



UNIVERSITY *of* DELAWARE

Nanotechnology Research

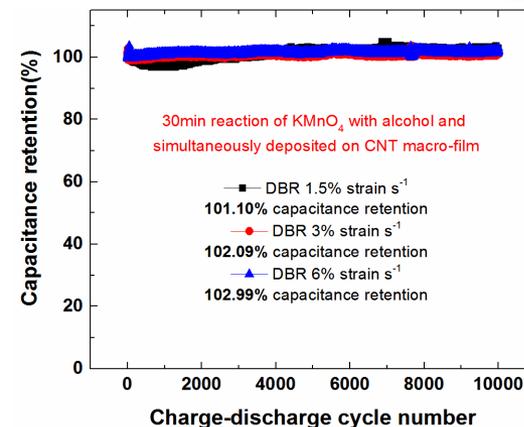
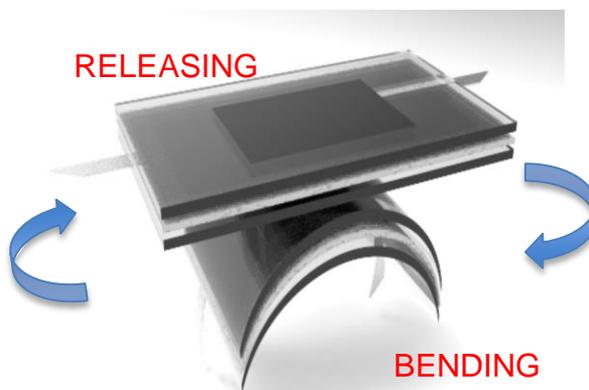
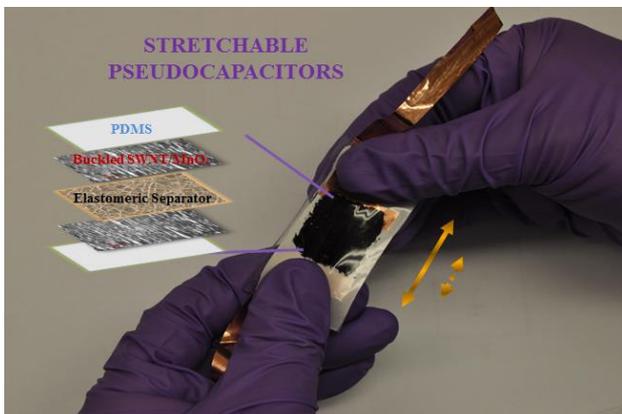
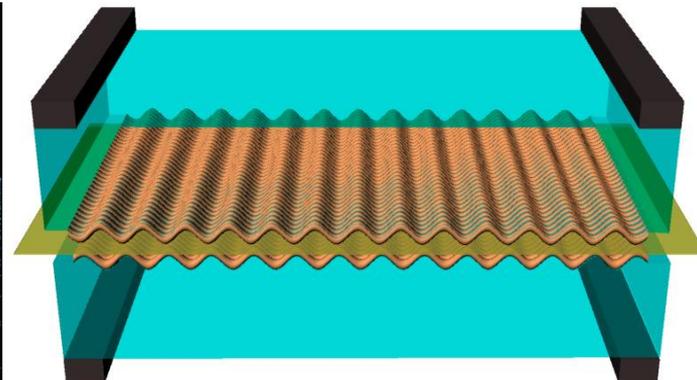
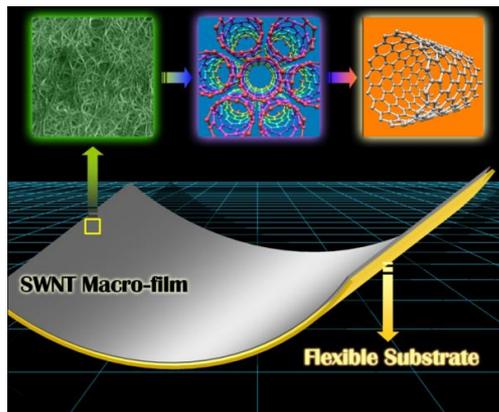
Mechanical Engineering
University of Delaware



Nanomaterial-Enabled Flexible/Stretchable Energy Storage Devices

- Flexible and Stretchable Double-layered Supercapacitors
- Stretchable and Bendable Pseudocapacitors
- Stable, Repeatable, Static and Dynamic Bending and Stretching

