

**Vibration reduction evaluation of a linear system with
a nonlinear energy sink under a harmonic
and random excitation***

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Abstract The nonlinear behaviors and vibration reduction of a linear system with a nonlinear energy sink (NES) are investigated. The linear system is excited by a harmonic and random base excitation, consisting of a mass block, a linear spring, and a linear viscous damper. The NES is composed of a mass block, a linear viscous damper, and a spring with ideal cubic nonlinear stiffness. Based on the generalized harmonic function method, the steady-state Fokker-Planck-Kolmogorov equation is presented to reveal the response of the system. The path integral method based on the Gauss-Legendre polynomial is used to achieve the numerical solutions. The performance of vibration reduction is evaluated by the displacement and velocity transition probability densities, the transmissibility transition probability density, and the percentage of the energy absorption transition probability density of the linear oscillator. The sensitivity of the parameters is analyzed for varying the nonlinear stiffness coefficient and the damper ratio. The investigation illustrates that a linear system with NES can also realize great vibration reduction under harmonic and random base excitations and random bifurcation may appear under different parameters, which will affect the stability of the system.

Key words nonlinear energy sink (NES), Gauss-Legendre polynomial, transmissibility, percentage of energy absorption

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1 Introduction

Malatkar and Nayfeh^[1] first proposed that adding nonlinearity to a traditional linear vibration reduction device could effectively increase the bandwidth of the vibration suppression and improve the robustness of the vibration reduction system. Recently, the principle of targeted energy transfer has been prospering as a method of research to achieve the energy transfer from a linear oscillator to a nonlinear oscillator, i.e., the nonlinear energy sink (NES). It is a representative passive vibration reduction device, which is able to achieve the targeted energy transfer. The NES is an effective device to reduce vibration passively^[2-3].

In the past years, the steady-state response of multi-degree-of-freedom systems under harmonic excitations has been predicted by numerical simulation methods^[4-10] and approximate analytical methods^[4,11-15], e.g., the complexification averaging method^[4,6,8,11,13-18] and the harmonic balance method^[19-22].

Currently, the vibration reduction efficiency of the linear system coupled with an NES system under complex external excitations is a hot research topic for scholars at home and abroad. Gendelman et al.^[23] investigated the dynamic response of a small mass with a nonlinear damper under impact loading. Kerschen et al.^[24-25] studied the vibration reduction mechanism of a two-degree-of-freedom system with an NES. Tsakirtzi et al.^[26] investigated the dynamic behavior of the multi-degree-of-freedom system coupled with an NES. Shiroky and Gendelman^[27] investigated the dynamic behavior of a single degree-of-freedom system with an NES under parametric excitations. Starosvetsky and Gendelman^[28] verified the vibration reduction effect of a single degree-of-freedom system coupled with an NES under narrow-band random excitations with a numerical method. Xiong et al.^[29] simply sketched out the response regimes of a single degree-of-freedom system coupled with an NES under a narrow-band random excitation. Huang et al.^[30] analyzed the steady-state response, random jump, and bifurcation phenomenon of the Duffing system under harmonic and white noise excitations. Yan et al.^[31] studied the dynamic behavior of an axially moving beam excited by harmonic and parametric excitations with the Galerkin truncation method. Zhao et al.^[32] studied the vehicle random vibration with the pseudo-excitation method. Su et al.^[33] used the explicit time-domain method to analyze a coupled vehicle-bridge system.

However, the excitations in the aforementioned studies are the most common random ones, and no efficiency of vibration reduction under the excitations has been interpreted sufficiently and intuitively. In view of the above problems, in this paper, the nonlinear behaviors and vibration reduction of a linear system with an NES are investigated. The paper is arranged as follows. In Section 2, a simple model of a linear system with an NES is presented, and the steady-state Fokker-Planck-Kolmogorov equation is determined by the method of generalized harmonic function. In Section 3, the path integral method based on the Gauss-Legendre polynomial is used to calculate the numerical solutions of the steady-state Fokker-Planck-Kolmogorov equation. In Section 4, the responses and evaluation of vibration reduction are simulated under some given parameters. In Section 5, the nonlinear stiffness coefficient and damper ratio are proven to have a significant effect on the vibration reduction. At last, in Section 6, the main conclusions are summarized.

2 A linear single degree-of-freedom system coupled with an NES

Consider a linear system with an NES under a harmonic and random excitation (see Fig. 1). The model consists of a mass block m_1 , a linear stiffness k_1 , and a linear viscous damper c_1 . The harmonic and random excitation is $u(t) = A_u \sin(\omega t) + W(t)$. The NES is composed of a mass block m_2 , a spring with an ideal cubic nonlinear stiffness k_2 , and a linear viscous damper c_2 .

