersphere electron transfer reaction: Applications to various experimental situations*, S. Roy and **B. Bagchi**, *J. Chem. Phys.* **100**, 8802 (1994).
174. *Solvation dynamics, energy distribution and trapping of a light solute ion in dipolar liquids*, S. Roy and **B. Bagchi**, *Chem. Phys.* **183**, 207 (1994).
175. *Molecular theory of nonpolar solvation dynamics*, B. Bagchi, *J. Chem. Phys.* **100**, 6658 (1994).
176. *Rank dependence of orientational relaxation in dipolar systems*, S. Ravichandran and **B. Bagchi**, *J. Phys. Chem.* **98**, 2729 (1994).
177. *Dielectric relaxation in dipolar solid rotator phases*, G. V. Vijayadamodar, S. S. Komath, S. Roy, and **B. Bagchi**, *Phase Transitions* **50**, 21 (1994).
178. *Solvation dynamics in liquid water: A novel interplay between librational and diffusive modes*, S. Roy and **B. Bagchi**, *J. Chem. Phys.* **99**, 9938 (1993).
179. *Ultra-fast solvation dynamics from far-infrared spectroscopy and Kerr relaxation*, **B. Bagchi** and S. Roy in *Ultra-fast Chemical Reaction and Solvent Response*", ed. Y. Gauduel and P.J. Rossky (American Institute of Physics, New York, 1993).
180. *Molecular Theory of Solvation of an Ion*, A. Chandra and **B. Bagchi**, *Proc. Indian Acad. Sci. (Chem. Sci.)* (1993).
181. *Microscopic theory of solvation of an ion in a binary dipolar liquid*, A. Chandra and **B. Bagchi**, *J. Mol. Liq.* **57**, 39 (1993).
182. *Molecular theory of ion solvation dynamics in liquid water, acetonitrile and methano: A unified microscopic description of collective dynamics in dipolar liquids*, S. Roy and **B. Bagchi**, *Proc. Indian Acad. Sci. (Chem. Sci.)* **105**, 295 (1993).
183. *Molecular theory of ultrafast solvation in liquid acetonitrile*, S. Roy, S. Komath, and **B. Bagchi**, *J. Chem. Phys.* **99**, 3139 (1993).
184. *Ultrafast underdamped solvation: Agreement between computer simulation and various theories of solvation dynamics*, S. Roy and **B. Bagchi**, *J. Chem. Phys.* **99**, 1310 (1993).
185. *Microscopic study of inertial and viscoelastic effects in dipolar solvation dynamics*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **99**, 553 (1993).
186. *Molecular theory of underdamped dielectric relaxation: Understanding collective effects in dipolar liquids*, **B. Bagchi** and A. Chandra, *Chem. Phys.* **173**, 133 (1993).

187. *Solvation dynamics in a Brownian dipolar lattice: Comparison between computer simulation and various molecular theories of solvation dynamics*, S. S. Komath and **B. Bagchi**, *J. Chem. Phys.* **98**, 8987 (1993).
188. *Molecular expression for dielectric friction on a rotating dipole: Reduction to the continuum limit*, **B. Bagchi** and G.V. Vijayadamodar, *J. Chem. Phys.* **98**, 3351 (1993).
189. *Dielectric friction and solvation dynamics: Novel results on relaxation in dipolar liquids*, S. Roy, S. Komath, and **B. Bagchi**, *Proc. Ind. Acad. Sci. (Chem. Sci.)* **105**, 79 (1993).
190. *Solvation dynamics in a Brownian dipolar lattice: A comparison between theory and computer simulation*, H.-X. Zhou, **B. Bagchi**, A. Papazyan, and M. Maroncelli, *J. Chem. Phys.* **97**, 9311 (1992).
191. *Ultrafast solvation dynamics: Molecular explanation of computer simulation results in a simple dipolar solvent*, **B. Bagchi** and A. Chandra, *J. Chem. Phys.* **97**, 5126 (1992).
192. *Dielectric and orientational relaxation in a Brownian dipolar lattice*, H.-X. Zhou and **B. Bagchi**, *J. Chem. Phys.* **97**, 3610 (1992).
193. *Leventhal paradox*, R. Zwanzig, A. Szabo, and **B. Bagchi**, *Proc. Natl. Acad. Sci. U. S. A.* **89**, 20 (1992).
194. *Molecular interpretation of the linear relationship between the entropy and enthalpy of activation of charge transfer reactions in polar liquids*, A. Chandra and **B. Bagchi**, *Proc. Indian Acad. Sci. (Chem. Sci.)* **104**, 399 (1992).
