

Mark N. Kobrak

January 2009

Department of Chemistry
Brooklyn College of the City University of New York
2900 Bedford Ave.
Brooklyn, NY 11210-2889

Tel: (718) 951-5758
Fax: (718) 951-4607
mkobrak@brooklyn.cuny.edu
<http://userhome.brooklyn.cuny.edu/mkobrak>

Education:

- **University of Chicago** Ph.D., Physical Chemistry (1997)
Thesis: "Aspects of Selective Photochemistry"
Advisor: Stuart A. Rice
- **Northwestern University** B.A., Chemistry and Integrated Science, with Honors (1992)

Research Experience:

- Jan 2006-present: **Brooklyn College of the City University of New York** Associate Professor of Chemistry
- Sept 2001-Jan 2006: **Brooklyn College of the City University of New York** Assistant Professor of Chemistry
- Jan 2000-Aug 2001: **The University of Notre Dame** and **The Pennsylvania State University** Post-doctoral fellow with Sharon Hammes-Schiffer, Department of Chemistry
- 1998-Jan 2000: **The University of Houston** Post-doctoral fellow with Eric R. Bittner, Department of Chemistry
- 1992-1998: **The University of Chicago** Graduate student with Stuart A. Rice, Department of Chemistry
- 1990-1992 **Northwestern University** Undergraduate researcher with Mark A. Ratner, Department of Chemistry
- 1989 **Northwestern University** Undergraduate researcher with Emile Okal, Department of Geology

Publications:

- *Coupled Ion Complexation and Exchange between Aqueous and Ionic Liquid Phases: A Thermodynamic Interpretation*, M.N. Kobrak, *Solv. Extr. Ion. Exch.* **26** 735 (2008).
- *The Relationship between Ionic Structure and Viscosity in Room-Temperature Ionic Liquids*, H. Li, M. Ibrahim, I. Agberemi and M.N. Kobrak, *J. Chem. Phys.* **129** 124507 (2008).
- *Laboratory Manual in General Chemistry*, M.N. Kobrak, *ed.*, Kendall-Hunt, Dubuque, IA 2008.
- *The Chemical Environment of Ionic Liquids: Links Between Liquid Structure, Dynamics and Solvation*, M.N. Kobrak, *Adv. Chem. Phys.* **139** 85 (2008).
- *The Relationship Between Solvent Polarity and Molar Volume in Room-Temperature Ionic Liquids* M.N. Kobrak, *Green Chem.* **9** 80 (2008).
- *A Comparative Study of Solvation Dynamics in Room Temperature Ionic Liquids*, M.N. Kobrak, *J. Chem. Phys.* **127** 184507 (2007).
- *Electrostatic Interactions of a Neutral Dipolar Solute with a Fused Salt: A New Model for Solvation in Ionic Liquids* M.N. Kobrak, *J. Phys. Chem. B* **111** 4755 (2007).

- *Lewis Structure Representation of Free Radicals Similar to ClO*, W. Hirsch and M. Kobrak, *J. Chem. Ed.* **84** 1360 (2007).
- *SmartTutor: A Unified Approach for Enhancing Science Education*, K. Harrow, R. Eckhardt, D. Kopec, M. Kobrak and P. Whitlock, *J. Comp. Sci. in Coll.* **22** 29 (2007).
- *Characterization of the Solvation Dynamics of a Room-Temperature Ionic Liquid via Molecular Dynamics Simulation*, M.N. Kobrak, *J. Chem. Phys.*, **125** 064502 (2006).
- *Understanding Organic Processes in Ionic Liquids: Achievements So Far and Challenges Remaining*, J. B. Harper and M.N. Kobrak, *Mini-Rev. in Org. Chem.* **3** 253 (2006).
- *An Electrostatic Interpretation of Structure-Property Relationships in Ionic Liquids*, M.N. Kobrak and N. Sandalow, in “*Molten Salts XIV*, R. Mantz, ed., The Electrochemical Society, Pennington, NJ, 2006.
- *Solvation Dynamics of Room-Temperature Ionic Liquids: Evidence for Collective Solvent Motion on Sub-Picosecond Timescales*, M.N. Kobrak and V. Znamenskiy, *Chem. Phys. Lett.* **395** 127 (2004).
- *A Molecular Dynamics Study of Polarity in Room-Temperature Ionic Liquids*, V. Znamenskiy and M.N. Kobrak, *J. Phys. Chem. B* **108** 1072 (2004).
- *Error Estimation in Histogram-Based Free Energy Calculations*, M.N. Kobrak, *J. Comp. Chem.* **24** 1437 (2003).
- *Doppler Shift and Energy Transfer to a Solar Sail*, W. Hirsch and M.N. Kobrak, *Physics Ed.* **37** 422 (2002).
- *Molecular Dynamics Simulation of Proton-Coupled Electron Transfer in Solution*, M.N. Kobrak and S. Hammes-Schiffer, *J. Phys. Chem. A* **105** 10435 (2001).
- *Reaction Path Hamiltonian Analysis of Dynamical Solvent Effects for a Claisen Rearrangement and a Diels-Alder Reaction*, H. Hu, M.N. Kobrak, C. Xu, and S. Hammes-Schiffer, *J. Phys. Chem. A* **104** 8058 (2000).
- *Quantum Simulations of Polaron Recombination Dynamics in Linear Polyenes*, E. R. Bittner and M. N. Kobrak, *Synth. Metals* **121** 1635 (2001).
- *A Quantum Molecular Dynamics Study of Polaron Recombination in Conjugated Polymers*, M. N. Kobrak and E. R. Bittner, *Phys. Rev. B* **62** 11473 (2000).
- *A Quantum Molecular Dynamics Study of Exciton Self-Trapping in Conjugated Polymers: Temperature Dependence and Spectroscopy*, M. N. Kobrak and E. R. Bittner, *J. Chem. Phys.* **112** 7684 (2000).
- *A Dynamic Model for Exciton Self-Trapping in Conjugated Polymers I: Theory*, M. N. Kobrak and E. R. Bittner *J. Chem. Phys.* **112** 5399 (2000).
- *A Dynamic Model for Exciton Self-Trapping in Conjugated Polymers II: Implementation*, M. N. Kobrak and E. R. Bittner *J. Chem. Phys.* **112** 5410 (2000).
- *The Equivalence of Photoselective Adiabatic Passage and the Strong Field Brumer-Shapiro Approach*, M.N. Kobrak and S.A. Rice, *J. Chem. Phys.* **109** 1 (1998).
- *Selective Photochemistry via Adiabatic Passage: An Extension of StiRAP for Degenerate Final States*, M.N. Kobrak and S.A. Rice, *Phys. Rev. A* **57** 2885 (1998).
- *Coherent Population Transfer via a Resonant Intermediate State: The Breakdown of Adiabatic Passage*, M.N. Kobrak and S.A. Rice, *Phys. Rev. A* **57** 1158 (1998).
- *The Influence of High-Frequency Modes on Two Pulse Spectroscopy*, M.N. Kobrak and S.A. Rice, *J. Chem. Phys.* **107** 4091 (1997).

- *The Influence of High-Frequency Modes on Ultrashort Pulse Absorption Initiated Processes*, M.N. Kobrak, E.M. Hiller, and S.A. Rice, *J. Chem. Phys.* **105** 9403 (1996).

Awards:

- Petroleum Research Fund Type-G Grant (2005-2007)
- Department of Energy Faculty and Student Team Fellowship (2003, 2004)
- PSC-CUNY Research Grant (2002)
- PSC-CUNY Incentive Grant (2001)