The model moves in the horizontal direction, and the effects of gravity and friction are

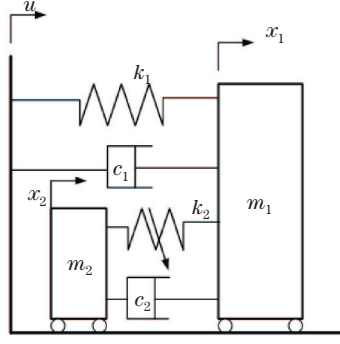


Fig. 1 A single degree-of-freedom system coupled with an NES

neglected. According to the second Newton's law, the governing dynamic equations are as follows:

$$\begin{cases} m_1 \ddot{x}_1 + c_1 \dot{x}_1 + k_1 x_1 + c_2 (\dot{x}_1 - \dot{x}_2) + k_2 (x_1 - x_2)^3 = c_1 \dot{u} + k_1 u, \\ m_2 \ddot{x}_2 + c_2 (\dot{x}_2 - \dot{x}_1) + k_2 (x_2 - x_1)^3 = 0, \end{cases} \quad (1)$$

where x_1 and x_2 are the displacements of the mass blocks m_1 and m_2 , respectively.

The dimensionless form of Eq. (1) can be obtained as follows:

$$\begin{cases} \ddot{y}_1 + \zeta_1 \dot{y}_1 + y_1 + \zeta_2 (\dot{y}_1 - \dot{y}_2) + \beta (y_1 - y_2)^3 = f_t \sin(\gamma \tau) + \xi(t), \\ \varepsilon \ddot{y}_2 + \zeta_2 (\dot{y}_2 - \dot{y}_1) + \beta (y_2 - y_1)^3 = 0, \end{cases} \quad (2)$$

where y_1 is the dimensionless displacement of the mass block m_1 , y_2 is the dimensionless displacement of the mass block m_2 , τ is the dimensionless time, and

$$y_1 = \frac{x_1}{l}, \quad y_2 = \frac{x_2}{l}, \quad u_0 = \frac{u}{l}, \quad \tau = \omega_0 t. \quad (3)$$

In the above equations, u_0 is the dimensionless displacement of the base, l represents the elongation or compression when the linear spring is subjected to an external force of 1 kN. The natural frequency of the linear oscillator ω_0 , the mass ratio ε , the damper ratios of the linear oscillator ζ_1 and the NES oscillator ζ_2 , the ideal cubic nonlinear stiffness β , the amplitude of the harmonic excitation f_t , and the frequency ratio γ are expressed as follows:

$$\begin{cases} \omega_0 = \sqrt{\frac{k_1}{m_1}}, \quad \varepsilon = \frac{m_2}{m_1}, \quad \zeta_1 = \frac{c_1}{m_2 \omega_0}, \\ \zeta_2 = \frac{c_2}{m_2 \omega_0}, \quad \beta = \frac{k_2 l^2}{m_1 \omega_0^2}, \quad f_t = \frac{A_u}{k_1 l}, \quad \gamma = \frac{\omega}{\omega_0}. \end{cases} \quad (4)$$

The random excitation is concretized with the Gaussian white noise, whose strength is two-dimensional, and the autocorrelation function is $E(\xi(t)\xi(t+\tau)) = \delta(\tau)$ in which $\delta(\tau)$ is the Dirac delta function.

Some new state variables are introduced as follows:

$$q_1 = y_1, \quad q_2 = y_1 - y_2, \quad p_1 = \dot{q}_1, \quad p_2 = \dot{q}_2. \quad (5)$$

Substitute Eq. (5) into Eq. (2). Then, after applying a few simple algebraic transformations, we can convert Eq. (2) into the following state-space equation:

$$\dot{q}_i + g_i(q_i) = -\varepsilon h_i(p_i; q_i) + \xi(t). \quad (6)$$

Based on the generalized harmonic function method, the solution to Eq. (6) can be assumed to be as follows:

$$q_i(t) = A_i(t) \cos \phi_i(t), \quad p_i(t) = -A_i(t) v_i(A_i, \phi_i) \sin \phi_i(t), \quad \phi_i(t) = \alpha_i(t) + \psi_i(t), \quad (7)$$

where $v_i = (A_i, \phi_i)$ is the instantaneous angular frequency of the i th oscillator. It converts a function of the generalized displacement and velocity into a function of magnitude and phase as follows:

$$v_i(A_i, \phi_i) = \frac{d\alpha_i}{dt} = \sqrt{\frac{2(U_i(A_i) - U_i(A_i \cos \phi_i))}{A_i^2 \sin^2 \phi_i}}. \quad (8)$$

Substitute Eq. (7) into Eq. (6). Then, we can obtain the standard random Itô differential equation as follows:

$$\begin{cases} \frac{dA_1}{dt} = F_{11}(A_1, \phi_1, \omega t) + H_{11}\xi(t), & \frac{dA_2}{dt} = F_{21}(A_2, \phi_2, \omega t) + H_{21}\xi(t), \\ \frac{d\phi_1}{dt} = F_{12}(A_1, \phi_1, \omega t) + H_{12}\xi(t), & \frac{d\phi_2}{dt} = F_{22}(A_2, \phi_2, \omega t) + H_{22}\xi(t), \end{cases} \quad (9)$$

where

$$\begin{cases} F_{11} = -\frac{\varepsilon A_1}{g_1(A_1)} h_1 v_1 \sin \phi_1, & H_{11} = -\frac{A_1}{g_1(A_1)} v_1 \sin \phi_1, \\ F_{21} = -\frac{\varepsilon A_2}{g_2(A_2)} h_2 v_2 \sin \phi_2, & H_{21} = -\frac{A_2}{g_2(A_2)} v_2 \sin \phi_2, \\ F_{12} = -\frac{\varepsilon}{g_1(A_1)} h_1 v_1 \cos \phi_1, & H_{12} = -\frac{1}{g_1(A_1)} v_1 \cos \phi_1, \\ F_{22} = -\frac{\varepsilon}{g_2(A_2)} h_2 v_2 \cos \phi_2, & H_{22} = -\frac{1}{g_2(A_2)} v_2 \cos \phi_2. \end{cases} \quad (10)$$