195. *Dynamics of solvation of an ion in a dense dipolar liquid*, A. Chandra and **B. Bagchi**, *Proc. Indian Acad. Sci. (Chem. Sci.)* **104**, 675 (1992).
196. *Molecular hydrodynamic theory of non-Markovian collective orientational relaxation in dense dipolar liquids*, G. V. Vijayadamodar and **B. Bagchi**, *J. Chem. Phys.* **95**, 5289 (1991).
197. *Effects of solvent viscoelasticity in the solvation dynamics of an ion in a dense dipolar liquid*, A. Chandra and **B. Bagchi**, *Chem. Phys.* **156**, 323 (1991).
198. *Electrostriction: A density functional theory*, G.V. Vijayadamodar and **B. Bagchi**, *J. Chem. Phys.* **95**, 1168 (1991).
199. *Microscopic expression for dielectric friction on a moving ion*, **B. Bagchi**, *J. Chem. Phys.* **95**, 467 (1991).
200. *Solvation and Barrierless electron transfer*, **B. Bagchi**, A. Chandra and G.R. Fleming, in *Ultrafast Phenomena VII* eds. C.B. Harris and E. Ippen (Springer Verlag, 1991).

201. *Molecular theory of solvation and solvation dynamics in a binary dipolar liquid*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **94**, 8367 (1991).
202. *Collective orientational relaxation in dense dipolar liquids*, **B. Bagchi** and A. Chandra, *Adv. Chem. Phys.* **80**, 1 (1991).
203. *Microscopic free energy functional for polarization fluctuations: Generalization of Marcus-Felderhof expression*, A. Chandra, and **B. Bagchi**; *J. Chem. Phys.* **94**, 2258 (1991).
204. *Molecular theory of dielectric relaxation in a dense binary dipolar liquid* A. Chandra, **B. Bagchi**; *J. Phys. Chem.* **95**, 2529 (1991).
205. *Dielectric relaxation in dipolar liquids*, **B. Bagchi** and A. Chandra, *Int. Rev. Mod. Phys.* **5**, 461 (1991).
206. *Inertial effects in solvation dynamics*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.*, **94**, 3177 (1991).
207. *Relaxation dominated by inertia: Solvation dynamics of a small ion in a dipolar solvent*, A. Chandra and **B. Bagchi**, *Proc. Ind. Acad. Sci. (Chem. Sci.)* **103**, 77 (1991).
208. *Dynamics of barrierless chemical reactions in solution*, **B. Bagchi**, *Reaction Dynamics; Recent Advances*, ed. N. Sathyamurthy (Narosa/Springer-Verlag, 1990), p.12-133.
209. *An interpretation of the bifurcation of orientational relaxation processes in a supercooled liquid*, **B. Bagchi**, A. Chandra, and S.A. Rice, *J. Chem. Phys.* **93**, 8991 (1990).
210. *Collective orientational relaxation in a dense liquid of ellipsoidal molecules*, A. Chandra and **B. Bagchi**, *Physica A* **169**, 246 (1990).
211. *Macro-micro relations in dipolar orientational relaxations: An exactly solvable model of dielectric relaxation*, **B. Bagchi** and A. Chandra *J. Chem. Phys.* **93**, 1955 (1990).
212. *Dynamics of a barrierless reaction on a two-dimensional potential surface in solution*, C. S. Poornimadevi and **B. Bagchi**, *Chem. Phys. Lett.* **168**, 276 (1990)
213. *Collective excitations in a dense dipolar liquid: How important are dipolarons in the polarization relaxation of common dipolar liquids?* A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **92**, 6833 (1990).
214. *Dynamic solvent effects in adiabatic electron-transfer reactions: Role of translational modes*, **B. Bagchi**, A. Chandra, and G.R. Fleming, *J. Phys. Chem.* **94**, 5197 (1990)

215. *Effects of molecular size in solvation dynamics*, A. Chandra and **B. Bagchi**, *J. Phys. Chem.* **94**, 1874 (1990).
216. *Dielectric relaxation in dipolar liquids: Route to Debye behavior via translational diffusion*, **B. Bagchi** and A. Chandra, *Phys. Rev. Lett.* **64**, 455 (1990).
217. *Relationship between energy gap time correlation and fluorescence Stokes shift correlation functions in solvation dynamics*, A. Chandra and **B. Bagchi**, *Chem. Phys. Lett.* **165**, 115 (1990).
