

Nonlinear nonlocal damped free and forced vibrations of piezoelectric SWCNTs under longitudinal magnetic field due to surface effects using a two steps perturbation method

Abstract

In the present work, damped free and forced vibrations of single-walled piezoelectric carbon nanotubes under longitudinal magnetic field due to surface effects surrounded on a non-linear viscoelastic medium using the nonlocal Euler-Bernoulli beam theory and multiple time scales method are investigated. Lorentz force equation is used to obtain the vertical force due to the applied voltage to the system. The surface effects as well as a combinational non-linear viscoelastic foundation are considered, and finally, the dynamic equilibrium equations are used, and non-linear equations of motion are extracted. In the following, the Galerkin and multiple time scales methods are used, and finally, analytical solutions are extracted as the non-linear free and forced vibrational responses of the system. The relevant coefficients of the extracted analytical solutions are discovered for two both simple support (S-S) and clamped (C-C) boundary conditions. In the following, , and the effects of the different parameters such as non-local parameter as well as electric-magnetic fields, effect of hardness-linear damping parameters of nonlinear considered viscoelastic foundation, applied magnetic field, base modes for different forms considering surface effects, and etc. will be studied. As one the results of this study, the presence of a non-local parameter has increased the curvature deviation to the right and the stiffening effect. In other words, the non-local parameter is a factor to increase the nonlinear effect of the system. Also, it is predictable that as the load affect position moves away from the center of the single-walled piezoelectric carbon nanotube toward the supports, the amplitude of the dynamic response decreases significantly, and this relative reduction is greater for the C-C boundary condition than for the S-S boundary condition. It is also important to note that the location of the load has no effect on the rate of deviation of the curve peak, and the degree of nonlinearity of the vibrational response of the system.

Keywords: nonlocal euler-bernoulli beam theory, non-linear non-local damped forced vibration, viscoelastic piezoelectric carbon nanotubes, magnetic field, non-linear viscoelastic foundation, surface effects

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Introduction

Having unique physical and chemical properties such as extremely high tensile strength and rigidity, combined with very low density, makes CNTs an essential part of all industrial fields. Nanotubes are critical members of the design and fabrication of nanoelectromechanical systems.¹ Microelectromechanical and nanoelectromechanical systems are the essential applications of piezoelectric generators.² Surface effects are other phenomena that occur due to the static balance of atoms on the surface.³ These effects on elastic bodies are divided into three categories: surface tension, residual surface stress, and surface density.⁴ Indeed, surface effects will play a serious role when the surface-to-volume ratio increases.⁵⁻⁶ Finally, surface effects may have an important role in the vibration analysis of nanotubes.⁷⁻⁸

The damped forced vibration of SWCNTs was analyzed using a new shear deformation beam theory.⁹ In this new shear deformation beam theory, there was no need to use any shear correction factor, and also the number of unknown variables was the only one that was similar to the Euler-Bernoulli beam hypothesis. The torque effect of an axial magnetic field on a functionally graded (FG) nano-rod has been studied to capture size effects under magnetic field using Maxwell's relation.¹⁰ As an important result of this study could be mentioned that

an FG nano-rod model based on the nonlocal elasticity theory behaves softer and has a smaller natural frequency. Investigation of free vibration of viscoelastic nanotube under longitudinal magnetic field also was studied in recent years.¹¹ According to a deep investigation on the natural frequencies and effect of different parameters such as the nonlocal parameter, structural damping coefficient, material length scale parameter, and the longitudinal magnetic field, the results of this research may be helpful for understanding the potential applications of nanotubes in Nano-Electromechanical System. Many studies in the literature are focused on the investigation of the vibration analysis of nanostructures in an elastic medium. The vibrational response of a SWCNT member considered in an elastic medium to transport a viscous fluid,¹² Investigating the vibration of a DWCNT embedded in an elastic medium despite initial axial forces,^{13,14} Investigating the vibration of nanotubes in conditions where the surrounding medium is considered elastic,¹⁵ investigating the nonlinear free vibration of a DWCNT member considering the von Karman assumption in order to apply nonlinear effects and nonlinear,¹⁶ nonlinear vibration of a DWCNT by C-C boundary condition in an elastic medium using the nonlinear van der Waals forces, and also the von Kármán geometric to consider nonlinearity,¹⁷ Investigating the forced vibration of a DWCNT with the potential to carry a moving nanoparticle,¹⁸ investigation of the nonlinear vibration of an MWCNT in thermal

environments,¹⁹ vibration analysis of an MWCNT by thermal effects, and also considering the size effects on the large amplitude,²⁰ studying free vibration of SWCNT with elastic effects for different boundary conditions,²¹ vibration of SWCNT with an elastic medium due thermal conditions,²² thermal-mechanical effects consideration to study vibration and buckling instability of a SWCNT that carry out fluid, and also is rested on an elastic foundation,²³ electro-thermo-mechanical effects consideration to study vibration of boron nitride nanorod that is in elastic medium by non-uniform and non-homogeneous properties,²⁴ buckling analysis of SWCNT on a viscoelastic foundation for different boundary conditions,²⁵ buckling investigation of SWCNT under thermal effects that is in an elastic medium with one elasticity parameter,²⁶ and buckling behavior of single-walled CNT considering thermal effect in an elastic medium²⁷ were studied using non-local theory of elasticity. another theory that can be considered as size-dependent continuum theories, and also can be used to study the electro-thermal transverse vibration of CNT in an elastic medium was presented based on the non-local shell theory as well as piezoelectricity properties of system.²⁸ Continuum shell theory by considering nonlocal effects was used to study free vibration of single-walled carbon nanotubes by non-homogenous elastic medium.²⁹ Using multiple elastic models of beams as well as continuum mechanics theory, non-linear free vibration of multi-walled carbon nanotubes was studied.³⁰ Nonlinear vibration as well as thermal stability and of pre/post buckling due to temperature effects, and also size-dependent FG beams designed on an elastic foundation was studied based on the modified couple stress theory.³¹ The multiple time scales method, as a perturbation method, is known as an efficient technique in nonlinear differential equations analysis. Multiple scale method was used to study free and forced vibration of beams as well as DWCNT on an elastic medium by considering geometric nonlinearity and S-S boundary condition.³²⁻³⁴ Non-local elasticity theory was used to study tensioned nanobeam nonlinear vibration with considering various boundary conditions.^{35,36} The non-local Euler-Bernoulli theory of elasticity³⁷ can be considered as a theory to modify nanoscale as well as nanostructures models that greatly has been used in modelling carbon nanotubes due to the size-effects in nanoscale behavior. Several references in the literatures can be found on the basis of the nonlocal theory of elasticity.^{38,39} In some of the recent relevant papers, the nonlinear frequency response of single-walled carbon nanotubes to primary resonance has been studied based on the nonlocal Euler-Bernoulli beam theory.⁴⁰⁻⁴⁸ The nonlocal Euler-Bernoulli elasticity theory also has been greatly used to study Nonlocal instability of cantilever piezoelectric carbon nanotubes by considering surface effects subjected to axial flow,⁴⁹ Linear free vibration analysis of piezoelectric SWCNTs that the results are extracts considering linear part of a new general nonlinear viscoelastic foundation without considering external force,⁵⁰ Vibration analysis of piezoelectric nanowires with surface and small scale effects,⁵¹ Flexoelectric and surface effects on size-dependent flow-induced vibration and instability analysis of fluid-conveying nanotubes based on flexoelectricity beam model,⁵² and Size-dependent nonlinear vibration of functionally graded composite micro-beams reinforced by carbon nanotubes with piezoelectric layers in thermal environments.⁵³ Consequently, due to lots of the presented works in the literature, the objective of the present work is to study damped free and forced vibration of single-walled piezoelectric carbon nanotubes under longitudinal magnetic field considering surface effects resting on a non-linear viscoelastic foundation based on the nonlocal Euler-Bernoulli beam theory as well as multiple time scales and Galerkin methods. A theoretical and numerical study on non-linear nonlocal free and forced vibration responses are performed. This study, considering surface effects, develops

a theoretical response for the non-linear nonlocal vibration of piezoelectric carbon nanotubes located on a non-linear viscoelastic foundation, and attenuation factors due to the external magnetic field and piezoelectric voltage Using Galerkin and multiple time scales methods. Surface effects, which harden the surface of nanotubes, are created by the formation of homogeneous masses by Van forces in the waltz. In the following, considering different boundary conditions of the system (S-S and C-C), the relevant coefficients of the developed analytical response are determined. Finally, the vibrational response parameters such as Amplitude-frequency response curves of nonlinear forced vibration, and etc. will be generally investigated for different boundary conditions, and the effects of the different parameters such as non-local parameter and electric-magnetic fields, Effect of hardness-linear damping parameters of viscoelastic foundation, applied magnetic field, base modes for different forms considering surface effects, and etc. will be studied.