According to the Itô differential law, since the amplitude A_i is the slow variable and the phase ϕ_i is the fast variable, if the phases are averaged over time, we can denote the deterministic random average Itô differential equation for amplitude and phase difference as follows:

$$\begin{cases} dA_1 = \bar{m}_{11}(A_1, A_2, \Theta)dt + \bar{\sigma}_{11}dB(t), \\ dA_2 = \bar{m}_{21}(A_1, A_2, \Theta)dt + \bar{\sigma}_{21}dB(t), \\ d\Theta = \bar{m}_{12}(A_1, A_2, \Theta)dt + \bar{\sigma}_{12}dB(t), \end{cases} \quad (11)$$

where \bar{m}_{11} , \bar{m}_{21} , and \bar{m}_{12} are drift coefficients, and $\bar{\sigma}_{11}$, $\bar{\sigma}_{21}$, and $\bar{\sigma}_{12}$ are diffusion coefficients. $B(t)$ is a standard Wiener process.

The phase difference can be expressed as follows:

$$\Theta = \varepsilon \sigma \tau_1 - \phi_1. \quad (12)$$

The averaged drift coefficients can be expressed as follows:

$$\begin{cases} \bar{m}_{11} = F_{11} + D\left(H_{11} \frac{\partial H_{11}}{\partial A_1} + H_{12} \frac{\partial H_{11}}{\partial \phi_1}\right), \\ \bar{m}_{21} = F_{21} + D\left(H_{21} \frac{\partial H_{21}}{\partial A_2} + H_{22} \frac{\partial H_{21}}{\partial \phi_2}\right), \\ \bar{m}_{12} = F_{12} + D\left(H_{11} \frac{\partial H_{12}}{\partial A_1} + H_{12} \frac{\partial H_{12}}{\partial \phi_1}\right). \end{cases} \quad (13)$$

The averaged diffusion coefficients can be expressed as follows:

$$\begin{cases} \bar{b}_{111} = 2DH_{11}^2, \\ \bar{b}_{211} = 2DH_{21}^2, \\ \bar{b}_{112} = 2DH_{11}H_{12}, \\ \bar{b}_{122} = 2DH_{12}^2. \end{cases}$$

According to Eq. (11), the steady-state Fokker-Planck-Kolmogorov equation can be established as follows:

$$\begin{aligned} \frac{\partial p}{\partial t} = & -\frac{\partial(\bar{m}_{11}p)}{\partial A_1} - \frac{\partial(\bar{m}_{21}p)}{\partial A_2} - \frac{\partial(\bar{m}_{12}p)}{\partial \Theta} \\ & + \frac{1}{2} \frac{\partial^2(\bar{b}_{111}p)}{\partial A_1^2} + \frac{1}{2} \frac{\partial^2(\bar{b}_{211}p)}{\partial A_2^2} + \frac{1}{2} \frac{\partial^2(\bar{b}_{122}p)}{\partial \Theta^2} = 0, \end{aligned} \quad (14)$$

where A_1 , A_2 , and Θ constitute a three-dimensional (3D) state vector $X(t)$, m_{11} , m_{21} , and m_{12} are 3D averaged drift coefficient function vectors, and b_{111} , b_{211} , and b_{122} are averaged diffusion coefficient matrix functions.

3 Solving the steady-state Fokker-Planck-Kolmogorov equation

It is known that for the general linear systems and some specific single degree-of-freedom nonlinear systems, e.g., the Duffing system and the Van der Pol system, it is easy to get the exact stationary solutions to a transition probability density function; but for complex multi-degree-of-freedom nonlinear systems, it is difficult to get the exact stationary solutions. The transition probability density function can be used to indicate the probability of a stochastic variable appearing at a certain time after a period of time. In terms of vibration reduction, if the transition probability density of the system at the equilibrium position is larger, the performance of vibration reduction will be better. The approximate numerical solution to a transition probability density function can only be obtained by a numerical method, e.g., the finite difference method^[34], the cell mapping method^[37], and the finite element method^[35–36]. In this paper, a new revised path integral method is adopted to calculate the numerical solution to the steady-state Fokker-Planck-Kolmogorov equation of our proposed system.

The revised path integral method replaces some complex interpolation integral functions with some discrete Gaussian integration points, and enhances the computational efficiency in solving the high-dimensional high-order partial differential equations without boundary singularity.