Statically and Dynamically Stretchable Double-layered Supercapacitors from CNTs

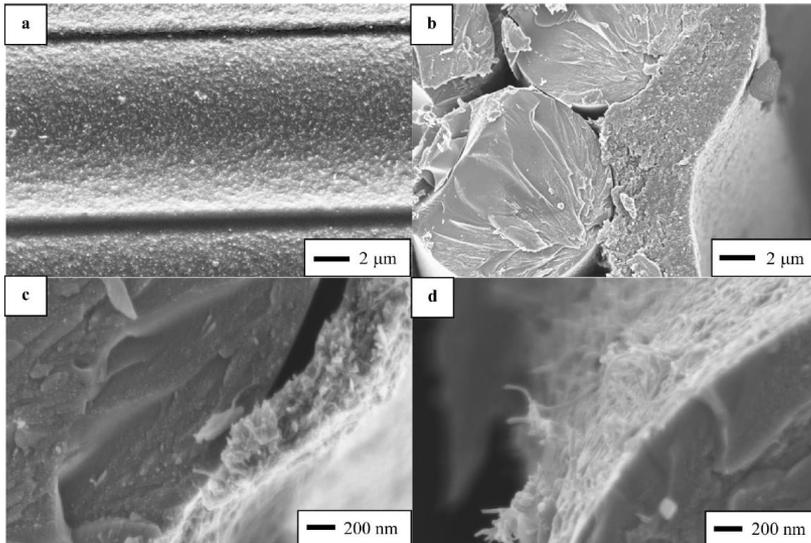
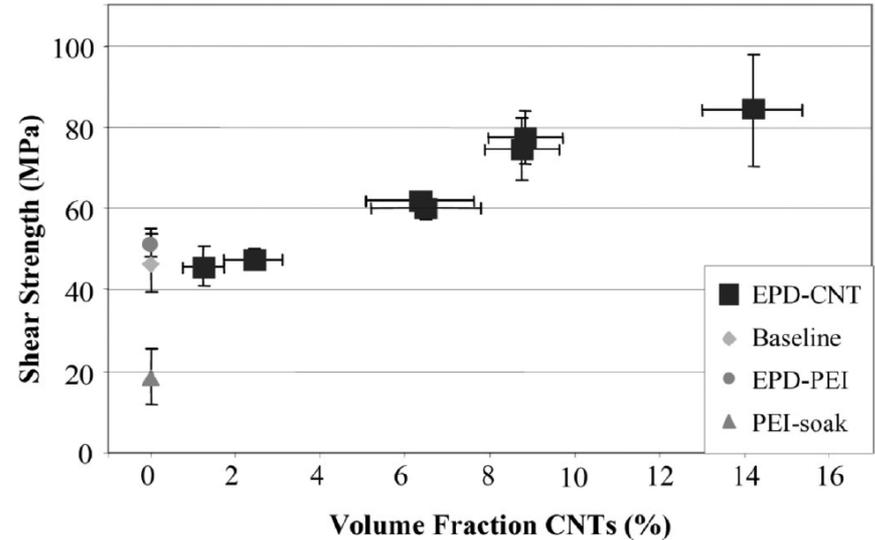
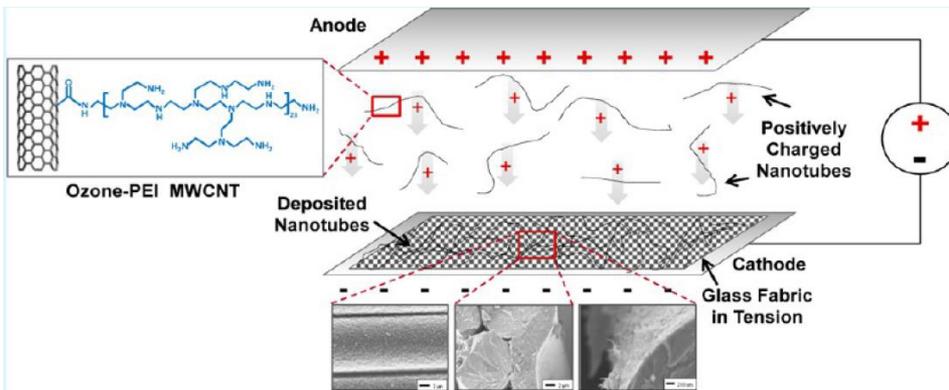