Mathematical modelling and formulations

Scheme of the considered system

Figure 1 shows structural schematics of a piezoelectric single-walled carbon nanotube with an inner diameter d , outer diameter D , thickness h , length L , mass per unit length m , and elastic modulus E , subjected to voltage V and a magnetic field of magnitude H_x . This nanotube, subjected to a harmonic external point load, resting on a nonlinear viscoelastic foundation. Figure 2 shows discrete model of the single walled carbon nanotube by coordinate system.

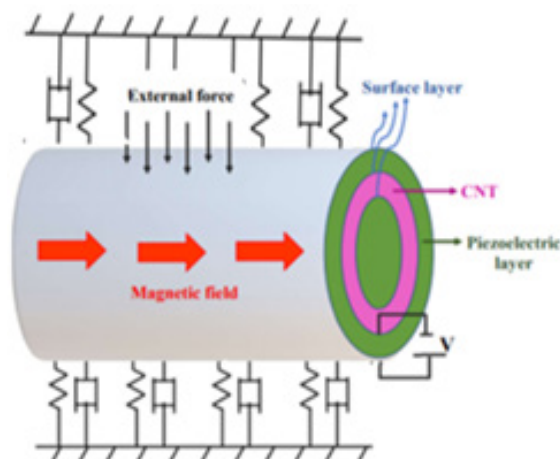


Figure 1 Scheme of the single-walled piezoelectric carbon nanotubes under longitudinal magnetic field subjected to voltage V and resting on nonlinear viscoelastic foundation with surface layer.

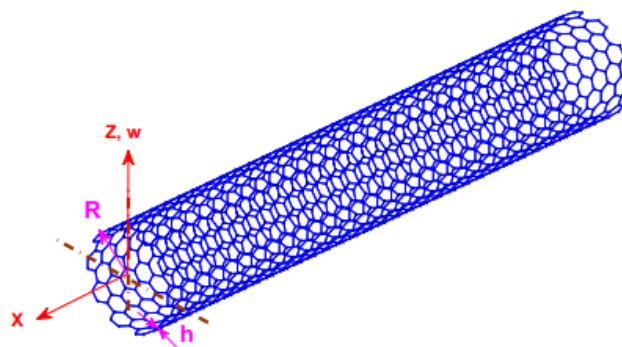


Figure 2 Scheme of the discrete model of the single walled carbon nanotube by coordinate system.

To adequately understand, it is necessary to investigate some fundamental issues and identify the primary mechanism of these advanced systems.

Non-local effects

Generally, nanoscale or microstructure analysis methods can be classified into three categories: molecular simulation, non-classical continuous environment, and multi-scale.⁵⁴ The molecular simulation methods, which simulated atoms and their bonds, investigate nanostructures' behavior more profound than other methods. The high computational volume of this method makes it possible to simulate only small nanostructured dimensions using this method.^{55,56} Molecular methods are divided into quantum mechanics, molecular dynamics, and molecular mechanics.⁵⁷ Complex energy functions are used to explain the motion of elementary particles in quantum mechanics. However, this method is associated with computational limitations due to the complexity of the equations used. In the molecular dynamics method, the motion of atoms is studied concerning their adjacent bonds. Atoms are known as rigid particles that are affected by the potential field of neighbouring atoms. The last one, the molecular mechanical method, is based on the displacement of interatomic bonds with beam and spring elements. Potential energy can express the field of interatomic forces in atomic structures. Equalizing the beam and spring elements' potential energy and strain energy, the mechanical properties are obtained equally. In classical theory, the stress state at a particular point can be calculated if the strain state is known at that point. Whereas, in non-classical theory, such as non-local tension theory, the strain state must be known throughout the range to obtain the stress state at a particular point. All other proposed theories based on non-classical continuum mechanics, including strain gradients and coupled stress theories, relate to two-dimensional elasticity.⁵⁸⁻⁶⁰ Therefore, nonlocal stress tensor can be written as follows:^{61,62}

$$\sigma(X) = \int_V k(|X' - X|, \tau) T(X') dV \quad (1)$$

$$T(X) = C(X) : \varepsilon(X) \quad (2)$$

That $\sigma(X)$ is nonlocal stress tensor at point X , $(|X' - X|, \tau)$ is nonlocal modulus, τ is a material constant, $T(X)$ is classical stress tensor, is $\varepsilon(X)$ strain tensor and $C(X)$ is elasticity tensor. Because of that solving of the integral constitutive Equation (1) is complicated, a simplified differential form is greatly used as follows:

$$T = (1 - \mu \nabla^2) \sigma, \mu = \tau^2 l^2 \quad (3)$$

Where μ is non-local coefficient and is ∇^2 Laplacian operator.

Also, τ and l are material constant.⁶¹ Finally, nonlocal stress tensor for a beam must satisfy equation (4):

$$\sigma(X) - (\tau l)^2 \frac{\partial^2 \sigma(X)}{\partial X^2} = E \varepsilon(X) \quad (4)$$

Where τ is a constant that known as the classical elasticity modulus.

Now, using the existing stress relations in the non-local theory, the relationship between the local bending torque $M^l(x, t)$ and the non-local torque $M^{nl}(x, t)$ at the cross section of the nanotube is written as follows⁶³:

$$\left(1 - (\tau l)^2 \frac{\partial^2}{\partial x^2}\right) M^{nl}(x, t) = M^l(x, t) \quad (5)$$

In equation (5), l is the parameter of the nanotube size scale, which is also known as the non-local parameter.

External force related to Lorentz force and the effect of the magnetic field

Assume $\vec{U} = (u, v, w)$ as the displacement field of the nanostructure; the Lorentz force f_i is equal to⁶⁴:

$$\vec{f}_i = \eta (\vec{J} \times \vec{H}) \quad (6)$$

Where \vec{J} stands for the current density and \vec{H} is the magnetic field strength vector in the environment. Also, in equation (6):

$$\vec{J} = \nabla \times \vec{h} \quad (7)$$

That:

$$\vec{h} = \nabla \times (\vec{U} \times \vec{H}) \quad (8)$$

Therefore, by considering longitudinal magnetic field:

$$\vec{h} = -H_x \left(\frac{\partial v}{\partial y} + \frac{\partial w}{\partial z} \right) \vec{i} + H_x \frac{\partial v}{\partial x} \vec{j} + H_x \frac{\partial w}{\partial x} \vec{k} \quad (9)$$

$$\vec{J} = H_x \left(-\frac{\partial^2 v}{\partial x \partial z} + \frac{\partial^2 w}{\partial x \partial y} \right) \vec{i} - H_x \left(\frac{\partial^2 v}{\partial y \partial z} + \frac{\partial^2 w}{\partial x^2} + \frac{\partial^2 w}{\partial z^2} \right) \vec{j} + H_x \left(\frac{\partial^2 v}{\partial x^2} + \frac{\partial^2 v}{\partial y^2} + \frac{\partial^2 w}{\partial y \partial z} \right) \vec{k} \quad (10)$$

Finally, the Lorentz force is:

$$\vec{f}_i = f_x \vec{i} + f_y \vec{j} + f_z \vec{k} = \eta \left[0 \vec{i} + H_x^2 \left(\frac{\partial^2 v}{\partial x^2} + \frac{\partial^2 v}{\partial y^2} + \frac{\partial^2 w}{\partial y \partial z} \right) \vec{j} + H_x^2 \left(\frac{\partial^2 w}{\partial x^2} + \frac{\partial^2 w}{\partial y^2} + \frac{\partial^2 v}{\partial y \partial z} \right) \vec{k} \right] \quad (11)$$

Therefore, along the x, y, z directions:

$$f_x = 0$$

$$f_y = H_x^2 \left(\frac{\partial^2 v}{\partial x^2} + \frac{\partial^2 v}{\partial y^2} + \frac{\partial^2 w}{\partial y \partial z} \right)$$

$$f_z = H_x^2 \left(\frac{\partial^2 w}{\partial x^2} + \frac{\partial^2 w}{\partial y^2} + \frac{\partial^2 v}{\partial y \partial z} \right), \quad (12)$$

Therefore, considering $w(x, y, z, t) = w(x, t)$, the vertical force affecting the lateral vibration per unit length of the Euler-Bernoulli nanotube is according to the following equation:

$$F_{Lz} = \int_A f_z dA = (\eta A H_x^2) \frac{\partial^2 w}{\partial x^2} \quad (13)$$

Applied force from nonlinear viscoelastic foundation

In this study, vibration analysis of SWCNTs embedded in viscoelastic medium is presented. In this case, it is assumed that the chemical bonds of SWCNT to be formed generally between the external surface of the carbon nanotube and the viscoelastic medium (see Figures 1 and 2). In this study, the present viscoelastic foundation has linear Winkler stiffness, nonlinear stiffness, linear viscosity damping, and the attenuation is nonlinear. Therefore, in the present work, the force coming from the nonlinear viscoelastic medium could be considered in the general form as follows:

$$F_{medium} = F_{linear} + F_{nonlinear} =$$

$$\left(F_{linear_k} + F_{nonlinear_k} \right) + \left(F_{linear_c} + F_{nonlinear_c} \right) =$$

$$\left[\left(k_1 w(x, t) + (k_3 w^3(x, t)) \right) \right] + \left(c_0 + c_2 w^2(x, t) \right) \frac{\partial w(x, t)}{\partial t}, \quad (14)$$

That includes spring element k_1 , nonlinear spring k_3 , linear damper c_0 , and nonlinear damper c_2 .