In a Markov process, the following equation, called the transition probability density function, can be used to describe the probability of the stochastic variables passing a period of time

s at the moment t :

$$P(X(t+s) = j | X(t) = i) = P_{ij}(t, t+s). \quad (15)$$

Then, Eq. (14) can be rewritten as follows:

$$\begin{aligned} & \frac{\partial p(X, t | X^0, t_0)}{\partial t} \\ = & - \frac{\partial(\bar{m}_{ij}(X, t | X^0, t_0)p(X, t | X^0, t_0))}{\partial x_i} + \frac{1}{2} \frac{\partial(\bar{b}_{ijk}(X, t | X^0, t_0)p(X, t | X^0, t_0))}{\partial x_i \partial x_j}. \end{aligned} \quad (16)$$

Given an initial probability condition and some appropriate boundary conditions as follows:

$$\lim_{t \rightarrow 0} p(X, t | X^0, t_0) = \delta(X - X^0), \quad (17)$$

$$p(X_i, t) \rightarrow 0 \quad \text{as} \quad \{X\} \rightarrow \pm\infty, \quad (18)$$

the transition probability density function can be used to completely define the statistical steady-state solution.

According to the initial probability condition and the appropriate boundary conditions, the solution to the transition probability density function can be expressed as follows:

$$p(X, t) = \int_{\Omega} p(X, t | X^0, t_0) p(X^0, t_0) dX, \quad (19)$$

where Ω is the range of the state vector for $X(t)$. The interval $[t_0, t]$ is divided into M sub-intervals, and the probability density function can be derived as follows:

$$\begin{aligned} p(X, t) = & \int_{\Omega} p(X, t | X^{M-1}, t_{M-1}) dX^{M-1} \times \int_{\Omega} p(X^{M-1}, t_{M-1} | X^{M-2}, t_{M-2}) dX^{M-2} \times \dots \\ & \times \int_{\Omega} p(X^2, t_2 | X^1, t_1) dX^1 \times \int_{\Omega} p(X^1, t_1 | X^0, t_0) dX^0. \end{aligned} \quad (20)$$

The integral equation (20) can be discretized at Gauss-Legendre orthogonal points and embodied as follows:

$$p(x^i, t_i) = \sum_{k=1}^K \frac{z_k}{2} \sum_{l=1}^L c_{kl} p(x_{kl}^{(i-1)}, t_{i-1}) p(x^i, t_i | x_{kl}^{(i-1)}, t_{i-1}), \quad (21)$$

where K is the quantity of the sub-intervals, L is the quantity of the Gauss-Legendre orthogonal points in sub-intervals, z_k is the length of the sub-intervals, each x_{kl} is a location of a Gauss-Legendre orthogonal point, and c_{kl} is the corresponding weight coefficient. According to the initial conditions and boundary conditions, Eq. (21) can be used to calculate the transition probability density function of some point in a certain time point. Given some point in a previous time point, it is convenient to calculate some point in a next time point. However, only the following transition probability density function at the Gauss-Legendre orthogonal points is essential:

$$p(x_{mn}^i, t_i) = \sum_{k=1}^K \frac{z_k}{2} \sum_{l=1}^L c_{kl} p(x_{kl}^{(i-1)}, t_{i-1}) p(x_{mn}^i, t_i | x_{kl}^{(i-1)}, t_{i-1}), \quad (22)$$

where $m = 1, 2, \dots, K$, and $n = 1, 2, \dots, L$.

It is worth observing that it is generally assumed that the short-term probability density function is approximately Gaussian, but the approximation of the short-term transition probability density function is different at any Gauss-Legendre orthogonal points. Sun and Hsu^[38] proposed to use the moment equations to derive the first moment and the second moment of the short-term transition probability density function. This method does not require a step size. However, the moment equations of nonlinear stochastic systems are usually infinite and non-closed. Based on the Gaussian truncation method, the short-term displacement and velocity transition probability density functions of the linear and NES oscillators and their joint transition probability density functions can be presented as follows:

$$p(x_{mn}^i, t_i | x_{kl}^{i-1}, t_{i-1}) = \frac{1}{\sqrt{2\pi}\sigma(t_i)} \exp\left(-\frac{(x_{mn}^i - m_1(t_i))^2}{2\sigma^2(t_i)}\right), \quad (23)$$

$$\begin{aligned} & p(x_{mn}^i, y_{mn}^i, t_i | x_{kl}^{i-1}, y_{kl}^{i-1}, t_{i-1}) \\ &= \frac{1}{\sqrt{2\pi}\sigma_1(t_i)\sigma_2(t_i)\sqrt{1-\rho_{12}(t_i)}} \exp(-((\sigma_2^2(t_i)(x_{mn}^i - m_{10}(t_i))^2 \\ &\quad - 2\sigma_1(t_i)\sigma_2(t_i)(x_{mn}^i - m_{10}(t_i))(y_{mn}^i - m_{01}(t_i)) \\ &\quad + \sigma_1^2(t_i)(y_{mn}^i - m_{01}(t_i))^2)(2\sigma_1^2(t_i)\sigma_2^2(t_i)(1-\rho_{12}(t_i))^2)), \end{aligned} \quad (24)$$

where

$$\begin{cases} m_{ij} = E[X^i \dot{X}^j], & \sigma_1^2(t_i) = m_{20}(t_i) - (m_{10}(t_i))^2, \\ \sigma_2^2(t_i) = m_{02}(t_i) - (m_{01}(t_i))^2, & \sigma_1(t_i)\sigma_2(t_i)\rho_{12}(t_i) = m_{11}(t_i) - m_{10}(t_i)m_{01}(t_i). \end{cases} \quad (25)$$

Substitute Eq. (25) into Eqs. (23) and (24). Then, the global transition probability density function can be obtained

4 Simulation

The reduced interval for the path integral method is selected from -0.5 to 0.5 , and is divided into 500 consistent sub-intervals with 10 Gauss-Legendre orthogonal points in each sub-interval, i.e., $K = 500$, $z_k = 1/500$, and $L = 10$. The time step is 0.01.

