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Shape transition of unstrained flattest single-walled carbon nanotubes under pressure

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Single walled carbon nanotube’s (SWCNT’s) cross section can be flattened under hydrostatic pressure. One example is the cross section of a single walled carbon nanotube successively deforms from the original round shape to oval shape, then to peanut-like shape. At the transition point of reversible deformation between convex shape and concave shape, the side wall of nanotube is flattest. This flattest tube has many attractive properties. In the present work, an approximate approach is developed to determine the equilibrium shape of this unstrained flattest tube and the curvature distribution of this tube. Our results are in good agreement with recent numerical results, and can be applied to the study of pressure controlled electric properties of single walled carbon nanotubes. The present method can also be used to study other deformed inorganic and organic tube-like structures. © 2014 AIP Publishing LLC. [http://dx.doi.org/10.1063/1.4863455]

I. INTRODUCTION

Carbon nanotubes (CNTs) have superior mechanical properties due to strong carbon-carbon atomic interactions in their honeycomb lattices.1 CNTs’ high elastic modulus, exceptional axial stiffness, and low density, make them ideal for nanotechnology applications.2 CNTs’ mechanical properties are highly anisotropic, and have been studied extensively.3,4 In contrast to the high tensile strength,5 CNTs are susceptible to mechanical distortion in their radial directions under applied hydrostatic pressure on the order of GPa.6–17 The radial deformation controlled by the applied pressure provides an approach to modify the electronic properties of single walled carbon nanotube (SWCNTs),18 consequently, radial deformation of SWCNTs can be observed by optical spectroscopy since the electronic band structure of a SWCNT is sensitive to its morphological transition.4,19 In addition, a SWCNT’s chemical reactivity depends on its mechanical deformations,20–22 which plays the key role in the design of CNT-based gas sensors.23

Among the radially deformed SWCNTs, the fully collapsed structure with two strained edges bridged by a flat middle section24–26 attracts intense interest because of its physical and chemical properties.28–30 Associated with its flat ribbon-like middle part,30 However, the fully collapsed SWCNTs are stabilized by van der Walls (vdW) interaction between two opposing flat walls (with typical interlayer distance $d_0 \approx 3.4 \text{ Å}$) leads to irreversible collapsing. The vdw interaction may also induce twist and bending of the flat section in a fully collapsed SWCNT. Moreover, the edge section of a collapsed SWCNT is highly strained. As it is widely known, with the increase of hydrostatic pressure, a cylindrical SWCNT at first becomes oval-shaped, and then becomes peanut-shaped.15 Between these two shapes, there exists a critical shape under certain pressure which is the unstrained flattest configuration of SWCNTs. The radially deformed SWCNT with a flat section that is similar to fully collapsed SWCNTs have advantages over fully collapsed ones: (1) it can be shifted back to the state with a circular cross section reversibly; (2) there is no twist in the flat section since vdW interaction can be neglected; (3) there is no strain in the two edges, therefore avoiding the strain-induced change of the electronic band structure.

II. MODEL

In the present manuscript, we will theoretically determine the shapes of the unstrained flattest SWCNTs. Although the problem has been studied by molecular dynamical simulations and numerical calculations,15 there still lacks the analytical explicit expressions which can provide a design principle for CNT-based devices. We will accomplish this task by carefully studying the transition of the SWCNT deforming from a convex shape to a concave shape, and giving an analytical expression of the critical shape for an unstrained flattest tube. Based on this analytical expression, we can calculate the critical pressure for the unstrained flattest SWCNTs. We will also discuss the curvature effect for the orbital hybridization of carbon atoms on the sidewall of the SWCNT. The latter is important for studying the absorption of molecules on SWCNTs. Although the deformation of the SWCNT investigated in the present work is in ideal conditions, the results can be used as the theoretical limits for the actual CNT-devices.

