Atomistic and electronic modeling of materials science and strength

# Umeno Lab.

### [Atomistic and electronic modeling analysis of strength and physical properties of solids]

**Center for Research on Innovative Simulation Software** 

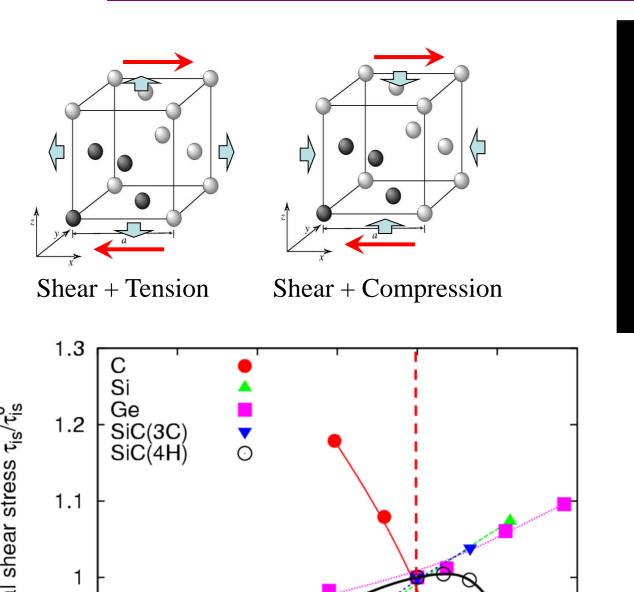
http://www.cmsm.iis.u-tokyo.ac.jp

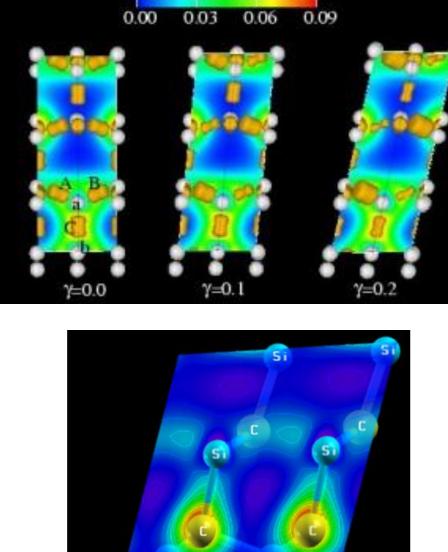
**Nanostructured Materials Strength and Science** 

**Dept. of Mechanical Engineering** 

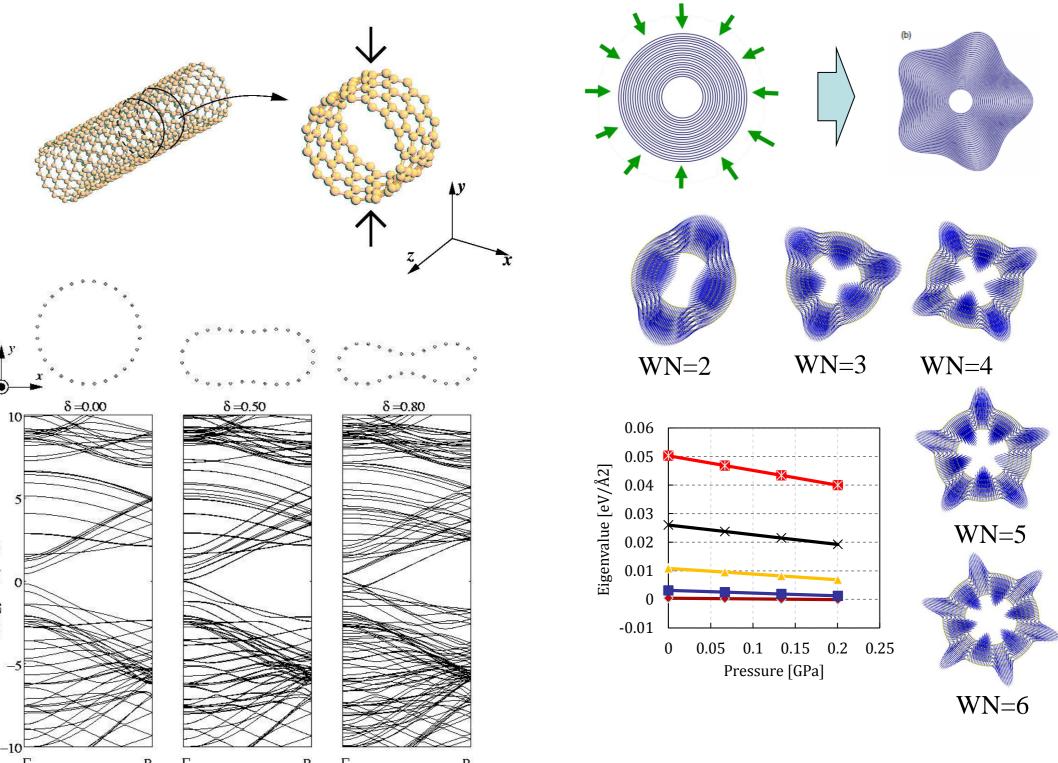
## **Theoretical Prediction of Strength and Physics of Nanomaterials and Multiscale Simulation**

