

**Curriculum Vitae:** *Dmitry Bedrov*

Associate Professor, Department of Materials Science and Engineering, University of Utah, 122 S. Central Campus Drive, Rm. 304, University of Utah, Salt Lake City, UT 84112.  
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**Professional Preparation**

Odessa State Academy of Refrigeration, Odessa, Ukraine	Thermophysics	<b>B.S. with Honors</b> 1995.
University of Utah, Salt Lake City, Utah	Chemical & Fuels Engineering	<b>Ph.D</b> 1999.
University of Utah, Salt Lake City, Utah	Computational Modeling of Materials	1999-2000 (post-doctoral fellow)

**Appointments**

University of Utah, Department of Materials Science and Engineering	Associate Professor	08/2013-present
University of Utah, Department of Materials Science and Engineering	Research Associate Professor	06/2011-08/2013
University of Utah, Department of Materials Science and Engineering	Research Assistant Professor	06/2000-06/2011
Wasatch Molecular Inc.	President	11/2011-present time
Wasatch Molecular Inc.	Vice President of Research	05/2006-10/2011

**Awards**

Alexander von Humboldt Experienced Researcher Fellowship to study "Multiscale Modeling of Non-equilibrium Block-Copolymer Morphologies for Energy Applications" (2011-2014)

**Professional Activities and Services:**

- Assistant Program Manager, Member of the Executive Committee, and the Lead for the Area of Electrochemical Energy Devices in the ***Collaborative Research Alliance*** “Computationally Guided Design of Energy Efficient Electronic Materials” between the University of Utah and the Army Research Lab.
- Since 2012 a member of the University of Utah ***Center for High Performance Computing*** allocation committee.
- Regular reviewer for variety of premier scientific journals such as *Macromolecules*, *Journal of Chemical Physics*, *Journal of Physical Chemistry*, *Langmuir*, *Journal of American Chemical Society*, *Physical Review (E,B, and Letters)*, *Polymer*, *Journal of Polymer Science Part B: Polymer Physics*, *Nanoletters*, *Journal of Non-crystalline Solids*, and *Philosophical Magazine*.

### Selected publications:

Chen, D.; Porada, J.H.; Hooper, J.B.; Klittnick, A.; Shen, Y.; Korblova, E.; **Bedrov, D.**; Walba, D.M.; Glaser, M.A.; MacLennan, J.E.; Clark, N.A.; "Chiral Helical Ground State of Nanoscale Pitch in a Nematic Liquid Crystal of Achiral Molecular Dimers" *Proc. Nat. Acad. Sci.* **2013**, *110*, 15931-15936.

Vatamanu, J.; Hu, Z.; **Bedrov, D.**; Perez, C.; Gogotsi, Y.; "Increasing Energy Storage in Electrochemical Capacitors with Ionic Liquid Electrolytes and Nanostructured Carbon Electrodes" *J. Phys. Chem. Lett.* **2013**, *4*, 2829–2837.

Xing, L.; Vatamanu, J.; Borodin, O.; **Bedrov, D.**; "On the Atomistic Nature of Capacitance Enhancement Generated by Ionic Liquid Electrolyte Confined in Subnanometer Pores" *J. Phys. Chem. Lett.* **2013**, *4*, 132-140.

Steinmüller, B.; Mueller, M.; Hambrecht, K.; **Bedrov, D.**; "Random Block Copolymers: Structure, dynamics and mechanical properties in the bulk and at selective substrates" *Macromolecules* **2012**, *45*, 9841-9853.

Xing, L.; Vatamanu, J.; Borodin, O.; Smith, G.D.; **Bedrov, D.**; "Electrode/Electrolyte Interface in Sulfolane-based Electrolytes for Li-Ion Batteries: A Molecular Dynamics Simulation Study" *J. Phys. Chem. C* **2012** *116*, 23871-23881.