218. *Dynamics of activationless reactions in solution*, **B. Bagchi** and G. R. Fleming, *J. Phys. Chem.* **94**, 9 (1990).
219. *Relaxation of intermediate wave-vector density fluctuations in dense binary liquids*, G.V. Vijayadamodar and **B. Bagchi**, *J. Chem. Phys.* **92**, 6833 (1990).
220. *Relationship between microscopic and macroscopic orientational relaxational times in polar liquids*, A. Chandra and **Biman Bagchi**, *J. Phys. Chem.* **94**, 3152 (1990).
221. *Molecular theory of solvation and solvation dynamics of a classical ion in a dipolar liquid*, A. Chandra and **B. Bagchi**, *J. Phys. Chem.* **93**, 6996 (1989).
222. *Effects of translational diffusion on dielectric friction in a dipolar liquid*, G. V. Vijayadamodar, A. Chandra, and **B. Bagchi**, *Chem. Phys. Lett.* **161**, 413 (1989).
223. *Force constants of solvent polarization fluctuations: Softening at intermediate wave vectors*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **91**, 7181 (1989).
224. *Exotic dielectric behavior of polar liquids*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **91**, 3056 (1989).
225. *Breakdown of Onsager's conjecture on distance dependent polarization relaxation in solvation dynamics*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **91**, 2594 (1989).
226. *Microscopic expression for time-dependent solvation energy of ions and dipoles in dense polar liquids*, A. Chandra and **B. Bagchi**, *Proc. Indian Acad. Sci. (Chem. Sci.)* **101**, 83 (1989).
227. *Analysis of differing experimental results in barrierless reactions in solution*, **B. Bagchi**, U. Aberg, and V. Sundstrom, *Chem. Phys. Lett.* **162**, 227 (1989).
228. *Dynamics of solvation and charge transfer reactions in dipolar liquids*, **B. Bagchi**, *Ann. Rev. Phys. Chem.* **40**, 115 (1989).
229. *Non-linear structural relaxation in glassy systems: An interpretation of Narayanaswamy model*, **B. Bagchi**, *Rev. Solid State Sci.* **3**, 207 (1989).

230. *Solvation of an ion of a dipole in a dipolar liquid: How different are the dynamics?* **B. Bagchi** and A. Chandra, *Chem. Phys. Lett.* **155**, 533 (1989).
231. *On the generalized continuum model of dipolar solvation dynamics*, **B. Bagchi**, E.W. Castner and G. R. Fleming, *J. Mol. Structure* **194**, 171 (1989).
232. *A molecular theory of collective orientational relaxation in pure and binary dipolar liquids*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **91**, 1829 (1989).
233. *Polarization relaxation, dielectric dispersion, and solvation dynamics in dense dipolar liquid*, **B. Bagchi** and A. Chandra, *J. Chem. Phys.* **90**, 7338 (1989).
234. *Microscopic expression for frequency and wave vector dependent dielectric constant of a dipolar liquid*, A. Chandra and **B. Bagchi**, *J. Chem. Phys.* **90**, 1832 (1989).
235. *Dynamics of Polar Solvation : Comparison between theory and Experiment*, M. Maroncelli, E.W. Castner, **B. Bagchi** and G.R. Fleming, *Ultrafast Phenomena VI* (Springer, 1988).
236. *Dynamics of polar solvation: Route to single exponential relaxation via translational diffusion*, **B. Bagchi** and A. Chandra, *Proc. Indian Acad. Sci (Chem. Sci.)*, **100**, 353 (1988).
237. *The role of translational diffusion in the polarization relaxation in dense polar liquids*, A. Chandra and **B. Bagchi**, *Chem. Phys. Lett.* **151**, 47 (1988).
238. *The dynamics of polar solvation: Inhomogeneous dielectric continuum models*, E.W. Castner, Jr., G.R. Fleming, **B. Bagchi**, and M. Maroncelli, *J. Chem. Phys.* **89**, 3519 (1988).
239. *Fractional power dependence of rate on activation energy for reactions with very low internal barriers*, C. S. Poornimadevi and **B. Bagchi**, *Chem. Phys. Lett.* **149**, 411 (1988).
240. *Dynamics of Dipolar Solvation*, M. Maroncelli, E.W. Castner, **B. Bagchi** and G.R. Fleming", *Faraday Discussions* **85**, 1 (1988).
241. *On the stability of infinite dimensional fluid of hard hyperspheres: A statistical mechanical estimate of the density of closest packing of simple hypercubic lattices in spaces of large dimensionality*, **Biman Bagchi** and S. A. Rice, *J. Chem. Phys.* **88**, 1177 (1988).
