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1 ORIGINAL ARTICLE

2 **Transient in a two-DOF nonlinear system**

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Abstract Dynamics of a system containing a linear oscillator, linearly coupled to an essentially nonlinear attachment, is considered. A damping is taken into account. It is assumed that some initial excitation implies vibrations of the linear oscillator. Envelops of the subsystem's kinetic energies are selected to use the numerical investigation of transient in the system. The parametrical optimization approach is used to obtain regions of effective energy transfer in the system parameter space. It is demonstrated that this efficient energy transfer may be obtained for a rather small value of the attachment mass.

Keywords Transient. Nonlinear energy transfer • Parametric optimization

20 **1 Introduction**

The absorption problem is very important in engineering. The classical passive absorber in the form of the single-DOF linear oscillator was used first by Frahm [1] to reduce forced oscillations. He discovered that an extinguishing of the resonance vibration is possible

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if the fundamental frequency of the absorber is equal to the frequency of the external periodical force. To avoid the resonance vibrations near the fundamental frequencies of the 2-DOF system containing coupled oscillators, Den Hartog [2] used the linear damped absorber. In many cases the absorption can be made effective by using linear absorbers with big masses, but this is impossible to realize in most concrete systems, particularly in civil engineering.

Roberson and Arnold [3, 4] considered the non-damped single-DOF absorber connected with a main system by using the spring with a cubic characteristic. They showed that the soft connecting spring allows to extend the external frequencies interval where the forced vibration of the main system can be reduced.

Over the past years numerous new devices have been used for the vibration absorption and for the reduction of the transient responses of structures. It seems interesting to study nonlinear passive absorbers for this reduction in nonlinear structures. In literature, nonlinear absorbers have been already used like in [3, 4], or [5] where a pendulum-type centrifugal vibration absorber is used to reduce torsion oscillations, or in [6] where the pendulum is also used as a vibration absorber to reduce response of a flexible cantilever beam, etc. Impact absorber was also considered [7, 8].

A description of the energy transfer was first presented in work by Witt and Gorelik [9]. Authors of this work tried to explain in the point of view of the classical mechanics the effect of the splitting in the spectrum of combination scattering. They considered the spring



57 pendulum as an example of two nonlinear coupling lin-
 58 ear partial subsystems. It was shown that for any small
 59 connection the absolute transient of energy of angular
 60 oscillations to energy of vertical oscillations and back
 61 takes place for the frequency ratio which is equal to 2:1.
 62 Quickness and intensity of this transient is dependent
 63 on initial conditions. For some initial conditions, this
 64 transient is absent (this is a case of periodical solutions),
 65 and the partial subsystems are synchronized.

66 Transfer of energy in a case of bending and tor-
 67 sional vibrations of elastic systems is described by
 68 Kononenko [10]. Transfer of energy was investigated
 69 too by Struble and Heinbockel [11, 12], Cheshankov
 70 [13], Celman [14], and Mercer et al. [15]. In a book
 71 by Starzhinsky [16], the energy transient in Lyapunov
 72 systems, and systems close to ones, was analyzed. Dy-
 73 namics of the spring pendulum, pendulum on the elastic
 74 suspension, and elementary particles in cyclic acceler-
 75 ators were considered. It was shown that the problem
 76 of the energy transfer groups together with a problem
 77 of dynamics of oscillation chain. First stage of analysis
 78 is to determine a periodic solution and regions of its
 79 instability in the system parameter space by using the
 80 general mathematical theory of parametric resonance.
 81 Second stage is a determination of bifurcation periodic
 82 solutions which appear for some critical parameter val-
 83 ues. Asymptotic methods permit to analyze the energy
 84 transfer.

85 The energy transient from one nonlinear vibration
 86 mode to another, is considered in a book by Nayfeh and
 87 Mook [17]. In [18], it was shown that for some condi-
 88 tions this transfer from high-frequency modes to low-
 89 frequency mode is possible in quasilinear systems. But,
 90 this transfer is caused by the modes interaction, and the
 91 one-way channeling of energy is not obligatory. Trans-
 92 fer of energy from high-frequency to low-frequency
 93 modes of the elastic system nonlinear vibrations was
 94 described by Hedrih [19, 20].

95 In papers by Vakakis, Manevitch, Gendelman,
 96 Bergman, and others [21-25], theoretical investigation
 97 and some experimental verification on the use of non-
 98 linear localization for reducing the transmitted vibra-
 99 tions in structures subjected to transient base motions
 100 has been presented. Conditions for the effective energy
 101 transfer and localization are discussed. The experimen-
 102 tal assembly, containing the main linear subsystem and
 103 the nonlinear absorber is described in [24]. Authors of
 104 the works [21-25] define the controlling process of the
 105 space energy transfer from the place of their initial ap-

pearance in the system under consideration to the other
 specific place where the energy, as the final result, is lo-
 calized, as the “energy pumping.” It means a controlled,
 unidirectional process of the energy transfer to passive
 nonlinear receiver where the energy is localized and di-
 minishes in time due to damping dissipation. Energetic
 criterion of this pumping is not proposed. Regions of
 this energy pumping in parameter space of mechanical
 systems were also not obtained.