Stretchable pseudocapacitors based on buckled SWNT/MnO₂ hybrid films

Dynamic Bending/Releasing (DBR) Charge/Discharge Retention



Processing of Multifunctional Nanostructured Composites



- **Electrochemical assembly**
- **Novel AC approach for nanotube integration**
- **Demonstrated for the first time on glass**
- **Formation of stiff and strong interphase**

Funding: NSF CAREER

An, Rider and Thostenson, Carbon (2012)

An, Rider and Thostenson, ACS Applied Materials and Interfaces (2013)

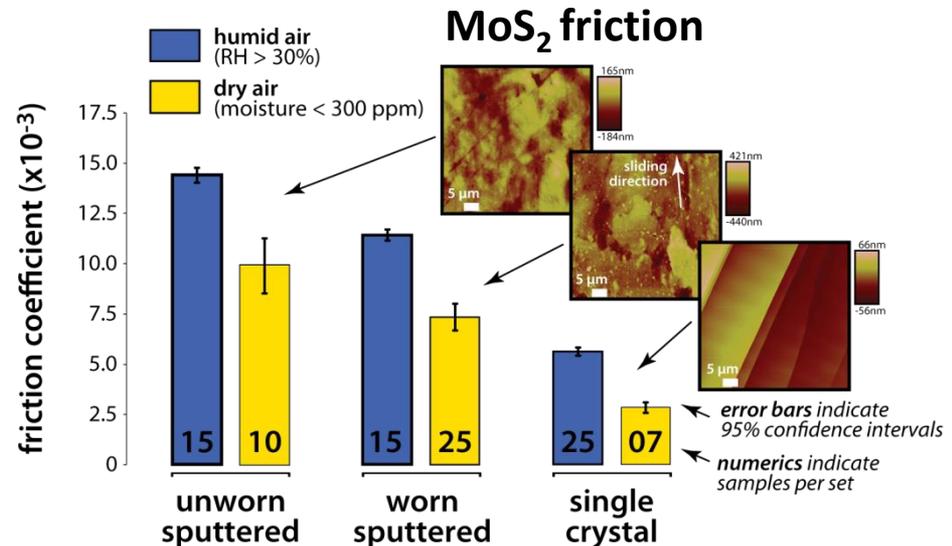
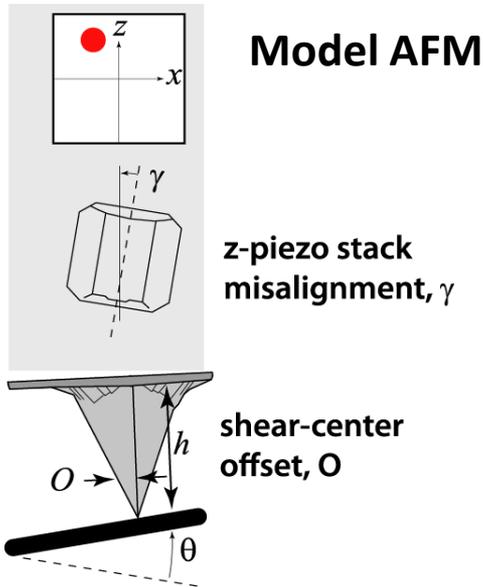
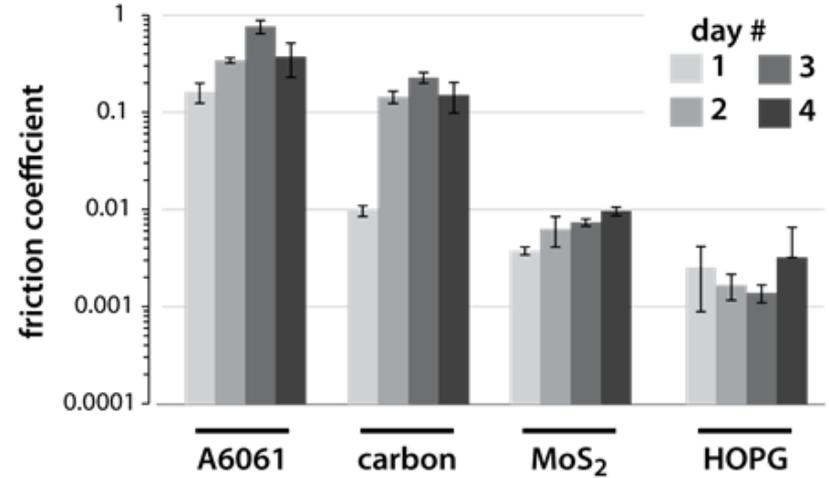