Piezoelectric equations considering surface effects

One of the characteristics that distinguish nanomaterials is their surface effects. There are two essential and mechanically distinct

surface effects in nanostructures: surface stress and surface elasticity. Atoms on the surface have different bonds than atoms in the bulk, and surface stresses lead to these atoms being at the minimum energy level. Also, in nanostructures, surface elasticity occurs due to the lack of bonded neighbors in surface atoms.^{65,66}

According to Euler-Bernoulli beam theory, the axial and transverse displacement fields can be considered as follows:

$$u(x, y, z, t) = U^0 - z \frac{\partial w}{\partial x}$$

$$w(x, y, z, t) = w(x, t), \quad (15)$$

Therefore, the only nonzero (axial) strain of this beam at any point is expressed by the following equation:

$$\epsilon_{xx} = \frac{\partial U^0}{\partial x} - z \frac{\partial^2 w(x, t)}{\partial x^2} \quad (16)$$

Where ϵ_0 is the axial strain applied to the structure and $w(x, t)$ is also the transverse displacement (bending) at the desired location x on the center plate of the nanotube. Let's suppose E_z is the electric field created by the electric potential Φ assuming that this field exists only in the direction of the z -axis where Z indicates the distance from the center of the arrow, then the relationship between the electric current created and the existing electric potential is:⁶⁷

$$E_z = -\frac{\partial \Phi}{\partial z} \quad (17)$$

For the surface piezoelectric model, the surface-related structural relationships are different from the bulk volume equations. If the direction of polarization of the piezoelectric medium is assumed in the direction of the z -axis, the stress-strain structural equations related to the surface of the piezoelectric carbon nanotube-based on the relationships presented are written as follows:³⁻⁵

$$\sigma_x^s = \sigma_x^0 + c_{11}^s \epsilon_x - e_{31}^s E_z \quad (18)$$

$$D_x^s = D_x^0 \quad (19)$$

In equations (18) and (19), σ_x^s and D_x^s are axial surface stress and electric surface displacement, respectively. The parameters σ_x^0 , D_x^0 , e_{31}^s and c_{11}^s indicate the residual surface stress, the residual electrical displacement, the surface piezoelectric constant, and the elastic surface constant, respectively. On the other hand, in piezoelectric carbon nanotubes, the structural relationship between bulk stress and bulk strain would be obtained as follows:⁶⁸

$$\sigma_x = C_{11} \epsilon_x - e_{31} E_z \quad (20)$$

$$D_z = e_{31} \epsilon_x + k_{33} E_z \quad (21)$$

That σ_x and D_z represent the classical stress tensor component and the electrical displacement, respectively. Also, C_{11} , e_{31} , and k_{33} are the elastic constant, piezoelectric constant, and dielectric constant of nanotubes, respectively.

In the absence of free electric charge $\frac{\partial D_z}{\partial z} = 0$, and the potential energy boundary conditions related to the voltage applied to the external environment of the nanotube is obtained from the following equation:

$$\Phi \left(\frac{D}{2} \right) = V, \quad \Phi \left(-\frac{D}{2} \right) = 0 \quad (22)$$

By placing the relation (16) in equation (18), and then placing them

in the constitutive equations (20) and (21) as well as using Gauss's law, the axial stress equations (in the local area) on the surface and volume of bulk nanotubes in the local space will be written as follows:

$$\sigma_x = C_{11} \epsilon_0 - e_{31} \frac{V}{D} - \left(C_{11} + \frac{e_{31}^2}{k_{33}} \right) z \frac{\partial^2 w(x, t)}{\partial x^2} \quad (23)$$

$$\sigma_x^s = \sigma_x^0 + c_{11}^s \epsilon_0 - e_{31}^s \frac{V}{D} - \left(C_{11} + \frac{e_{31}^s e_{31}}{k_{33}} \right) z \frac{\partial^2 w(x, t)}{\partial x^2} \quad (24)$$

Nonlocal governing differential equation of motion for equivalent continuum structure of embedded SWCNT

The equations of motion for the considered SWCNT, as shown in Figure 1, can be extracted as follows:

$$\left[(V(x, t)) - \left(V(x, t) + \frac{\partial V(x, t)}{\partial x} dx \right) + F_{medium} dx + F_{ext} dx \right] = (\rho A dx) \frac{\partial^2 w(x, t)}{\partial t^2}$$

$$-\frac{\partial V(x, t)}{\partial x} + F_{medium} + F_{ext} = (\rho A) \frac{\partial^2 w(x, t)}{\partial t^2}, \quad (25)$$

And also for the momentum components:

$$F_{va_{eff}} \frac{\partial w(x, t)}{\partial x} + V(x, t) - \frac{\partial M(x, t)}{\partial x} = 0 \quad (26)$$

Where $F_{va_{eff}} = P_{eff} + F_{medium}$; that P_{eff} is the equivalent axial force which includes the force induced in the system due to the application of axial strain, surface electric charge and the electric field created in the whole nanotube structure, $V(x, t)$ is the shear force at the cross section and $M(x, t)$ is the bending moment at the cross section. Also, F_{medium} is the force applied to the system by the viscoelastic foundation and F_{Lz} is the force applied in the nanotube by the magnetic field to the nanotube per unit length of the nanotube. m is also the mass per unit nanotube length. Based on volume and surface stresses of SWCN, the resultant bending moment can be written as follows for a beam model:

$$M = \int_A \sigma_x z dA + \int_S \sigma_x z dS = -EI_{eff} \frac{\partial^2 w(x, t)}{\partial x^2} = \left[\frac{\pi D^4}{64} \left(E + \frac{e_{31}^2}{k_{33}} \right) + \frac{\pi D^3}{8} \left(c_{11}^s + e_{31}^s \frac{e_{31}}{k_{33}} \right) \right] \frac{\partial^2 w(x, t)}{\partial x^2}, \quad (27)$$

where A , S are the circumference and surface area of the circular point of the nanotube, respectively, and EI_{eff} is the equivalent bending stiffness of the whole piezoelectric nanotube. The equivalent axial force to the nanotube due to the application of strain and electric charge on the volume and surface of the nanotube can be considered as follows:³⁻⁵

$$P_{eff} = \frac{\pi D^2}{4} \left(E \epsilon_0 + e_{31} \frac{V}{D} \right) + 2 \left(D(\sigma_x^0 + c_{11}^s \epsilon_x) + e_{31}^s V \right), \quad (28)$$

So, the equation of motion in local theory is developed as follows:

$$\frac{\partial^2 M(x, t)}{\partial x^2} - P_{eff} \frac{\partial^2 w(x, t)}{\partial x^2} - F_{medium} - (\eta A H_x^2) \frac{\partial^2 w}{\partial x^2} = -(\rho A) \frac{\partial^2 w(x, t)}{\partial t^2} \quad (29)$$

To apply the non-local effects, the equation (5) is used. First, equation (5) is rewritten as follows:

$$\frac{\partial^2 M(x, t)}{\partial x^2} = \frac{\partial^2}{\partial x^2} \left(1 - (\tau l)^2 \frac{\partial^2}{\partial x^2} \right) M^{nl}(x, t) \quad (30)$$

$$\text{So: } \frac{\partial^2 M(x, t)}{\partial x^2} = -(\tau l)^2 \frac{\partial^4 M^{nl}(x, t)}{\partial x^4} + \frac{\partial^2 M^{nl}(x, t)}{\partial x^2} \quad (31)$$

