Contribution ID: **133** Type: **Oral Presentation**

Relative stability of graphene and carbon nanotube structures

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Abstract content
   (Max 300 words)

Abstract.
Molecular dynamics simulations were used to study properties of graphene and single walled carbon nanotube. The formulation of the Tersoff bond-order potential was used to investigate the structural properties and thermodynamics properties of these two carbon polymorphs through a range of temperature. The structural properties were studied using the radial distribution functions effects and the thermodynamics effects were studied using the energy-temperature and volume-temperature plots. Similarities and differences in graphene and carbon nanotubes are discussed.

Apply to be
 considered for a student
   award (Yes / No)?

yes

Level for award
 (Hons, MSc,
   PhD)?

Msc

Main supervisor (name and email)
and his / her institution

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Would you like to
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 Proceedings (Yes / No)?

Yes

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