David Burris

Quantifying Friction at the Nanoscale

Motivation

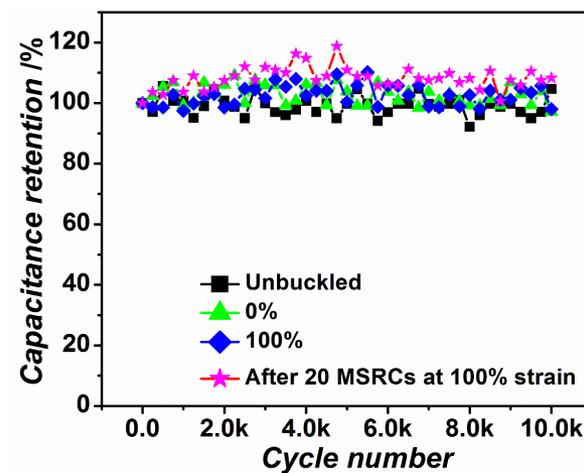
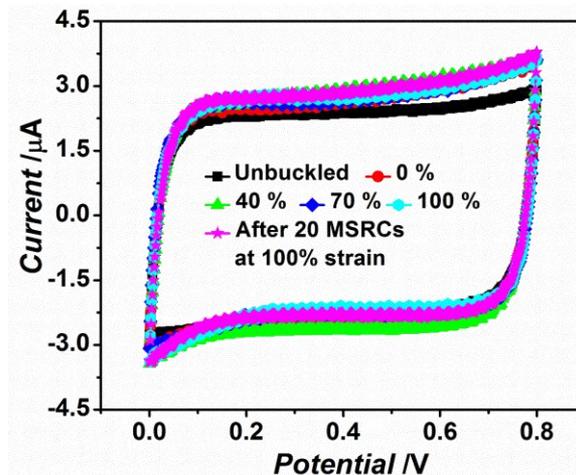
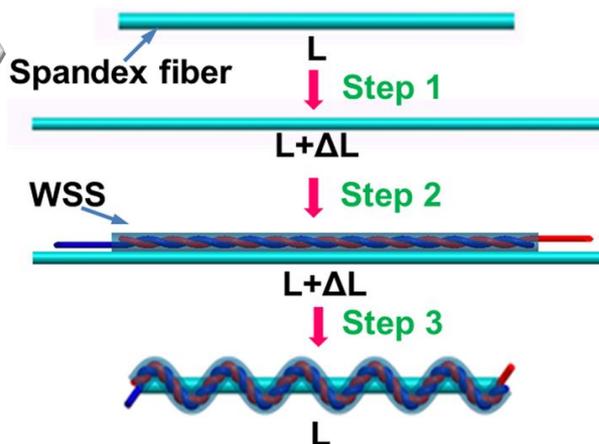
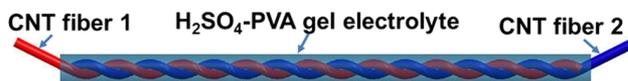
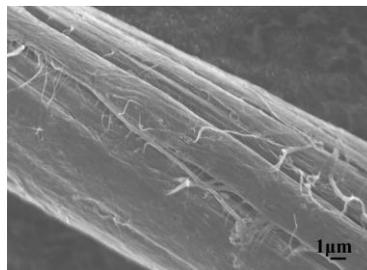
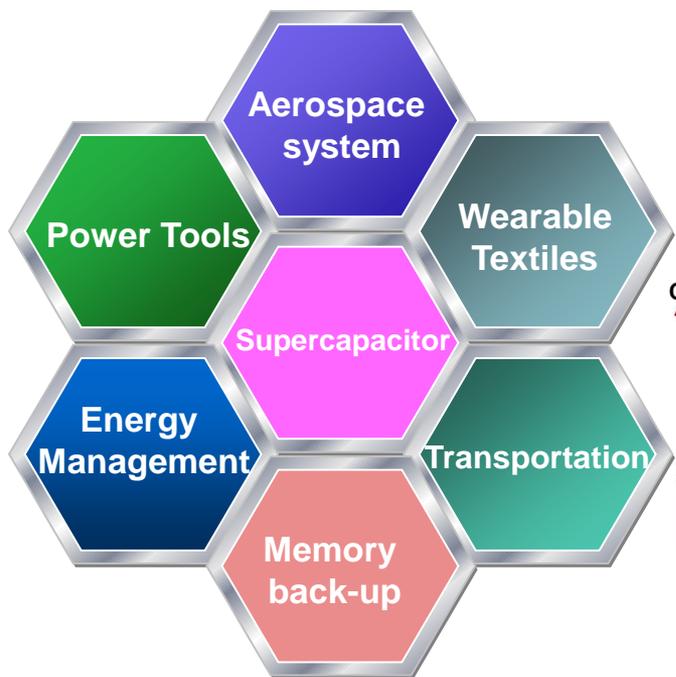
- Macroscale friction is due to an ensemble of fundamental frictional interactions
- AFM: probes fundamental interactions, calibration is difficult, setup adjustments can change calibration constants
- Khare and Burris, *RSI* 2013: Developed '**extended wedge method**' to calibrate AFM voltage signals **during** measurement