242. *Influence of non-Debye relaxation and of molecular shape on the time dependence of the Stokes shift in Polar Media*, E. W. Castner Jr., G. R. Fleming, and **B. Bagchi**, *Chem. Phys. Lett.* **143**, 270 (1988).
243. *The dynamics of polar solvation*, E. W. Castner, Jr., **B. Bagchi**, M. Maroncelli, S. P. Webb, A. J. Ruggiero, and G. R. Fleming, *Ber. Bunsenges Phys. Chem.* **92**, 363 (1988).

244. *Freezing of a colloidal liquid subject to shear flow*, **B. Bagchi** and D. Thirumalai, *Phys. Rev. A* **37**, 2530 (1988).
245. *Numerical studies of phonon localization in a bond-disordered solid*, C. S. Poornima Devi and **B. Bagchi**, *Rev. Solid State Science*, **1**, 551 (1988).
246. *On the theory of barrierless electronic relaxation in solution*, **B. Bagchi**, *J. Chem. Phys.* **87**, 5393 (1987).
247. *On the behavior of small clusters near the spinodal decomposition*, **B. Bagchi**, *Proc. Indian Acad. Sci. (Chem. Sci.)* **99**, 243 (1987).
248. *On the temperature dependence of the rate in barrierless reactions in solutions*, **B. Bagchi**, *Chem. Phys. Lett.* **139**, 119 (1987).
249. *Fractional viscosity dependence of relaxation rates and non-steady-state dynamics in barrierless reactions in solution*, **B. Bagchi**, *Chem. Phys. Lett.* **138**, 315 (1987).
250. *Effects of non-equilibrium solvation dynamics on barrierless electronic relaxation in solution*, **B. Bagchi**, *Chem. Phys. Letts.* **135**, 553 (1987).
251. *Stability of a supercooled liquid to periodic density waves and dynamics of freezing*, **B. Bagchi**, *Physica A* **145**, 273 (1987).
252. *New results in the theory of barrierless electronic relaxation in solution*, **B. Bagchi**, *Chem. Phys. Lett.* **135**, 558 (1987).
253. *Dynamics of freezing and liquid instability*, **B. Bagchi**, *Phys. Lett. A* **121**, 29 (1987).
254. *On the breakdown of the most probable distribution to Mayer clusters*, **B. Bagchi**, *Chem. Phys. Lett.* **134**, 121 (1987).
255. *Isomerization dynamics in solution*, **B. Bagchi**, *Int. Rev. Phys. Chem.* **6**, 1 (1987).
256. *Debye-Waller factor of the solid from the self-diffusion coefficient at the solid-liquid interface*, **B. Bagchi**, *J. Chem. Phys.* **85**, 4667 (1986).
257. *Excitation wavelength and viscosity dependence of Landau-Zener electronic transition in condensed media*, **B. Bagchi**, *Chem. Phys. Lett.* **128**, 521 (1986).
258. *Dynamics of freezing*, **B. Bagchi**, *Curr. Sci.* **55**, 691 (1986).
259. *On the kinetics of crystal growth from a supercooled melt*, **B. Bagchi** and T. R. Kirkpatrick, *Proc. Ind. Acad. Sci. (Chem. Sci.)*, **96**, 465 (1986).
260. *Dynamic structure factor across the liquid-solid interface: Appearance of a delta-function elastic peak*, **B. Bagchi**, *Chem. Phys. Lett.* **125**, 91 (1986).



261. *Effect of non-equilibrium solvation dynamics on photochemical reactions in solution*, **B. Bagchi**, *J. Ind. Chem. Soc.* **LXIII**, 168 (1986).
262. *A “mean-field” theoretic analysis of the spinodal decomposition in the liquid  $\rightleftharpoons$  crystal transformation*, **B. Bagchi**, *Ind. J. Phys.* **59A**, 356 (1985).
263. *A Comment on the consistency of truncated nonlinear integral equation based theories of freezing*, C. Cerjan, **B. Bagchi**, and S. A. Rice, *J. Chem. Phys.* **83**, 2376 (1985).
264. *Self-diffusion across the liquid-crystal interface*, **B. Bagchi**, *J. Chem. Phys.* **82**, 5677 (1985)
265. *A conjecture concerning transformation of a supercooled hard sphere liquid to a metastable disordered solid*, S. A. Rice, C. Cerjan, and **B. Bagchi**, *J. Chem. Phys.* **82**, 3350 (1985).