Therefore, equation (29) can be rewritten as follows:

$$-(\rho A) \frac{\partial^4 M^w(x,t)}{\partial x^4} + \frac{\partial^2 M^w(x,t)}{\partial x^2} - P_{eff} \frac{\partial^2 w(x,t)}{\partial x^2} - \left[(k_1 w(x,t) + k_3 w^3(x,t)) + (c_0 + c_2 w^2(x,t)) \frac{\partial w(x,t)}{\partial t} \right] - (\eta A H_x^2) \frac{\partial^2 w}{\partial x^2} - (\rho A) \frac{\partial^2 w}{\partial t^2} - F_{ext} - (\tau l)^2 \dots \quad (32)$$

Finally, based on the equations (25) and (26), and for a flexural moment and cross-sectional shear force in non-local theory space, the differential equation of motion of the piezoelectric single-walled carbon nanotube considering non-local effects is developed as follows:

$$EI_{eff} \frac{\partial^4 w}{\partial x^4} + \left[(k_1 w + k_3 w^3) + (c_0 + c_2 w^2) \frac{\partial w}{\partial t} \right] - (\eta A H_x^2) \frac{\partial^2 w}{\partial x^2} - P_{eff} \frac{\partial^2 w}{\partial x^2} + (\rho A) \frac{\partial^2 w}{\partial t^2} - F_{ext} - (\tau l)^2 \left\{ \left[k_1 \frac{\partial^2 w}{\partial x^2} + k_3 \frac{\partial^2 w^3}{\partial x^2} \right] + \left[c_0 \frac{\partial^3 w}{\partial x^2 \partial t} + c_2 \frac{\partial^2}{\partial x^2} \left(\frac{\partial w}{\partial t} w^2 \right) \right] \right\} - \left(\eta A H_x^2 \right) \frac{\partial^4 w}{\partial x^4} - P_{eff} \frac{\partial^4 w}{\partial x^4} + (\rho A) \frac{\partial^4 w}{\partial x^2 \partial t^2} - \frac{\partial^2}{\partial x^2} F_{ext} \} = 0, \quad (33)$$

The boundary conditions of nanotubes are also expressed according to the following relations, as shown in figures 3:

For both side simple support:

$$x = 0, L; w = 0, \frac{\partial^2 w}{\partial x^2} = 0 \quad (34)$$

For both side clamp support:

$$x = 0, L; w = 0, \frac{\partial w}{\partial x} = 0 \quad (35)$$

Next, the equation (33) is investigated by nonlinear equation solving methods, and an analytical solution for nonlinear free and force vibration responses are extracted

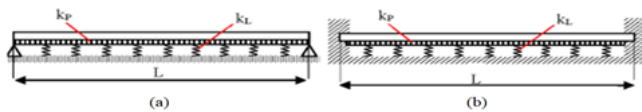


Figure 3 Boundary conditions of the system:

(a) S-S boundary condition

(b) C-C boundary condition.

Analytical solution

Reducing the order of the equations of motion using Galerkin method

The Galerkin method is used to reduce the differential equations obtained. Therefore, the general form of vibrational response is considered as follows:

$$w(x,t) = \sum_{i=1}^{i=N} \varphi_i(x) q_i(t) \quad (36)$$

Where $\varphi_i(x)$ is a mode shape function that satisfies the geometric boundary conditions, and $q_i(t)$ is the generalized time coordinates of the system.

By placing equation (36) in equation (33), and applying the Galerkin method, equation (37) is extracted:

$$\left[\bar{m} \frac{d^2 q_i}{dt^2} + \beta_0 \frac{dq_i}{dt} + \alpha_1 q_i \right] + \left[\beta_2 q_i^2 \frac{dq_i}{dt} + \alpha_3 q_i^3 \right] = \bar{f} e^{i\Omega t} \quad (37)$$

Where:

$$\alpha_1 = EI_{eff} \int_0^L \varphi_i^{(4)} dx + k_1 \int_0^L \varphi_i^2 dx - (F_{vertical}) \int_0^L \varphi_i \varphi_i^{(2)} dx - \mu k_1 \int_0^L \varphi_i \varphi_i^{(2)} dx + \mu (F_{vertical}) \int_0^L \varphi_i \varphi_i^{(4)} dx$$

$$\alpha_3 = k_3 \int_0^L \varphi_i^4 dx - \mu k_3 \left(6 \int_0^L \varphi_i^2 \varphi_i^2 dx + 3 \int_0^L \varphi_i^3 \varphi_i^{(2)} dx \right)$$

$$\beta_0 = c_0 \int_0^L \varphi_i^2 dx - \mu c_0 \int_0^L \varphi_i \varphi_i^{(2)} dx$$

$$\beta_2 = c_2 \int_0^L \varphi_i^3 dx - \mu c_2 \left(6 \int_0^L \varphi_i^2 \varphi_i^2 dx + 3 \int_0^L \varphi_i^3 \varphi_i^{(2)} dx \right)$$

$$\bar{m} = (\rho A) \int_0^L \varphi_i^2 dx - \mu (\rho A) \int_0^L \varphi_i \varphi_i^{(2)} dx \quad (38)$$

Solving nonlinear differential equation (the multiple time scales method)

The multiple time scales method, as a perturbation method, is known as an efficient technique in nonlinear differential equations analysis. In this part, the multiple time scales method is used to develop the nonlinear vibrational response of the system.

Free vibration analysis

To be able to apply this method, equation (37) is rewritten in the following form:

$$\left[\frac{d^2 q_i}{dt^2} + \varepsilon (\beta_0 + \beta_2 q_i^2) \frac{dq_i}{dt} + \omega_n^2 q_i \right] + [\varepsilon (\alpha_3) q_i^3] = 0 \quad (39)$$

In equation (39), ε is a small and dimensionless parameter. In the multiple time scales method, we first define the following parameters and equations:

$$T_n = \varepsilon^n t, \quad \frac{\partial}{\partial T_n} = D_n, \quad n = 0, 1, 2, \dots$$

$$\frac{d}{dt} = (D_0 + \varepsilon D_1 + \varepsilon^2 D_2 + \dots)$$

$$\frac{d^2}{dt^2} = (D_0^2 + 2\varepsilon D_0 D_1 + \varepsilon^2 D_1^2 + \dots) \quad (40)$$

The general form of the response is also expressed as follows:

$$q(t) = q_0(T_0, T_1, T_2, \dots) + \varepsilon q_1(T_0, T_1, T_2, \dots) + \varepsilon^2 q_2(T_0, T_1, T_2, \dots) + \dots = q_0(T_0, T_1, \dots) + \varepsilon q_1(T_0, T_1, \dots) + O(\varepsilon^2) \quad (41)$$

Due to the attenuation in the equations of motion, the amplitude is expanded as the following equation:

$$a(\varepsilon) = a_0 + \varepsilon a_1 + \varepsilon^2 a_2 + \dots \quad (42)$$

Therefore, the nonlinear equation of the system is rewritten as follows:

$$\left[\frac{d^2 q_1}{dt^2} + \varepsilon (\beta_0 + \beta_2 a^2(\varepsilon) q_1^2) \frac{dq_1}{dt} + \omega_n^2 q_1 \right] + \varepsilon [(\alpha_3 a^2(\varepsilon)) q_1^3] = 0 \quad (43)$$

Now, by placing equation (41) in the equation (43) and separating ε by different powers on both sides of the resultant equation, a set of differential equations will be obtained as follows:

$$O(\varepsilon^0): \left[\frac{\partial^2}{\partial T_0^2} + \omega_n^2 \right] q_0(T_0, T_1, T_2) \quad (44)$$

Therefore:

$$q_0 = \cos(\omega_n T_0 + \gamma(T_1, T_2)) \quad (45)$$

Also:

$$O(\varepsilon^1): \left[\frac{\partial^2}{\partial T_0^2} + \omega_n^2 \right] q_1(T_0, T_1, T_2) =$$

$$\left[-\beta_0 \frac{\partial}{\partial T_0} - 2 \frac{\partial^2}{\partial T_0 \partial T_1} \right] q_0(T_0, T_1, T_2) - (a_0^2 q_0^2(T_0, T_1, T_2)) \left[-\alpha_3 q_0(T_0, T_1, T_2) + \beta_2 \frac{\partial}{\partial T_0} q_0(T_0, T_1, T_2) \right], \quad (46)$$

By placing equation (46) in equation (43):

$$\begin{aligned} (D_0^2 + \omega_n^2)q_1 = & -\frac{1}{4}a_0^2\alpha_3 \cos(3\omega_n T_0 + 3\gamma(T_1, T_2)) + \frac{1}{4}\omega_n a_0^2 \beta_2 \sin(3\omega_n T_0 + 3\gamma(T_1, T_2)) \\ & + \frac{8}{4} \left(\frac{\partial}{\partial T_1} \gamma(T_1, T_2) \right) \omega_n \cos(\omega_n T_0 + \gamma(T_1, T_2)) - \frac{3}{4} a_0^2 \alpha_3 \cos(\omega_n T_0 + \gamma(T_1, T_2)) \\ & + (\beta_0 + \frac{1}{4} a_0^2 \beta_2) \omega_n \sin(\omega_n T_0 + \gamma(T_1, T_2)) \end{aligned} \quad (47)$$

The condition for the solvability of the high differential equation is to prevent the formation of very extensive terms in the time response. Therefore, the coefficients of the terms $\sin(\omega_n T_0 + \gamma(T_1, T_2))$ and $\cos(\omega_n T_0 + \gamma(T_1, T_2))$ in equation (47) should be considered equal to zero. This prevents the following terms from appearing in the system response, and does not take the vibrational response of the system to infinity over a long period time.