The equilibrium shape of deformed SWCNTs is determined by minimizing the free energy under certain constraints. In the present problem, the free energy contains the elastic
energy and the pressure term, \( F_b = E_b + p \Delta V \). Although both bond bending and bond stretching contribute to the elastic energy of SWCNTs, at the energy scale \( \Delta p \cdot V_0 \sim 1 \text{ eV} \approx k_c \) (\( k_c \) is the bending stiffness of SWCNTs\(^{11,33} \)), the bond bending effect predominates.\(^{33} \) The bond-bending energy of a SWCNT can be described by its curvatures\(^{24,33} \).

\[
E_b = \frac{k_c}{2} \int (2H)^2 \, dA + \int K \, dA. \tag{1}
\]

Here, \( H \) and \( K \) are the mean and Gaussian curvature of the surfaces of carbon atoms, and \( k_c = 1.17 \text{ eV} \) (Ref. 33) is consistent with the result of Tersoff et al.\(^{31} \). The expression of free energy can be mapped to 2D. Consider a straight SWCNT with radius \( r(\phi) \), without the inclusion of its two end-caps, the surface of the tube can be described in cylindrical coordinates as, \( r(\phi) = \{ r(\phi) \cos \phi, r(\phi) \sin \phi, l \} \). Here, \( 0 \leq \phi < 2\pi \), and \( 0 \leq l \leq L_0 \), with \( L_0 \) the length of straight tube axis. The origin of \( r \) is arbitrary. The surface’s mean and Gaussian curvature are \( 2H = -(r^2 + 2r'^2 - \rho r^2) / (r^2 + r'^2)^{3/2} \), \( K = 0 \). Comparing with the relative curvature \( k_r = [(r^2 + 2r'^2 - \rho r^2) / (r^2 + r'^2)^{3/2}] \) of a plane curve \( \rho = r(\phi) \), the bending energy can be rewritten as \( E_b = k_c L_0 / 2 \sqrt{k_r^2 ds} \), where \( s \) the arc parameter of boundary curve \( C: r(\phi) \), and line element \( ds = [r^2 + (\rho q / d\\phi)^2]^{1/2} d\\phi \).

The equilibrium shape of a deformed SWCNT is the solution of the 2D variation problem \( \delta^{(1)} F = 0, \delta^{(2)} F > 0 \), with

\[
\delta^{(1)} F = 0, \quad \delta^{(2)} F > 0,
\]

where \( \lambda \) is the Lagrange multiplier, which is introduced to keep the tube circumference at constant \( l_0 \). Distorting the curve \( C \) by a small perturbation \( \psi(s) \) along the normal direction of the curve, the variations in \( \delta^{(1)} F \) are

\[
\begin{align*}
\delta^{(1)} F & = p \int \psi(s) \, ds + \int \lambda \left[ l_0 - \int \psi(s) \, ds \right], \\
\delta^{(2)} F & > 0,
\end{align*}
\]

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\delta^{(1)} F & > 0,
\end{align*}
\]

which can be obtained in a simple way, as shown in Appendix A. Then, the equation for the equilibrium shape of SWCNTs is obtained

\[
p + \frac{k_c}{2} k_r^3 + k_c k_r' + \lambda k_r = 0. \tag{4}
\]

Obviously, there is a special solution corresponding to the tube of circular cross-section, \( k_r = 1 / l_0 \), which implies the necessary condition for maintaining SWCNT’s circular cross section. The initial equation of Eq. (4)

\[
p k_r + \frac{k_c}{2} k_r^3 + \frac{k_c}{2} k_r'^2 + \frac{\lambda}{2} k_r^2 = c_1,
\]

leads to the equation of \( k_r \).
The constants \( c_2(n) \) can be derived as

\[
c_2(n) = \frac{4\sqrt{\pi} \Gamma\left(\frac{3}{4}\right)}{\Gamma\left(\frac{1}{4}\right)} - \frac{4}{3} \left(2F_1\left(\frac{1}{2}, \frac{5}{2}; \frac{3}{2}; \cos^2\left(\frac{3\pi}{4n}\right) \cos^2\left(\frac{3\pi}{4n}\right)\right) - \frac{3\pi}{4n}\right).
\]