266. *Isomerization dynamics in solution in the absence of an activation barrier: Evaluation of the potential parameters*, **B. Bagchi**, *Chem. Phys. Lett.* **115**, 209 (1985).
267. *Theory of freezing in simple systems*, C. Cerjan and **B. Bagchi**. *Phys. Rev. A* **31**, 1647 (1985).
268. *Sol  $\rightarrow$  Gel transition with rings in a finite polycondensing system in solution*, **B. Bagchi**, *Chem. Phys. Lett.* **113**, 597 (1985).
269. *Diffusion in a two-dimensional periodic potential*, **B. Bagchi**, R. Zwanzig, and M.C. Marchetti, *Phys. Rev. A* **31**, 892 (1985).
270. *Freezing of the classical two-dimensional, one-component plasma*, P. L. Radloff, **B. Bagchi**, C. Cerjan, and S.A. Rice, *J. Chem. Phys.* **81**, 1406 (1984).
271. *Crystallization of the classical one-component plasma*, **B. Bagchi**, C. Cerjan, U. Mohanty, and S.A. Rice, *Phys. Rev. B* **29**, 2857 (1984).
272. *Theory of the time development of the Stokes shift in polar media*, **B. Bagchi**, D. W. Oxtoby and G. R. Fleming, *Chem. Phys.* **86**, 257 (1984).
273. *A study of the freezing transition in the Lennard-Jones system*, **B. Bagchi**, C. Cerjan, and S.A. Rice, *J. Chem. Phys.* **79**, 6222 (1983).
274. *Theory of non-Markovian exciton transport in a one-dimensional lattice*, **B. Bagchi** and D.W. Oxtoby, *J. Chem. Phys.* **79**, 6211 (1983).
275. *Theory of electronic relaxation in solution in the absence of an activation barrier*, **B. Bagchi**, G.R. Fleming, and D.W. Oxtoby, *J. Chem. Phys.* **78**, 7375 (1983).

276. *The effect of frequency dependent friction on isomerization dynamics in solution*, **B. Bagchi** and D. W. Oxtoby, *J.Chem. Phys.* **78**, 2735 (1983).
277. *Contribution to the theory of freezing*, **B. Bagchi**, C. Cerjan, and S. A. Rice, *J. Chem. Phys.* **79**, 5595 (1983).
278. *Theoretical analysis of the achievement of random close packing of hard spheres and a conjecture on spinodal decomposition*, **B. Bagchi**, C. Cerjan, and S.A. Rice, *Phys. Rev. B* **28**, 6411 (1983).
279. *Non-monotonic dependence of electronic relaxation rate on solvent viscosity*, **B. Bagchi**, S. Singer, and D.W. Oxtoby, *Chem. Phys. Lett.* **99**, 225 (1983).
280. *Crossover between Gaussian and self-avoiding limits via finite order self-avoiding walk: Conformation space renormalization for polymers. VII*, **B. Bagchi** and Y. Oono, *J. Chem. Phys.* **78**, 2044 (1983).
281. *Agreement between the gelation and molecular dynamics models of the hydrogen-bond network in water*, **B. Bagchi**, E. Donoghue and J. H. Gibbs, *Chem. Phys. Letts.* **94**, 253 (1983).
282. *Interference effects on relaxation in three-level systems: Breakdown of the rate equation description*, **B. Bagchi** and D.W. Oxtoby, *J. Chem. Phys.* **77**, 1391 (1982).
283. *Spatial Bose condensation: Universal features in size distribution of clusters*, **B. Bagchi** and U. Mohanty, *Phys. Lett.* **91A**, 77 (1982).
284. *Effect of excitation on non-Markovian vibrational energy relaxation*, **B. Bagchi** and D.W. Oxtoby, *J. Phys. Chem.* **86**, 2197 (1982).
285. *Bimodality and long-range order in ideal Bose systems*, U. Mohanty, **B. Bagchi** and J. H. Gibbs, *J. Stat. Phys.* **28**, 685 (1982).
286. *Bimodality of cluster-size distribution and condensation in a finite Lennard-Jonnes system*, J. H. Gibbs, **B. Bagchi**, and U. Mohanty, *Phys. Rev. B* **24**, 2893 (1981).
287. *From gelation to gas-liquid transition and some contributions to Bose-Einstein condensation in finite systems*, **B. Bagchi**. Ph.D. Thesis (1981), Brown University, Providence, RI 02912).
288. *New recursion relation for the moments of the random-flight problem*, **B. Bagchi**, *J. Phys. Chem.* **85**, 4039 (1981).

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