Therefore, the equations of the system are obtained as follows:

$$8 \left(\frac{\partial}{\partial T_1} \gamma(T_1, T_2) \right) \omega_n - 3a_0^2 \alpha_3 = 0 \quad (48)$$

$$4\beta_0 + a_0^2 \beta_2 = 0 \quad (49)$$

By solving the equations (48) and (49) together, the unknown variables are obtained in the following relation:

$$\gamma(T_1, T_2) = \frac{12\alpha_3 a_0^2}{32\omega_n} T_1 + \gamma_1(T_2) \quad (50)$$

So, the second-order estimation of the response will be obtained as follows $q_1 = \frac{a_0^2 \alpha_3}{32\omega_n^2} \cos(3\omega_n T_0 + 3\gamma(T_1, T_2)) - \frac{a_0^2 \beta_2}{32\omega_n} \sin(3\omega_n T_0 + 3\gamma(T_1, T_2))$ (51)

Higher estimates can be obtained similarly.

Forced vibration analysis (initial resonance)

To investigate forced vibrational response of the system, equation (37) is rewritten as follows:

$$\frac{d^2 q}{dt^2} + \varepsilon(\beta_0 + \beta_2 q^2) \frac{dq}{dt} + \omega_n^2 q + \varepsilon \alpha_3 q^3 = \varepsilon \bar{f} e^{i\Omega t} \quad (52)$$

In equation (52), the excitation frequency for the initial resonance is according to the relation $\Omega = \varepsilon \sigma + \omega_n$, where σ is the parameter of deviation from resonance.

So:

$$(D_0^2 + \omega_n^2)q_0 = 0 \rightarrow q_0 = A(T_1) e^{i T_0 \omega_n} \quad (53)$$

$$(D_0^2 + \omega_n^2)q_1 = -\beta_0 D_0 q_0 - 2D_0 D_1 q_0 - \beta_2 q_0^2 D_0 q_0 - \alpha_3 q_0^3 + \bar{f} e^{i(\sigma T_1 + \omega_n T_0)} \quad (54)$$

By placing equation (53) in equation (54), and to prevent significant time response terms, the coefficient of a term $e^{i T_0 \omega_n}$ must be set equal to zero, resulting is the following equation:

$$-2i\omega_n \frac{\partial A}{\partial T_1} - \left[\bar{A}(\alpha_3 + i\beta_2 \omega_n) A + i\beta_0 \omega_n \right] A + \bar{f} e^{i T_1 \sigma} = 0 \quad (55)$$

Now, assuming the polar form $A = \frac{a}{2} e^{i\gamma}$, and placing it in the equation (55), the following set of equations will be obtained:

$$\beta_0 \omega_n a + \frac{1}{4} \beta_2 \omega_n a^3 - 2\bar{f} \sin(\sigma T_1 - \gamma) = 0 \quad (56)$$

$$\frac{1}{8} \alpha_3 a^3 - \bar{f} \cos(\sigma T_1 - \gamma) = 0 \quad (57)$$

The frequency equation response of the system with the definition of $\theta = \sigma T_1 - \gamma$, and considering that in the case of uniform response,

the changes of a and θ will be zero. Finally, the nonlinear vibrational response equation is presented to the following final form in the form of a closed-form equation:

$$4\omega_n^2 \left[\left(\frac{1}{16} \beta_2 a^3 + \frac{1}{4} \beta_0 a \right)^2 + \left(\frac{1}{16} \frac{\alpha_3 a^3}{\omega_n} \right)^2 \right] - \bar{f}^2 = 0 \quad (58)$$

Numerical results, validations and discussions

Dimensionless form of input and output parameters are considered by the following equations:

$$\begin{aligned} \bar{x} = \frac{x}{L}, \quad \bar{w} = \frac{w}{L}, \quad \bar{f}_{exc} = \frac{f_{exc}}{EI} \\ \bar{EI}_{eff} = \frac{EI_{eff}}{EI}, \quad \bar{P}_{eff} = \frac{P_{eff} L^2}{EI}, \quad \bar{F}_{lz} = \frac{F_{lz} L^2}{EI} \\ K_1 = \frac{k_1 L^4}{EI}, \quad K_3 = \frac{k_3 L^6}{EI} \end{aligned} \quad (59)$$

$$C_0 = \frac{c_0 L^2}{\sqrt{EI m}}, \quad C_2 = \frac{c_2 L^4}{\sqrt{EI m}}$$

$$\mu = \frac{(e_0 a)^2}{L^2}, \quad T = \frac{t}{L^2} \sqrt{\frac{EI}{m}}, \quad \Omega = \omega L^2 \sqrt{\frac{m}{EI}}$$

Validation for the results of the free vibration analysis

In this case, to validate the results with the references,⁶⁹ the cross-sectional area of the nanowire is considered to be a rectangle with a small thickness and height h which its length is L . On the other hand, in this paper, using the classical local theory, the results of free vibration of a nanowire coated with piezoelectric crystals of PZT-5H for two boundary conditions, S-S and C-C, are obtained. Those properties are:

$$\begin{aligned} e_{31} = -6.5; k_{33} = 1.3 \times 10^{-8}; \\ c_{11}^s = 7.56; e_{31}^s = -3 \times 10^{-8}; \\ E = 126 \times 10^9; L = 20h; \end{aligned}$$

In the following figures, the dimensionless frequency of the first mode of nanowire vibration in terms of its cross-sectional height for different values of an external voltage applied in local and non-local theory is compared with the results obtained in local formulation references.⁷⁰ In the S-S boundary condition:

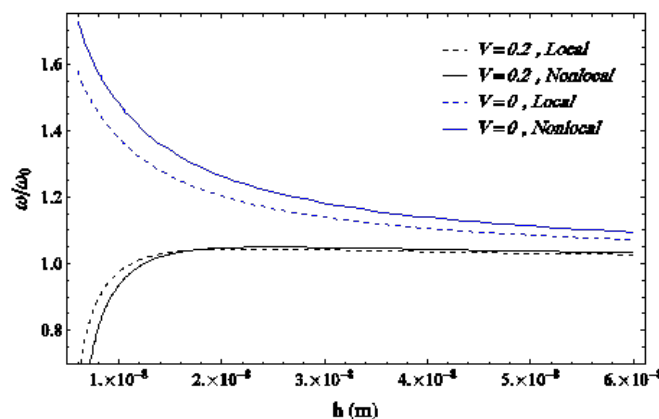


Figure 4 Validation for local⁷⁰ and non-local (present) theory in S-S boundary conditions

And also, for the C-C boundary condition:

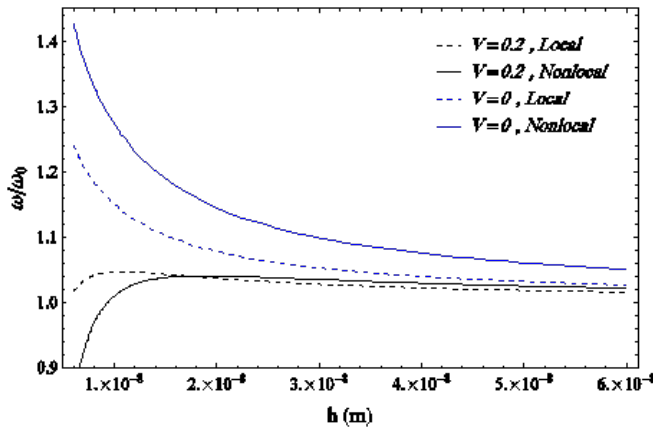


Figure 5 Validation for local⁷⁰ and non-local (present) theory in C-C boundary conditions