In order to better evaluate the performance of the vibration reduction of the system under the harmonic and random excitation, the transmissibility transition probability density defined by the standard deviation ratio of the passed force and the excitation and the percentage of the energy absorption transition probability density of the linear oscillator are used to evaluate the performance of the vibration reduction of the system.

$$T = \frac{S_d(y_1 + \zeta_1 \dot{y}_1)}{D + R_{MS}(f_t \sin(\gamma\tau))}, \quad (26)$$

$$\eta_p = A/B, \quad (27)$$

where

$$\begin{cases} A = S_d\left(\frac{1}{2}m_1\dot{y}_1^2 + \frac{1}{2}\omega_0^2 y_1^2 + \int_0^t c_1 \dot{y}_1^2(t) dt\right), \\ B = S_d\left(\frac{1}{2}m_1\dot{y}_1^2 + \frac{1}{2}\omega_0^2 y_1^2 + \int_0^t c_1 \dot{y}_1^2(t) dt + \frac{1}{2}m_2\dot{y}_2^2 \right. \\ \quad \left. + \frac{1}{2}k_n(y_1 - y_2)^4 + \int_0^t c_1(\dot{y}_1^2(t) - \dot{y}_2^2(t)) dt\right). \end{cases}$$

Table 1 Parameters and their values

Parameter	Symbol	Value
Nonlinear stiffness coefficient	k_2	300
Damper ratio of the linear oscillator	ζ_1	0.1
Damper ratio of the NES oscillator	ζ_2	0.15
Mass ratio	ε	0.1
Harmonic excitation amplitude	f_t	10
Frequency ratio	γ	0.5
Strength of Gaussian white noise	D	0.01

Figures 2 and 3 depict the transition probability densities of the displacement and velocity of the linear and the NES oscillators. The numerical result shows that the peak of the linear oscillator coupled with the NES in the equilibrium is higher than the peak of a two-degree-of-freedom linear system.

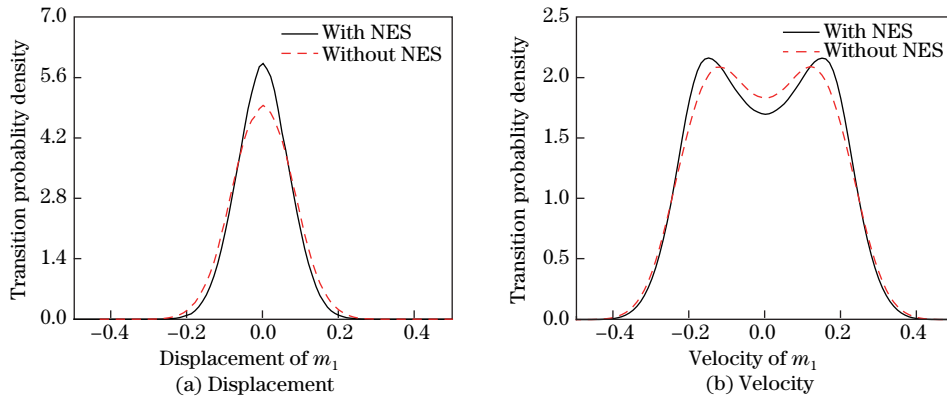
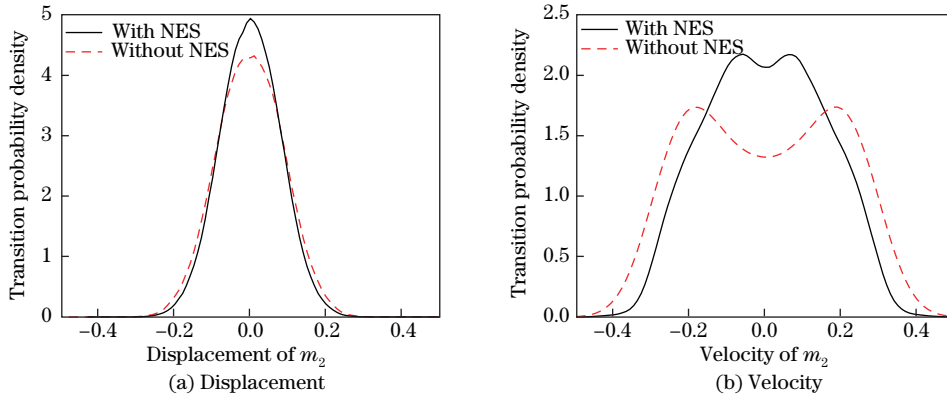
**Fig. 2** Transition probability densities of the displacement and velocity of the linear oscillator**Fig. 3** Transition probability densities of the displacement and velocity of the NES oscillator

Figure 4 shows the joint transition probability densities of the displacement and velocity of the linear and NES oscillators. From Fig. 4, we can see a typical response exhibiting a random jump phenomenon of the linear system coupled with the NES under the joint harmonic and Gaussian white noise. The essence of random jump is a response transition, and is near the equilibrium state.