Here, \( 2F_1(a, b, c; z) \) is the hypergeometric function.\(^\text{34}\) Some constants \( c_2(n) \) for an unstrained flattest tube with \( C_{24} \) symmetry are listed in Table I. The curvature \( k_r(n) \), \( n \geq 2 \) has the form

\[
k_{r,n}(\theta) = \left(\frac{n c_2(n)}{pr_0}\right) \sin^{2/3} \left[\frac{3}{4} \left(\theta + \frac{5\pi}{6} - \frac{\pi}{n}\right)\right]. \tag{11}\]

In particular, \( c_2(2) = 4.3244, \) predicts the critical shape for the transition from oval-shaped tube to peanut-like shape.

The equilibrium shape of the unstrained flattest SWCNTs with \( C_{24} \) symmetry can be described by parametric equations (see Appendix C for the details)

\[
\begin{align*}
x_2(\theta) &= \frac{2\pi r_0}{c_2} \sin \left(\frac{\theta}{4} + \frac{\pi}{4}\right) \sin^{1/3} \left(\frac{3\theta}{4} + \frac{\pi}{4}\right), \\
y_2(\theta) &= \frac{2\pi r_0}{c_2} \left[\frac{\sin^{2/3}\left(\frac{\pi}{8}\right)}{\sqrt{2}} + \sin \left(\frac{\theta}{4} + \frac{\pi}{4}\right) \sin^{1/3} \left(\frac{3\theta}{4} + \frac{\pi}{4}\right)\right].
\end{align*}
\tag{12}\]

The comparison of the explicit result and the exact result is shown in Fig. 3.

The cross-sectional area of the tube is

\[
S_2 = \frac{1}{2} \left(\frac{\pi r_0}{c_2}\right)^2 \int_{\pi/2}^{\pi} \sin \left(\theta \sin \left(\theta/4 + \pi/4\right)\right) \sin^{1/3} \left(\frac{3\theta}{4} + \frac{\pi}{4}\right) d\theta = 2.449r_0^2. \tag{13}\]

Comparing with the undistorted tube, \( S_0 = \pi r_0^2, \) the geometric constant is \( G = S_2/S_0 = 0.78, \) which is in good accordance with the exact value \( G = 0.8195.\text{15} \)

The curvature distribution characterizes the flatness of the deformed SWCNTs. The curvature distribution provides useful information about the bond hybridization of SWCNTs, which governs the gas-absorbing ability of SWCNTs. We compare the curvature distributions of our approximate result and the exact solution, as shown in Fig. 4. The exact curvature distribution can be calculated self-consistently according to Eqs. (7) and (8). The hybrid orbital of carbon atoms in the SWCNT is sensitive to the local curvature of the tube. For a SWCNT, in the local coordinate of a given carbon atom \( i, \) the three neighbours of the center atom \( i \) have coordinates\textsuperscript{33}

\[
\begin{align*}
x_{ij} &= \cos \theta_j - \frac{(a_0/R)^2}{3} \sin^2 \theta_j, \\
y_{ij} &= \sin \theta_j + \frac{(a_0/R)^2}{6} \sin^2 \theta_j \cos(2\theta_j), & j &= 1, 2, 3, \\
z_{ij} &= \frac{(a_0/R)}{2} \sin^2 \theta_j,
\end{align*}
\tag{14}\]

with \( \theta_1 = 0, \) \( \theta_2 = 2\pi/3, \) and \( \theta_3 = 4\pi/3 \) being the angles between bond curves and the direction of the tube axis, at the position of atom \( i.\text{33} \) Here, \( a_0 = 1.42 \text{ Å} \) is the equilibrium bond length of carbon-carbon bond, \( R \) is the
radius of the SWCNT, \( a_0/R \) characterizes the curvature dependent properties of SWCNTs. The distance between atom \( i \) and their three neighbours can be described as

\[
r_{ij}^0 = 1 - \frac{(a_0/R)^2}{2} \sin^4 \theta_j.
\]