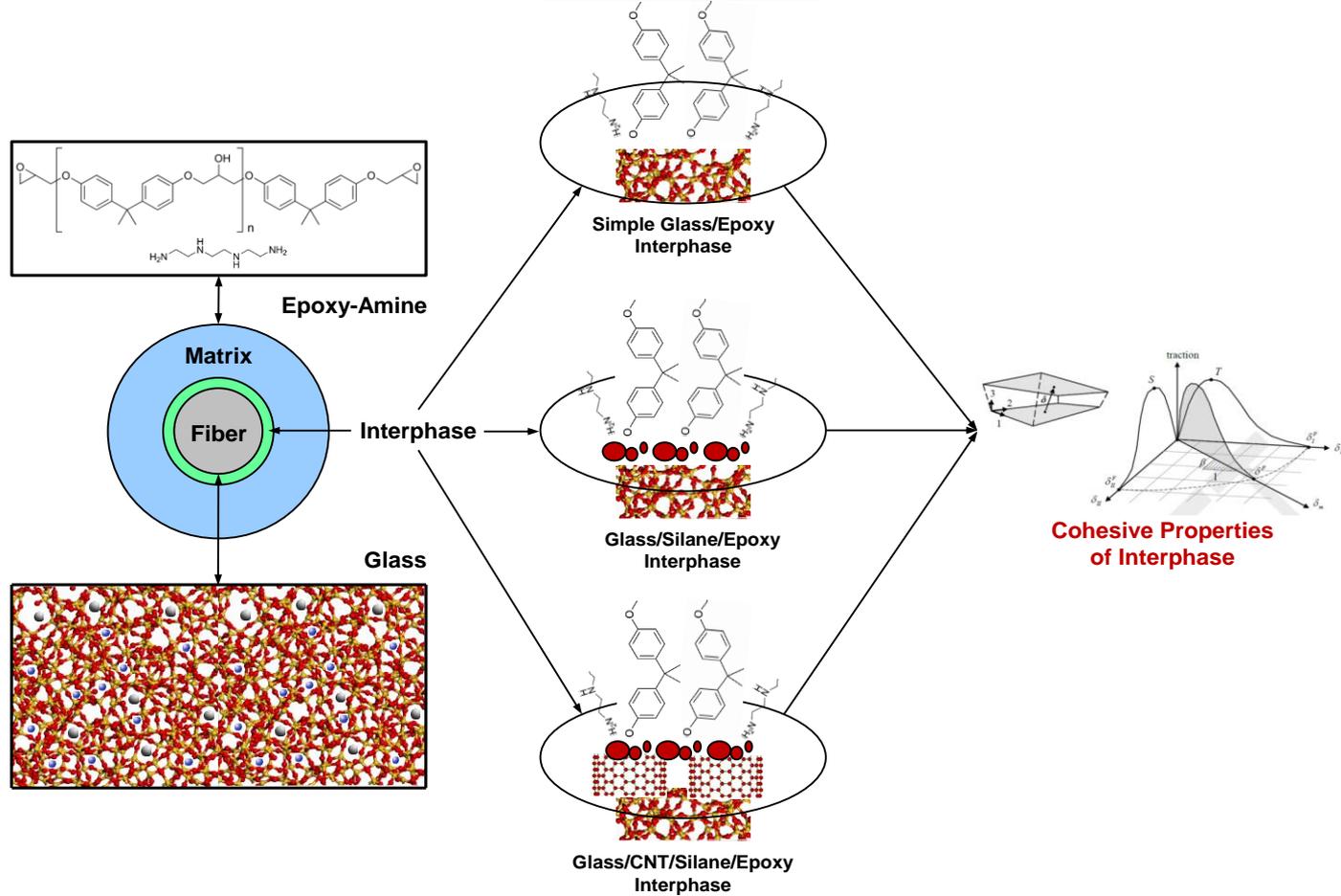
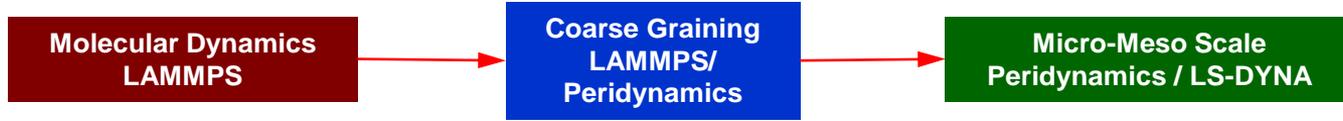
Tsu-Wei Chou