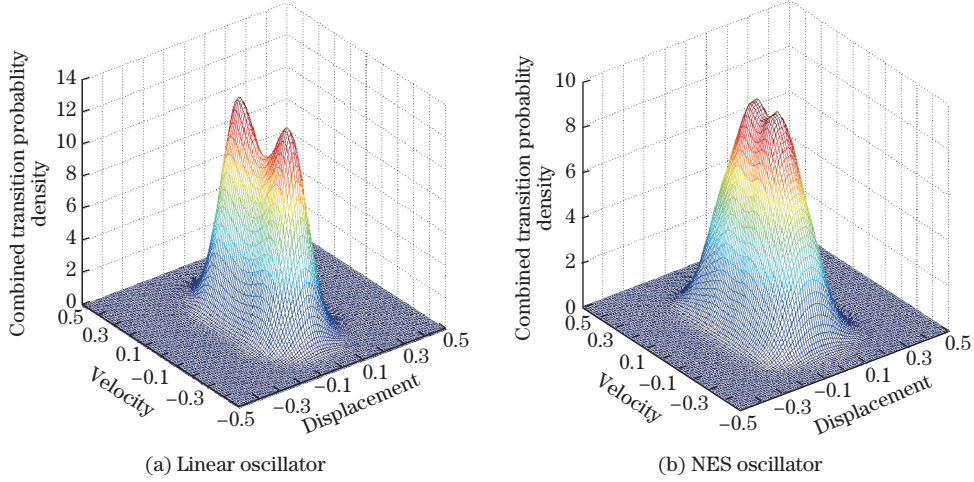


Fig. 4 Combined transition probability densities of the displacement and velocity of the linear and NES oscillators (color online)

Figures 5 and 6 reveal the transmissibility transition probability density and the energy absorption percentage of the linear oscillator transition probability density. From the figures, it can be seen that the transition probability density of the linear oscillator coupled with the NES is larger than that of the two-degree-of-freedom linear system when the transmissibility and the energy absorption percentage of the linear oscillator are close to zero.

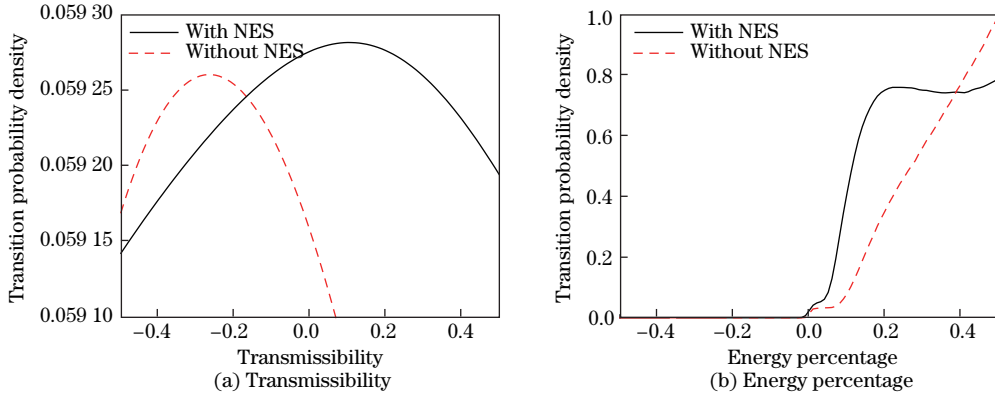


Fig. 5 Transition probability densities of transmissibility and energy percentage

5 Parameters analysis

In order to discuss and analyze the effects of the vibration reduction performance of the linear system coupled with the NES under harmonic and Gaussian white noise excitations with different parameters, the effects of the nonlinear stiffness k_2 and the damper ratio ζ_2 are investigated.

Figures 6 and 7 show the probability densities of the displacement and velocity of the linear and NES oscillators except the difference in the nonlinear stiffness. It can be seen that in all four cases shown in Figs. 6 and 7, the velocity transition probability density is bimodal, which indicates that random jump may occur. However, there is still a slight difference between them. The two peaks are slightly separated, and jump occurs more likely with the increase in the nonlinear stiffness.