The directions (the direction of symmetry axis) for the three \( sp^2 \) orbits can be written as

\[
\hat{e}_j = \{x_{ij}, y_{ij}, z_{ij}\}/r_{ij}^0.
\]

In the local coordinate system associated with the given carbon atom \( i \), the \( \pi \) orbit and three \( sp^2 \) orbits can be decomposed as,

\[
|\pi\rangle = c_1 |2\pi\rangle + \sqrt{1-c_1^2} |2p_z\rangle,
\]

and

\[
|sp^2\rangle_j = c_2 |2s\rangle + \sqrt{1-c_2^2} \langle x_{ij} |2p_x\rangle, k = x, y, z.
\]

with

\[
c_1 = \sqrt{\frac{6a_0^3}{32-3a_0^2}} \approx 0.2951a_0, \quad c_2 = \sqrt{\frac{32-9a_0^2}{32-3a_0^2}} \approx 0.2951a_0.
\]

Therefore,

\[
|\pi(x_0)\rangle = \frac{\sqrt{3}}{2} x_0 |2\pi\rangle + (1-\frac{3}{32}a_0^2) |2p_z\rangle,
\]

which plays a key role in the electronic and chemical properties of SWCNTs.3,20 For graphene, \( |\pi(x_0 = 0)\rangle = |2p_z\rangle \). By adjusting applied hydrostatic pressure, the \( |\pi\rangle \) at the main part of a SWCNT can be switched between \( |\pi(x_0)\rangle \) and \( |2p_z\rangle \) reversibly.

IV. CONCLUSION

In summary, we have theoretically investigated the unstrained flattest shape of SWCNT under hydrostatic pressure, which can recover to its original circular cross-section after withdrawing the pressure. We find a good approximate solution for the shape of this flattest SWCNT, and theoretically determine the curvature distribution. We also discuss the curvature-dependent hybrid orbital of the SWCNT. The present analytical solution is in good agreement with the exact numerical solution, and it can be used as the design guideline in CNT-based nano-electronic devices. Our approach can be generalized to investigate other inorganic and organic elastic membrane systems,35-37 including the self-assembled polymer materials and colloidal aggregations.38,39 In the actual devices, the deformed SWCNT may be supported by the substrate. The present results provide the theoretical boundary for this CNT-based devices. It is worth noting that the variations in Eq. (3) can also been obtained in other approaches.40,41

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APPENDIX A: THE DETAILS OF VARIATIONAL CALCULATIONS

We will show the details of deriving the variations in Eqs. (3). The curve is shown in Fig. 5.

1. \( \delta \int f ds \)

With the distortion \( \psi(s) \) along the normal direction of the closed boundary curve, the plane boundary curve \( C \) becomes

\[
\vec{r}(s) \rightarrow \vec{r}(s) + \psi(s)\hat{n}(s),
\]

with \( \hat{n} \) the normal direction of the curve \( C : \vec{r}(s) \), and \( s \) the arc parameter. The line element

\[
ds = |\vec{r}'(s)| ds = \left| \frac{d\vec{r}(s)}{ds} + \frac{d(\psi(s)\hat{n}(s))}{ds} \right| ds,
\]

(\ref{A1})

The length of the \( C \) becomes

\[
\hat{e}_y
\]

\[
\hat{e}_x
\]

FIG. 5. A general closed curve.
Here, we have used the relation between the tangent and normal vector of a plane curve
\[
\mathbf{\hat{r}}(s) = \frac{d\mathbf{r}(s)}{ds}, \quad \mathbf{\hat{r}}'(s) = k_r(s)\mathbf{\hat{n}}(s), \quad \mathbf{\hat{n}}(s) = -k_r(s)\mathbf{\hat{r}}(s),
\]
and \( \dot{\ } \) denoted \( d/ds \). Thus, the first order variation of the length of curve \( C \) has the form
\[
\delta(1) \int ds = \int k_r(s)\psi(s)ds.
\]