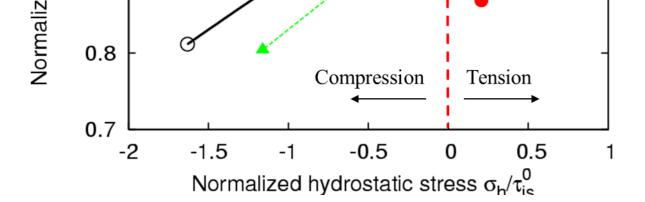
We aim to reveal the mechanical and physical properties of nanomaterials by ab initio density functional theory calculations and molecular dynamics simulations. We also work on multiscale simulation based on knowledge of nano-microscale phenomena.

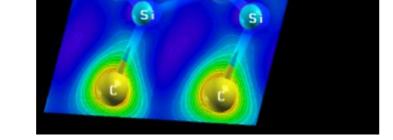




### Compression of carbon nanotubes



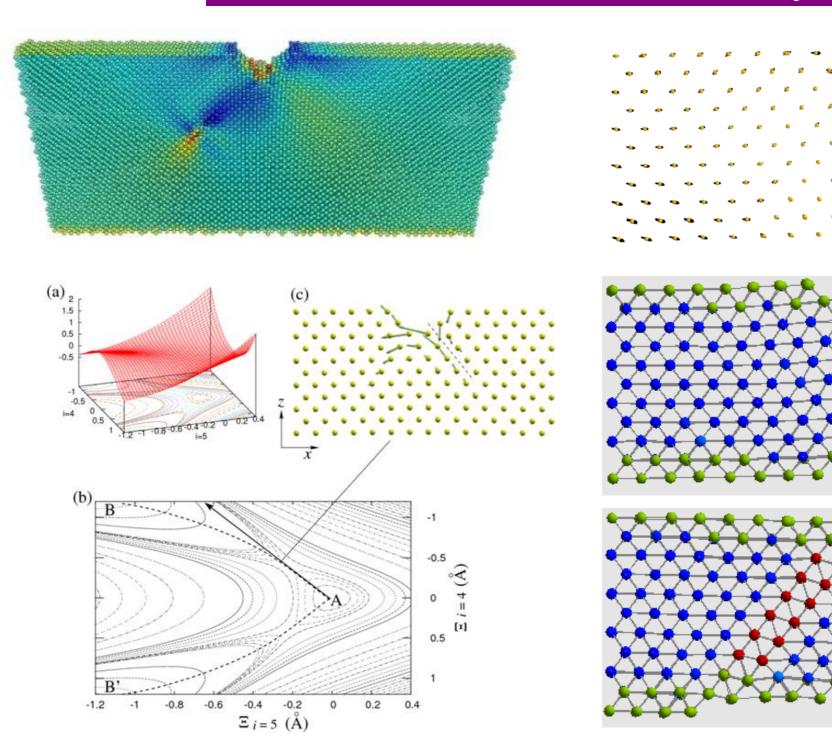




Ideal shear stress (ISS) under compression/tension, which is important to interpret experiments (e.g. nano-indentation tests), has been calculated. Response of ISS qualitatively differs. Note that compression always increases ISS in metals.