Furthermore, for the other forms of the surface effects:

- Form 1: $\sigma_x^0 \neq 0, c_{11}^s \neq 0, e_{31}^s \neq 0$
- Form 2: $\sigma_x^0 = 0, c_{11}^s = 0, e_{31}^s = 0$
- Form 3: $\sigma_x^0 \neq 0, c_{11}^s = 0, e_{31}^s = 0$
- Form 4: $\sigma_x^0 = 0, c_{11}^s \neq 0, e_{31}^s = 0$

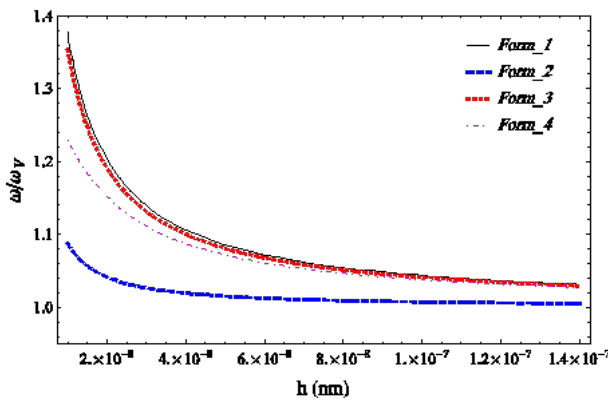


Figure 6 Validation for various forms of surface effects in S-S boundary conditions

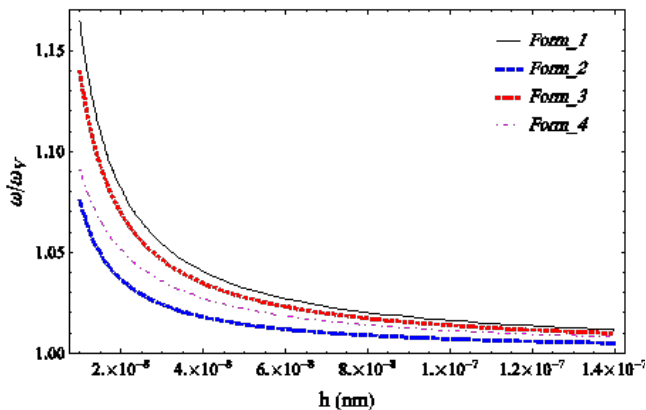


Figure 7 Validation for various forms of surface effects in C-C boundary conditions.

Results of the free vibrational analysis

The effect of non-local parameter and electric-magnetic fields

In the below curves, the changes of dimensionless frequency of the free oscillations at the first mode with two different boundary conditions in terms of the applied voltage and different values of dimensionless non-local parameter, with input parameters $\sigma_x^0 = 1, H_x = 0, C_0 = 0, K_1 = 10$, are shown.

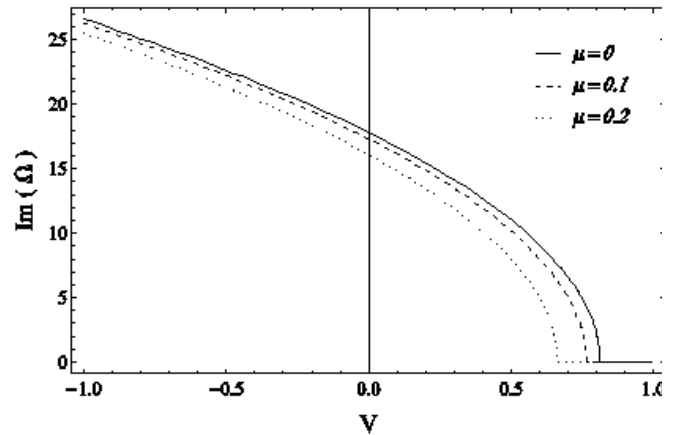


Figure 8 The dimensionless frequency change curve of the free oscillations of the first mode of the system in terms of the applied voltage for different values of the dimensionless non-local parameter of the nanotube, in SS boundary conditions.

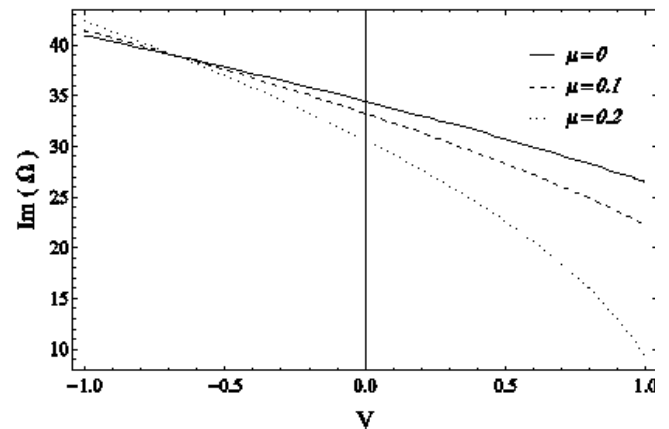


Figure 9 The dimensionless frequency change curve of the free oscillations of the first mode of the system in terms of the applied voltage for different values of the dimensionless non-local parameter of the nanotube, in CC boundary conditions.

In the below curves, the frequency changes of dimensionless free oscillations at the first mode of free vibration of the linear system for two different boundary conditions according to the magnitude of the applied magnetic field in different values of dimensionless non-local parameter, with input parameters $\sigma_x^0 = 1, V = 0.5, C_0 = 0, K_1 = 10$, are shown.

Effect of hardness-linear damping parameters of viscoelastic foundation

In the below curves, the frequency changes of the dimensionless free oscillations at the first mode of the free vibration of the linear system for two different boundary conditions according to the magnitude of the linear stiffness of the foundation in

different forms considering surface effects, with input parameters $\sigma_x^0 = 1, H_x = 1 \times 10^7, \mu = 0.1, C_0 = 0, V = 1$, are shown.

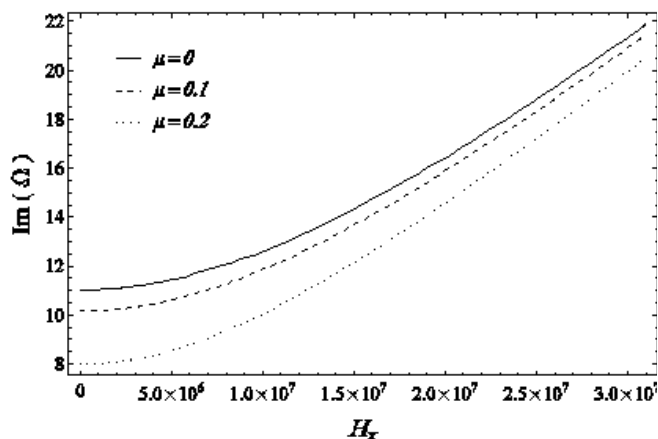


Figure 10 The dimensionless frequency change curve of the free oscillations of the first mode of the system in terms of the magnitude of the magnetic field for different values of the dimensionless non-local parameter of the nanotube, in SS boundary conditions.

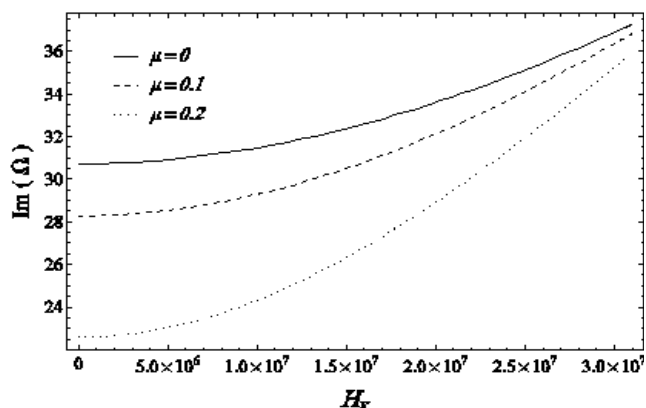


Figure 11 The dimensionless frequency change curve of the free oscillations of the first mode of the system in terms of the magnitude of the magnetic field for different values of the dimensionless non-local parameter of the nanotube, in CC boundary conditions.