2. \( \delta^{(1)} \int k_r^2 ds \)

With the distorsion \( \mathbf{\hat{r}}(s) \to \mathbf{\hat{r}}(s) + \psi(s)\mathbf{\hat{n}}(s) \), the \( s \) is not the arc parameter of the deformed curve \( \mathbf{\hat{r}}(s) \), the curvature changes to
\[
k_r \to \frac{|\mathbf{\hat{r}}'(s) \times \mathbf{\hat{r}}''(s)|}{|\mathbf{\hat{r}}'(s)|^3} = k_r + \psi''(s)c - 2k_r^2\psi(s) + \frac{1}{(1 - k_r(\psi(s))^2} + O(\psi^2(s)),
\]
\[
= k_r(s) + k_r^2\psi(s) + \psi''(s) + O(\psi^2(s)).
\]

Here, we have used the relations
\[
\mathbf{\hat{r}}'(s) \to (1 - k_r(s)\psi(s))ds + O(\psi^2(s)),
\]
\[
\mathbf{\hat{r}}''(s) \to (-k_r''(s) - 2k_r\psi'(s))\mathbf{\hat{r}}(s)
\]
\[
+ (k_r(s) + \psi''(s) - k_r^2(\psi(s)))\mathbf{\hat{n}}(s) + O(\psi^2(s)).
\]

Consider
\[
ds \to (1 - k_r(s)\psi(s))ds + O(\psi^2(s)),
\]
we have
\[
k_r^2(s)ds \to k_r^2(s)ds + k_r^2(s)\psi(s)ds + 2k_r(s)\psi''(s)ds
\]
\[
+ O(\psi^2(s)).
\]
The first order variation of \( \int k_r^2(s)ds \) is therefore
\[
\delta(1) \int k_r^2(s)ds = \int ds k_r^2(s)\psi(s) + \int ds k_r(s)\psi''(s),
\]
\[
= \int ds k_r^2(s)\psi(s) + \int ds k_r^2(s)\psi(s).
\]

3. \( \delta^{(1)} \int dA \)

As shown in Fig. 5, \( dA = \hat{\epsilon}_z \cdot |\mathbf{r}' \times \mathbf{r}"|/2 = \hat{\epsilon}_z \cdot |\mathbf{r}(s)\times \mathbf{\hat{r}}(s)|/2s, \) with the distorsion, the \( dA \) changes to
\[
dA \to \frac{\hat{\epsilon}_z}{2} \cdot ds \left[ (1 - 2k_r(s)\psi(s))\mathbf{\hat{r}}(s) \times \mathbf{\hat{r}}(s) + \psi'(s)\mathbf{\hat{r}}(s) \times \mathbf{\hat{n}}(s) + \psi(s)\mathbf{\hat{e}}_z + O(\psi^2(s)) \right].
\]

Thus,
\[
\delta(1) \int dA = \frac{\hat{\epsilon}_z}{2} \int ds \left[ (1 - 2k_r(s)\psi(s))\mathbf{\hat{r}}(s) \times \mathbf{\hat{r}}(s) + \psi'(s)\mathbf{\hat{r}}(s) \times \mathbf{\hat{n}}(s) + \psi(s)\mathbf{\hat{e}}_z + O(\psi^2(s)) \right].
\]
\[ d \tilde{s} = \frac{d \tilde{s}}{d \theta} d \theta = \frac{1}{k_r} d \theta, \]

the equation can be further reduced to

\[ x_2(\theta) = x_2(0) + \int_{\pi/2}^{\theta} d \tilde{\theta} \cos \frac{1}{k_r} \tilde{\theta}(\theta), \]
\[ y_2(\theta) = \int_{\pi/2}^{\theta} d \tilde{\theta} \sin \frac{1}{k_r} \tilde{\theta}(\theta). \]

(C2)

Substitute the explicit expression of \( k_{r2}(\theta) \) to the above equation and integrate, we can obtain Eq. (12).