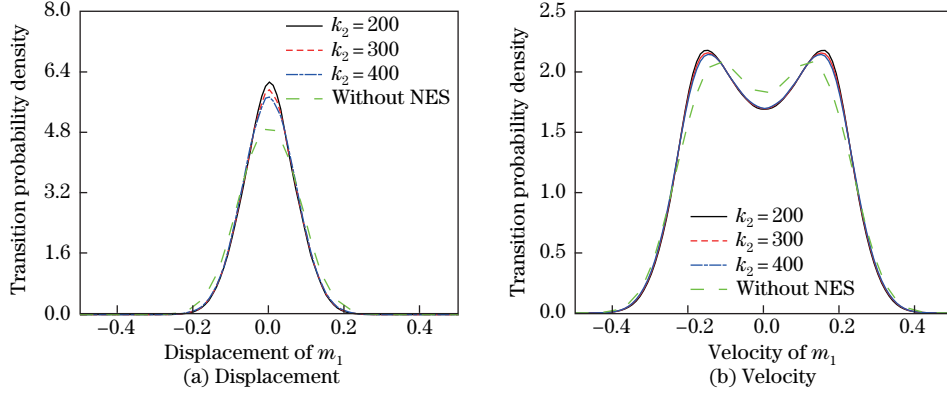


Fig. 6 Transition probability densities of the displacement and velocity of the linear oscillator except the difference in the nonlinear stiffness

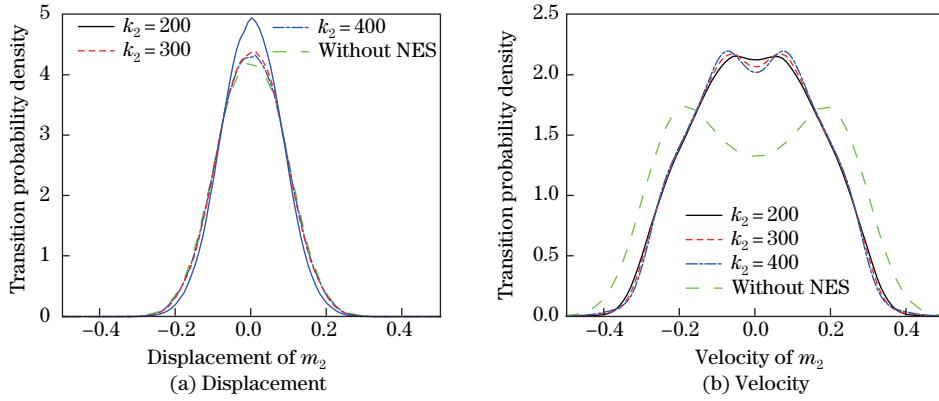


Fig. 7 Transition probability densities of the displacement and velocity of the NES oscillator except the difference in the nonlinear stiffness

Figure 8 presents the transition probability densities of the transmissibility and energy absorption percentage of the linear oscillator except the difference of the nonlinear stiffness. It can be seen that the transition probability density increases with the increase in the nonlinear stiffness when the transmissibility and the energy absorption percentage of the linear oscillator approaches to zero.

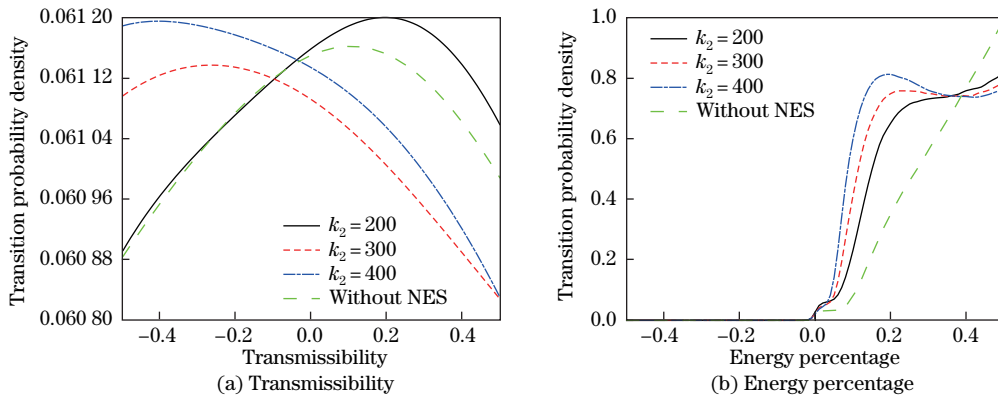


Fig. 8 Transition probability densities of the transmissibility and energy percentage except the difference in the nonlinear stiffness

Figures 9 and 10 present the probability densities of the displacement and velocity of the linear and NES oscillators except the difference of the damper ratio. In all four cases shown in Figs.9 and 10, the velocity transition probability density is bimodal. Not only it is possible for the occurrence of random jump, but also there is a large difference between them. The two peaks are obviously separated, and jump occurs more likely with the increase in the damper ratio. It can be expected that when the damper ratio approaches to zero, the random jump may disappear. The appearance and disappearance of random jump with the changes of the system parameters are called random jump bifurcations.