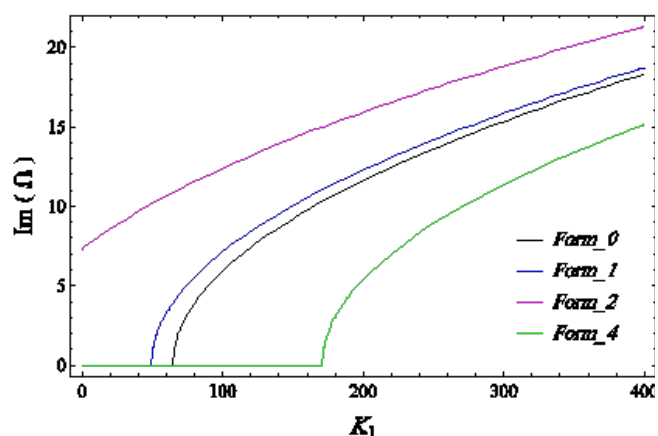


Figure 12 The dimensionless frequency change curve of free oscillations of the first mode of the system in terms of dimensionless linear stiffness of the substrate for different forms considering surface effects, in SS boundary conditions.

In the below curves, the frequency changes of the dimensionless free oscillations at the first mode of the free vibration of the

linear system for two different boundary conditions according to the linear damping coefficient of the foundation in different forms considering surface effects, with input parameters $\sigma_x^0 = 1, H_x = 1 \times 10^7, \mu = 0.1, K_1 = 100, V = 0.5$, are shown.

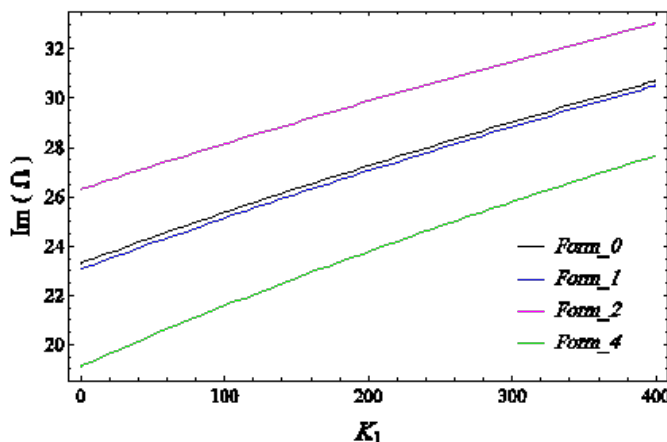


Figure 13 The dimensionless frequency change curve of free oscillations of the first mode of the system in terms of dimensionless linear stiffness of the substrate for different forms considering surface effects, in CC boundary conditions.

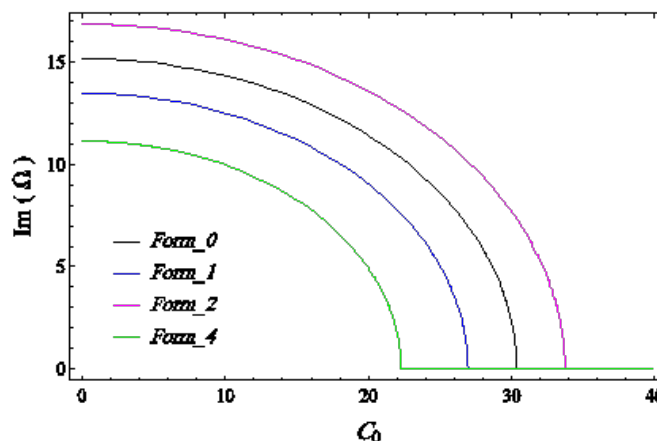


Figure 14 The dimensionless frequency change curve of real and imaginary parts of the special values of the first mode of the system in terms of dimensionless linear damping of the substrate for different forms considering surface effects, in SS boundary conditions.

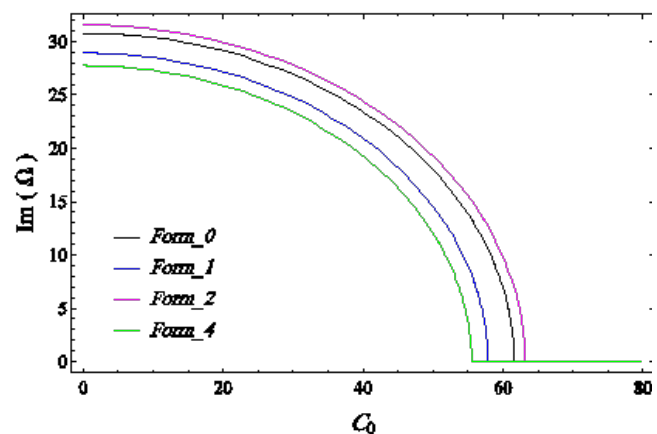


Figure 15 The dimensionless frequency change curve of real and imaginary parts of the special values of the first mode of the system in terms of dimensionless linear damping of the substrate for different forms considering surface effects, in CC boundary conditions.

Frequency results of nonlinear vibration of the non-damped system

In the below curves, the results of the nonlinear vibration for the non-damped system are presented.

For $H_x = 0, \mu = 0.1, K_3 = 100, V = 0.5$:

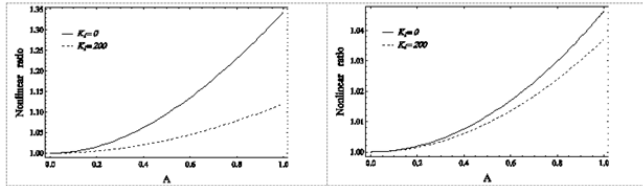


Figure 16 Graph of Nonlinear Frequency to Linear Frequency Ratio of the Immortal System Base Mode for Different Values of the foundation Linear Hardness Parameter, Left: SS boundary condition, right: CC boundary condition.

For $H_x = 0, \mu = 0.1, K_1 = 100, K_3 = 100$:

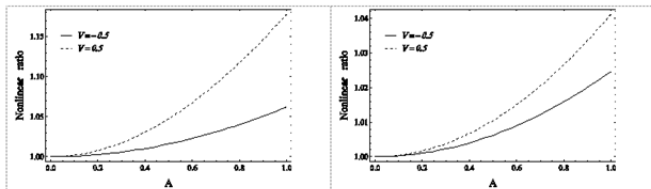


Figure 17 Graph of Nonlinear to Linear Frequency Ratio Changes of the Immortal System Base Mode for Different Values of Applied Voltage, Left: SS Boundary Conditions, Right: CC Boundary Conditions.

For $\mu = 0.1, K_1 = 100, K_3 = 100, V = 0.5$:

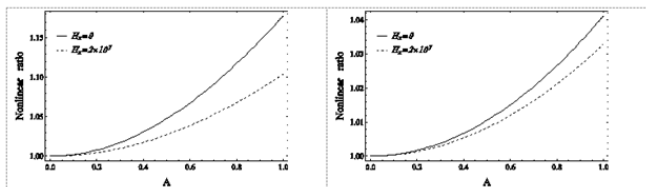


Figure 18 Graph of Nonlinear Frequency to Linear Frequency Ratio of the Immortal System Base Mode for Different Values of Applied Magnetic Field, Left: SS Boundary Conditions, Right: CC Boundary Conditions.

For $H_x = 1 \times 10^7, K_1 = 100, K_3 = 100, V = 0.5$

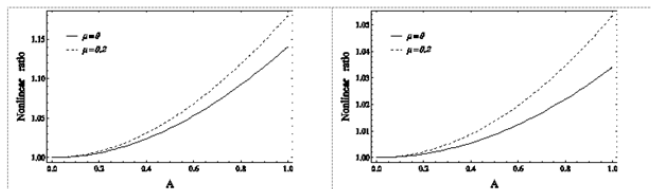


Figure 19 Graph of changes in the ratio of nonlinear to linear frequency ratio of the base mode of the immortal system for different values of the non-local parameter of the nanotube, left: SS boundary condition, right: CC boundary condition.

In the following two diagrams for zero voltage and negative voltage, nonlinear to linear frequency ratio diagrams are drawn for SS boundary conditions:

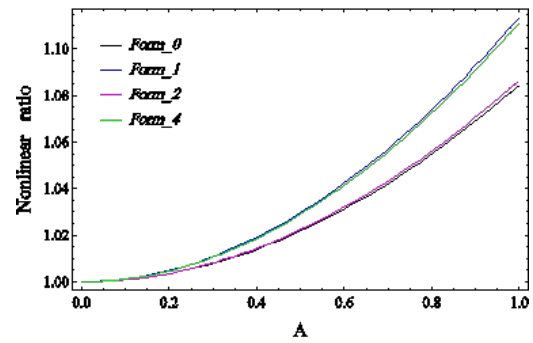


Figure 20 Graph of Nonlinear Frequency to Linear Frequency Ratio of Immortal System Base Mode for Different Forms Considering Surface Effects without Voltage Applied to SS Boundary Conditions.