Nanomaterials for Energy Storage Devices Wire-shaped Supercapacitor



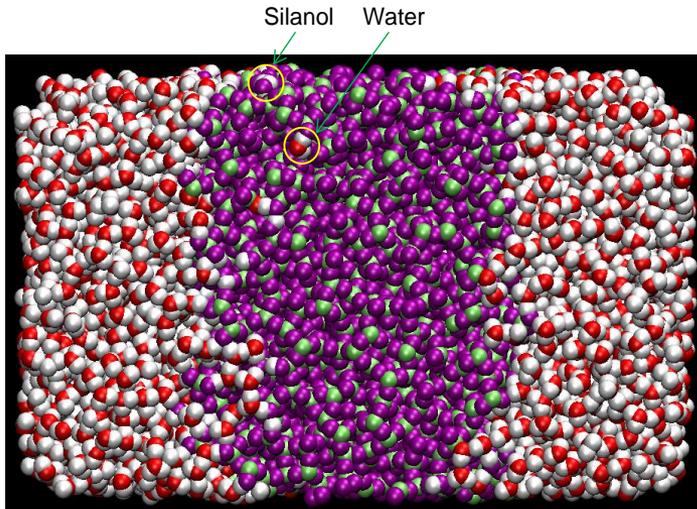


Molecular Dynamics Modeling of Interphase Formation

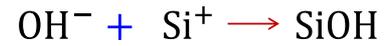
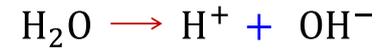
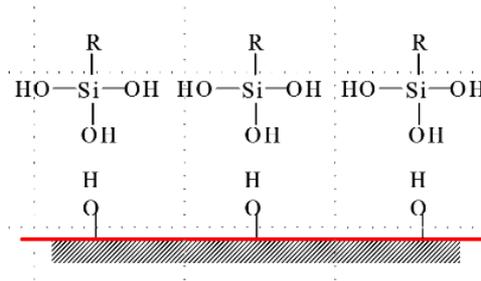