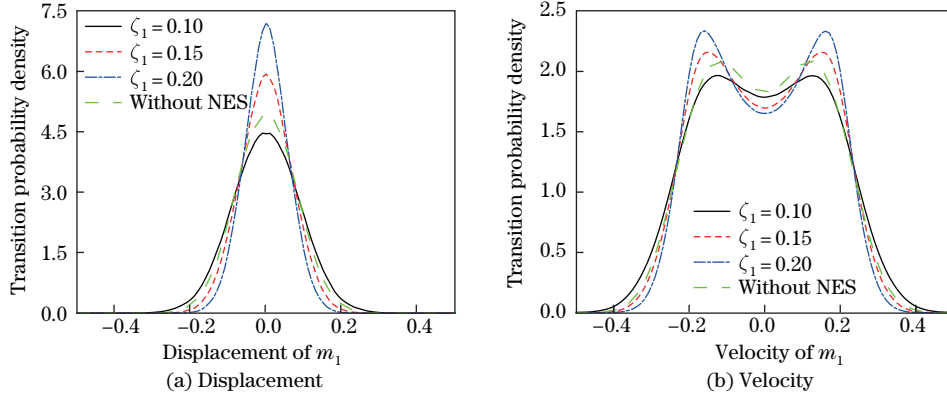


Fig. 9 Transition probability densities of the displacement and velocity of the linear oscillator except the difference in the damper ratio

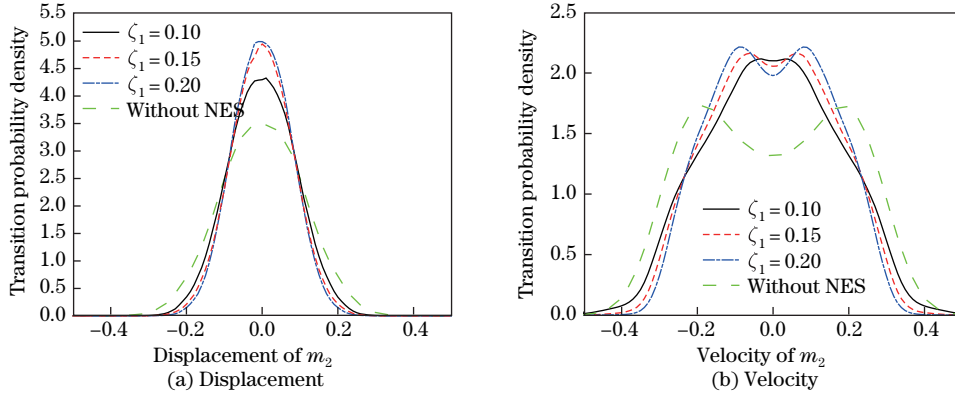


Fig. 10 Transition probability densities of the displacement and velocity of the NES oscillator except the difference in the damper ratio

Figure 11 presents the transition probability densities of the transmissibility and energy absorption percentage of the linear oscillator except the difference in the damper ratio. The transition probability density decreases with the increase in the damper ratio when the transmissibility or the energy absorption percentage of the linear oscillator is close to zero.

6 Conclusions

This paper mainly focuses on the numerical calculation of the response of the linear system with an NES under a harmonic and Gaussian white noise excitation. The path integral based

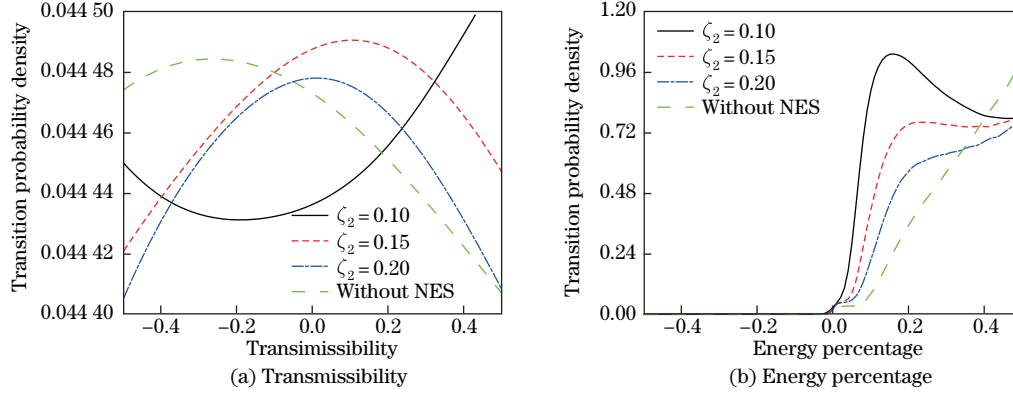


Fig. 11 Transition probability densities of the transmissibility and energy percentage of the linear oscillator except the difference in the damper ratio

on the Gauss-Legendre polynomial is used to solve the steady-state Fokker-Planck-Kolmogorov equation. The transition probability densities of displacement and velocity are used to determine the dynamical behaviors. The transition probability densities of the transmissibility and input energy absorption percentage of the linear oscillator are used as the indicators to evaluate the performance of vibration reduction. The investigation shows that an NES with reasonable parameters also has an efficient effect of vibration reduction. The conclusions lie in the following aspects.

(i) The linear system with an NES still has a good performance of vibration reduction under a harmonic and random excitation.

(ii) Random jump may appear in the linear system with an NES. The jump may occur from one peak to another peak randomly at any frequency of the harmonic excitation. It is harmful to the stability of the linear system with an NES system.

(iii) Increasing the nonlinear stiffness can effectively improve the vibration reduction performance of the linear system with an NES while ignoring its influence on the system stability, because the change is slight in the same system except the difference in the nonlinear stiffness. Decreasing the damping ratio can effectively improve the performance of vibration reduction of the linear system with an NES and enhance the stability at the same time.

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