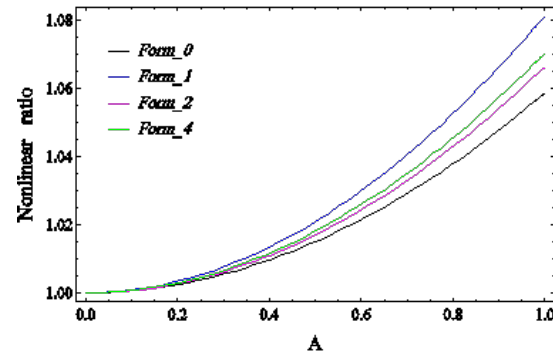


Figure 21 Graph of changes in the ratio of nonlinear to linear frequency ratio of the base mode of the immortal system for different forms Considering surface effects with applied voltage $V = -0.5$, for SS boundary conditions.

Amplitude-frequency response curves of nonlinear forced vibration

For $V = 0, H_x = 0, K_1 = 0, K_3 = 50, C_0 = 0.15, C_2 = 0, \bar{x}_0 = 0.5$:

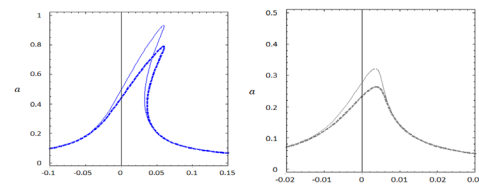


Figure 22 Amplitude-frequency vibration curves for different values of μ ($\mu = 0$ simple line, $\mu = 0.2$ bold fold line); Figure above: SS boundary conditions, bottom figure: CC boundary conditions.

For: $V = 0, H_x = 0, K_1 = 0, K_3 = 50, C_0 = 0.15, \mu = 0.1, \bar{x}_0 = 0.5$:

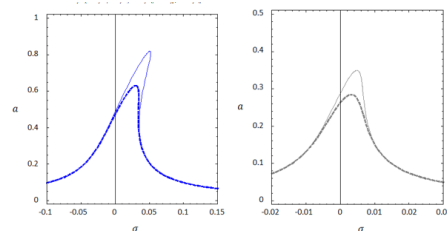


Figure 23 Amplitude-frequency vibration curves for different values of C_2 , ($C_2 = 1$ simple line, $C_2 = 3$ bold fold lines), Figure above: SS boundary conditions, bottom figure: CC boundary conditions.

For: $V = 0, H_x = 0, K_1 = 0, K_3 = 50, C_2 = 1, \mu = 0.1, \bar{x}_0 = 0.5$:

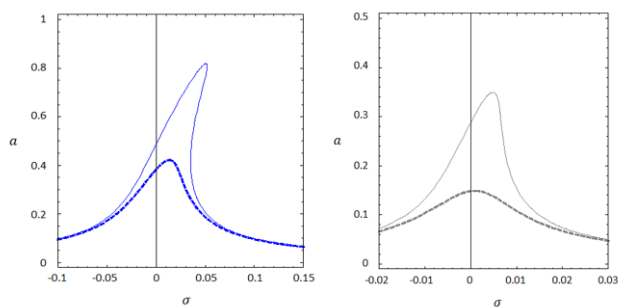


Figure 24 Amplitude-frequency vibration curves for different values of C_0 , ($C_0 = 0.1$ simple line, $C_0 = 0.3$ bold fold line), high: SS boundary conditions, low: CC boundary conditions.

For: $H_x = 0, K_1 = 100, K_3 = 50, C_0 = 0.1, C_2 = 1, \mu = 0.1, \bar{x}_0 = 0.5$:

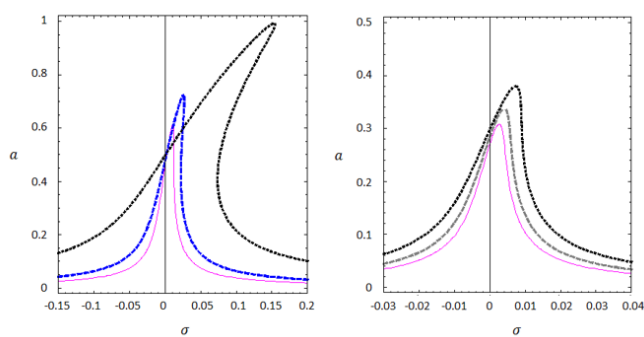


Figure 25 Amplitude-frequency vibration curves for different values of V , ($V = -0.5$ purple line, $V = 0$ blue fold line, $V = 0.5$ bold black), high: SS boundary conditions, low: CC boundary conditions.

According to the presented diagrams, the presence of a non-local parameter has increased the curvature deviation to the right and the stiffening effect. In other words, the non-local parameter is a factor to increase the nonlinear effect of the system. On the other hand, the presence of this factor reduces the maximum range of fluctuations.

Conclusions

By comparing the graphs related to two different boundary conditions, the following results are inferred:

- i. At the C-C boundary conditions, the diagram is collected and the amount of deviation to the right is reduced. For example, the deviation parameters for the maximum dynamic response for the S-S and C-C boundary condition are 0.065 and 0.0045, respectively. In addition, the dynamic response decreases with less degrees of freedom in boundary conditions. For example, in these diagrams, the maximum dynamic range for S-S and C-C are 0.79 and 0.26, respectively. Also, the damping of the system has a greater effect in the boundary conditions of C-C and has significantly reduced the sharpness of the diagram. Therefore, for the C-C boundary conditions, the effect of substrate damping in reducing the dynamic response at excitation frequencies close to the resonant state of the system was greater.
- ii. The higher the foundation attenuation, the lower the amplitude of the amplitude-frequency curve, and consequently the lower the maximum dynamic response. Of course, it should be noted that

the amplitude-excitation frequency curve for a system with C-C boundary conditions is more sensitive to the damping changes of the foundation than the amplitude-frequency curve of the system excitation of the S-S boundary conditions.

- iii. As the nonlinear hardness coefficient of the substrate increases, the amplitude-frequency curve deviates to the right, but the maximum dynamic response amplitude remains constant. By looking more closely at the two diagrams, it could be found that the effect of this parameter on deviating of the diagram is greater for S-S boundary conditions due to more flexibility, so the dynamic response in the CC boundary condition is less sensitive to the increasing nonlinear foundation stiffness parameter.
- iv. According to the obtained results, increasing the linear stiffness of the substrate has led to a decrease in the dynamic response as well as the peak deviation of the curve. Increasing the linear stiffness of the system reduces the nonlinear effects of the system. In addition, in the C-C boundary conditions, where the overall stiffness of the system is higher, this factor is less effective.
- v. The presence of a magnetic factor, in addition to reducing the amplitude of the dynamic response, has led to a decrease in the amplitude of the peak and also a reduction in its curvature to the right. However, these effects have attracted more attention for S-S boundary conditions, which have a more flexible system than for C-C boundary conditions.
- vi. According to the curves, it is quite clear that by changing the applied voltage from zero to a positive value, the response amplitude has increased sharply and the deviation to the right has increased as well. However, by changing the voltage from zero to a negative value, both the dynamic response amplitude and the peak deviation of the graph are reduced as well as nonlinear effects of the system. It is also important to note that under the same conditions these changes for C-C boundary conditions are much less than for S-S boundary conditions and in general it could be said that the sensitivity of the dynamic response curve of the system with C-C boundary condition to most investigated changes in the system input parameters such as foundation stiffness, voltage, and magnetic factor, non-local factor and surface effects are less than the sensitivity of the system with SS boundary conditions.
- vii. It could be seen that by considering the parameter of residual surface stresses for the system, the response amplitude, and the amount of curvature deviation to the right are reduced.
- viii. It is predictable that as the load affect position moves away from the center of the wire toward the supports, the amplitude of the dynamic response decreases significantly, and this relative reduction is greater for the C-C boundary condition than for the S-S boundary condition. It is also important to note that the location of the load has no effect on the rate of deviation of the curve peak and the degree of nonlinearity of the dynamic response of the system.
- ix. Given that the voltage applied to the system is positive for the last two numerical diagrams. It is expected that the graph with the highest deviation and the lowest peak amplitude is related to Form 4 and the graph with the lowest deviation and the lowest peak amplitude is also related to Form 2. In addition, in this case, by comparing the two curves of Form 1 and the main form, it can be concluded that by considering the surface parameters (Form 1), the maximum peak amplitude and the deviation of the graph have increased.

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Conflicts of interest

The authors declare that they have no competing interests.

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