Molecular Dynamics Modeling of Interphase Formation



SILANOL FORMATION

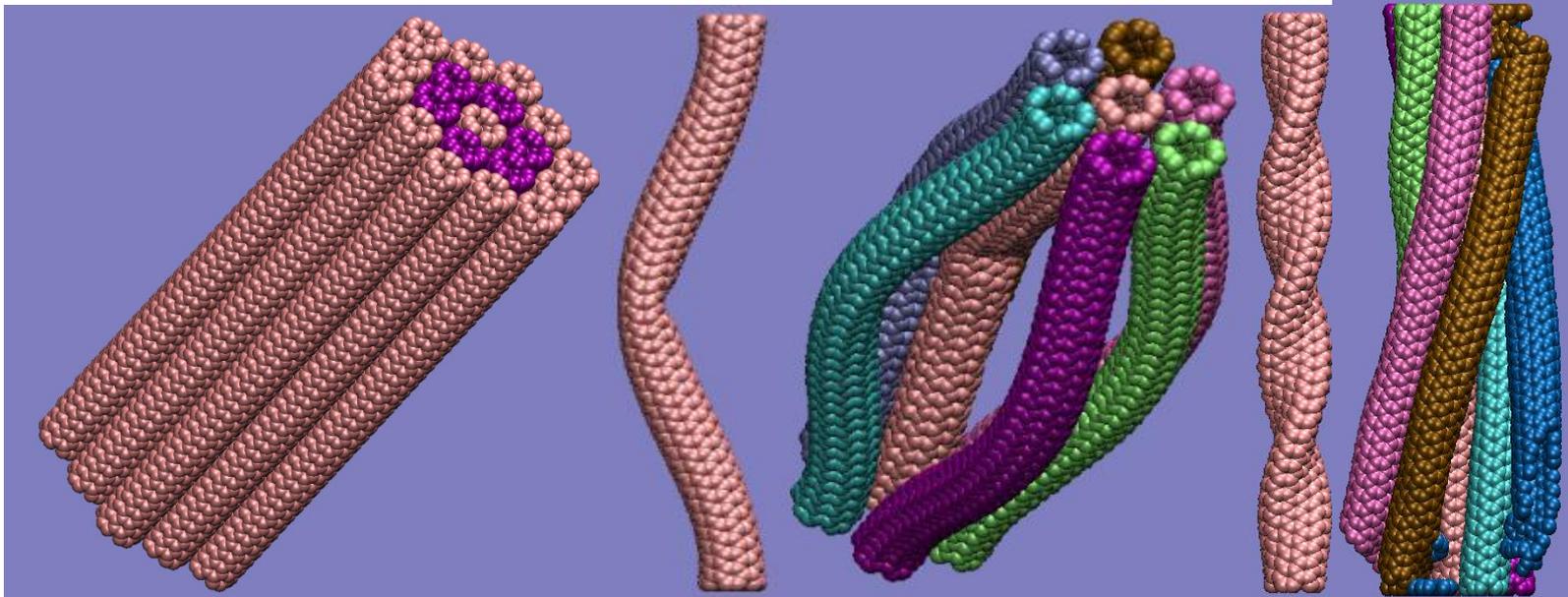


REACTIONS

- MD modeling of interphase formation during processing
- MD prediction of Mixed-Mode Traction Laws Based on Bond Breakage (strength and energy absorption)
- Bridge length scales using Peridynamics