Effect of normal stress on ideal shear strength

#### Atomic structure instability analysis



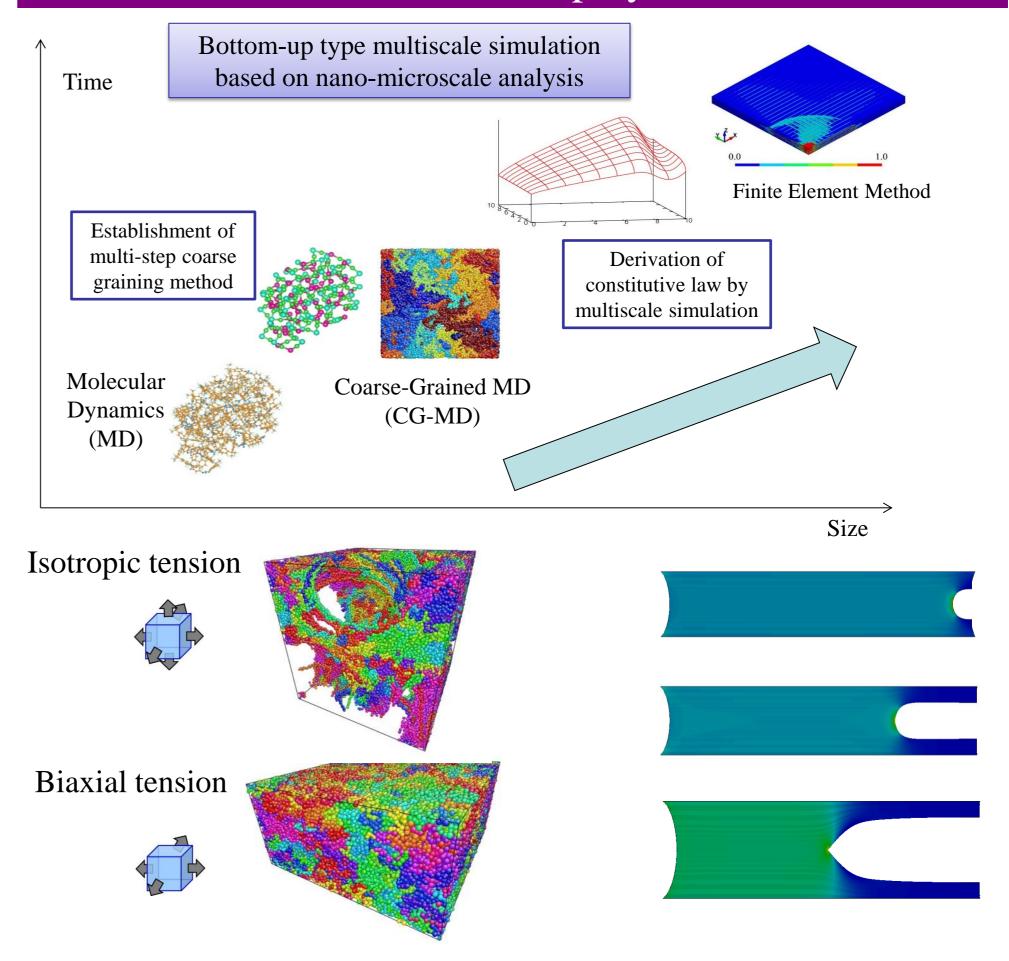
Analysis of band gap energy of carbon nanotubes under radial

compression

**Cw-302** 

Prediction of peculiar radial buckling under pressure (radial corrugation)

#### Multiscale simulation of polymer materials



#### Instability mode analysis of dislocation initiation from defect.





**Institute of Industrial Science**