Xing, L.; Vatamanu, J.; Smith G.D.; **Bedrov, D.**; "Nanopatterning of electrode surfaces as a route to improved energy storage in electrostatic capacitors" *J. Phys. Chem. Lett.* **2012**, *3*, 1124-1129.

**Bedrov, D.**; Borodin, O.; Li, Z.; Smith G.D.; "Influence of Polarization on Structural, Thermodynamic and Dynamic Properties of Ionic Liquids Obtained from Molecular Dynamics Simulations" *J. Phys. Chem. B* **2010**, *114*, 4984-4997.

Smith, G.D.; **Bedrov, D.**; "Dispersing Nanoparticles in Polymer Matrix: Are Long, Dense Polymer Tethers Really Necessary?" *Langmuir* **2009**, *25*, 11239–11243.

**Bedrov, D.** J.B. Hooper, G.D. Smith, T.D. Sewell " Shock-induced Transformations in Crystalline RDX: A Uniaxial Constant-Stress Hugoniotat Molecular Dynamics Simulation Study" *J. Chem. Phys.* **2009**, *131*, 034712.

Liu, W.; **Bedrov, D.**; Kumar, S.K.; Veytsman, B.; Colby, R.H. "Local Composition Distributions of Miscible Polymer Blends Probed by Molecular Dynamics Simulation" *Phys. Rev. Lett.* **2009**, *103*, 037801.

**Bedrov, D.**; Smith, G.D.; Davande, H.; Li, L. "Passive Transport of C<sub>60</sub> Fullerenes into Lipid Membrane. A Molecular Dynamics Simulation Study" *J. Phys. Chem. B* **2008**, *112*, 2078-2084.

**Bedrov, D.**; Smith, G.D.; Yoon, J.; "Structure and Interactions in Micellar Solutions: Molecular Simulations of Pluronic L64 Aqueous Solutions." *Langmuir* **2007**, *23*, 12032-12041.

**Bedrov, D.**; Smith, G.D.; "A Molecular Dynamics Simulation Study of Segmental Relaxation Processes in Miscible Polymer Blends" *Macromolecules* **2006**, *39*, 8526-8535.

Li, L.; **Bedrov, D.**; Smith, G.D. "Repulsive Solvent-induced Interaction between C<sub>60</sub> Fullerenes in Water" *Phys. Rev. E* **2005**, *71*, 011502 (1-4).

**Bedrov, D.**; Smith, G.D.; Li.L.; "A Molecular Dynamics Simulation Study of the Role of Evenly-spaced Poly(ethylene oxide) Tethers on the Aggregation of C<sub>60</sub> Fullerenes in Water", *Langmuir*, **2005**, *21*, 5251-5255.

He, Y.; Lutz, T.R.; Ediger, M.D.; Ayyagari, C.; **Bedrov, D.**; Smith, G.D. "NMR Experiments and Molecular Dynamics Simulations of the Segmental Dynamics of Polystyrene" *Macromolecules*, **2004**, *37*, 5032-5039.

**Bedrov, D.**; Smith, G.D., Smith, J.S. " Matrix-induced Nanoparticle Interactions in Polymer Melts. A Molecular Dynamics Simulation Study" *J. Chem. Phys.* **2003**, *119*, 10438-10447.

Sewell, T.D.; Menikoff, R.; **Bedrov, D.**; Smith G.D. " A Molecular Dynamics Simulation Study of Elastic Properties of HMX" *J. Chem. Phys.* **2003**, *119*, 7417-7426.

Smith, G.D., **Bedrov, D.**, Borodin, O., "A Molecular Dynamics Simulation Study of Hydrogen Bonding in Aqueous Poly(ethylene oxide) Solutions" *Phys. Rev. Lett.* **2000**, *85*, 5583-5586.

Smith, G.D., **Bedrov, D.**, Borodin, O., "Conformations and Chain Dimensions of Poly(ethylene oxide) in Aqueous Solution: A Molecular Dynamics Simulation Study" *J. Am. Chem. Soc.* **2000**, *122*, 9548-9549.