Molecular Dynamics Modeling of Nanotube Bundles

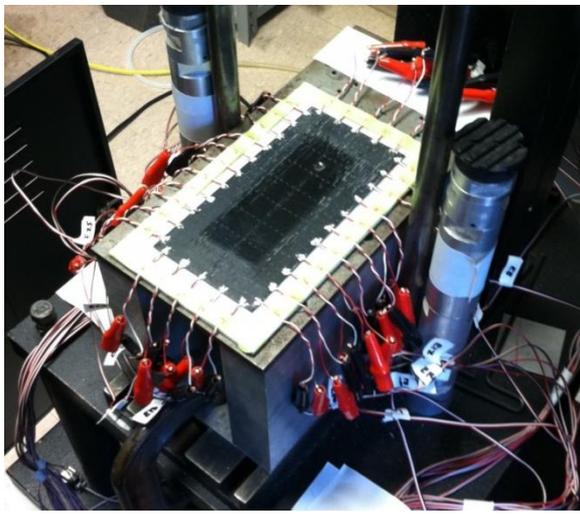
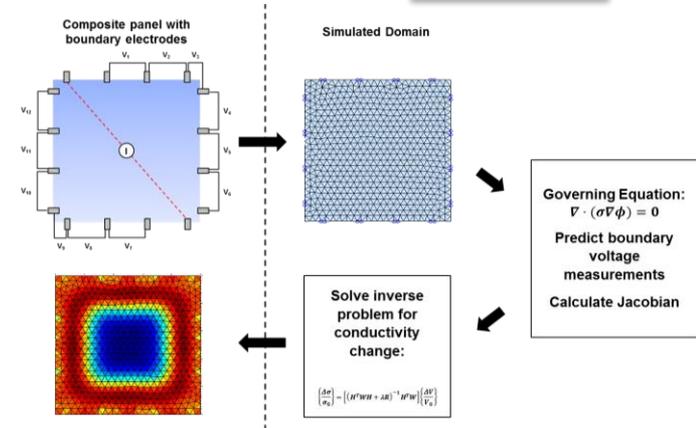
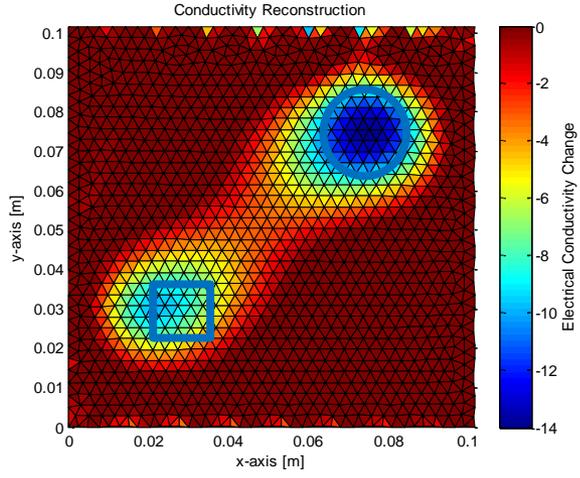
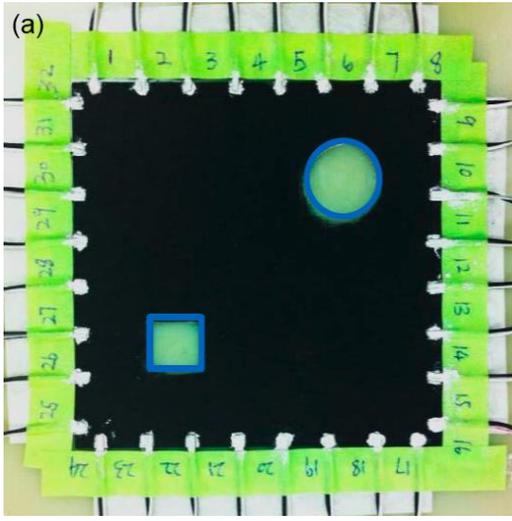


- Carbon nanotubes can form bundle where the axially aligned CNTs are packed closely together to create a larger diameter yarn.
- The inter-tube interaction in the bundle can be non-bonded van der Waals interaction or bonded sp^3 interaction.
- Such bonded and non-bonded inter-tube interaction may change the load carrying capability and damage modes at failure of the bundle as well as individual CNT within the bundle.

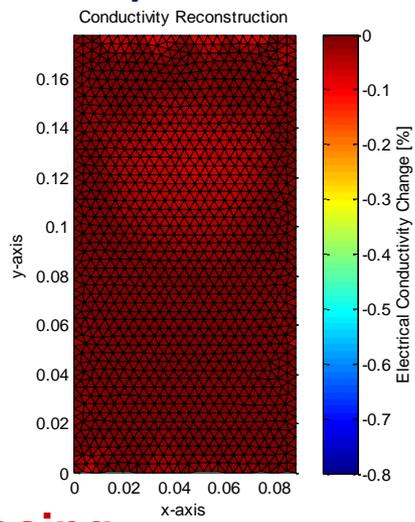


Erik Thostenson

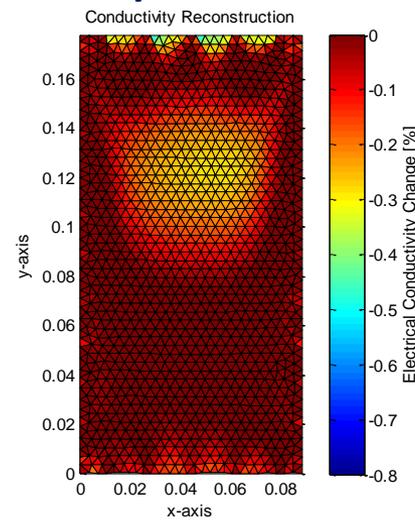
Carbon Nanotubes for *in situ* Sensing of Deformation and Damage



Impact #1



Impact #2



Impact #3