In the presented paper, the authors use effective
 methods of global optimization to obtain such regions.
 This paper is organized as follows: in Section 2, the re-
 cent methods of global optimization are presented, and
 the $\hat{\Lambda}$ -transformation method, used here, is described
 in detail. Then, in Section 3, a character of transient for
 two mechanical models, namely, for the spring pendu-
 lum and for a system of two linearly connected oscilla-
 tors, one of them linear, and other nonlinear, is consid-
 ered. In particular, envelopes of the subsystem’s kinetic
 energies are analyzed. Then, some criterion of the full
 energy transfer from the initially perturbed linear sub-
 system to the nonlinear attachment, is formulated. By
 using this criterion, the curves of the full energy transfer
 in place of the second model parameters, are obtained.
 This full energy transfer is illustrated by numerical sim-
 ulation. Additional criterion of optimization, implied
 with the energy, which returns to the linear subsystem,
 is discussed and used to obtain corresponding regions
 in the system parameter place. The character of the
 energy transfer in 2-DOF system, containing the es-
 sentially nonlinear oscillator with a small mass, which
 corresponds to concrete engineering practice, is studied
 in Section 4. Regions of the effective energy transfer
 in the parameter place are obtained. In this section, the
 influence of degree of nonlinearity and initial energy
 on the energy transfer is discussed too.

**2 Methods of global optimization: Global
 minimum search by using the method
 of $\hat{\Lambda}$ -transformation**

There are a lot of works on determination of optimal
 solutions of theoretical and applied problems, begin-
 ning from classic works by Bernoulli, Euler, Lagrange,
 Hamilton, and others. Considering concrete problems,
 it is difficult, or, may be, impossible to obtain properties
 of the input quantities and output parameters analytical
 dependence. So, there exists a problem of construction

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152 methods which permit to determine optimal solutions
 153 in a case when assumptions about a nature of function
 154 under consideration are absent.

155 2.1 Existent approaches to solve the global
 156 optimization problems

157 There are no universally effective algorithms to solve
 158 the global optimization problems.

159 All known global optimization methods (GOMs)
 160 can be divided into two categories: deterministic and
 161 stochastic [26, 27]. Deterministic methods permit to
 162 obtain the global solution by using the exhaustive find-
 163 ings over all the problem feasible sets. So, deterministic
 164 methods lose an effectiveness if the dimension of the
 165 problem increases. Moreover, such methods need in
 166 addition a limitation on the criterion function. Stochas-
 167 tic algorithms permit to get away from such problems.
 168 A majority of stochastic methods permit to estimate a
 169 value of the criterion function in randomized points of
 170 the feasible set with the next data processing.

171 The first method of global optimization was the
 172 *Monte-Carlo method*. On the basis of the Monte-Carlo
 All173 method, the *multistart method* [28] was created. In this
 174 method, the same subset consisting of the N points is
 175 chosen (stochastically or deterministically) from some
 176 set of the problem feasible points. The local descent
 177 algorithm is started from each point, and the best so-
 178 lution is chosen from all obtained local solutions. This
 179 method is not very effective, because one and the same
 180 local solution can be found sometimes. But some me-
 181 fective GOMs were based of the principal idea of the
 182 multistart method.

183 It is impossible to present here in detail all exist-
 184 ing GOMs whose description can be found in different
 185 publications. One mentions here some GOMs [26-33],
 186 namely, the *clustering method* which is a modification
 187 of the multistart method (here a careful selection of
 188 points of the local finding is made); the *topographical*
 189 *method*; the *bisection (covering) and interval (branch-*
 190 *and-bound) methods*; the *branch-and-bound strategy*
 191 which divides the region of finding to a set of many-
 192 dimensional cubes; the *tunneling method* (a principal
 193 deficiency of the method is a necessity to solve diffi-
 194 cult nonlinear differential equations); the principal idea
 195 of the method of *simulated annealing* comes from the
 196 physics of the liquid freezing; the main idea of the *evo-*
 197 *lutionary algorithms* is taken from the biological evolu-
 198 tion processes; the *controlled random search algorithm*

199 which is a method of the direct search; *trajectory meth-*
 200 *ods*, in particular, the *continuation method*.

201 2.2 The \wedge -transformation method

202 At present, it is impossible to predict which GOM
 203 is the most effective. In this paper, one of the most
 204 universal and effective GOMs is used, namely, the
 205 \wedge -transformation method [34].

206 The main peculiarity of this method is that an
 207 object of the analysis is not the criterion function
 $F(x_1, x_2, \dots, x_n)$, but some function $\wedge(Z)$, which is
 a result of the function $F(x_1, x_2, \dots, x_n)$ transforma-
 tion. This transformation is based on the definition of
 partition utilized for the Lebesgue integral construc-
 tion. This conception can be successfully used to solve
 problems the analysis of which was difficult up to
 now, in particular, problems with the criterion func-
 tion $F(x_1, x_2, \dots, x_n)$ which is not differentiable in a
 point of extremum.

217 Let E be some measurable set from the space R^N ,
 218 and $F(x)$ some function from the class $L_p(E)$. This set
 219 can be divided to a sequence of the linked subsets e_{oi}
 ($i = 1, 2, \dots, \nu$), and on each subset e_{oi} the function
 220 $F(x)$ is convex. For the three-dimensional space this
 221 situation is illustrated by Fig. 1.

222
 223 If $E^* = \{x : x \in E, F(x) > Z\}$. One defines then as
 224 $m(E^*)$ the measure of the set E^* . One introduces the
 225 following function:

$$m(E^*) = m(E^*). \tag{2.1}$$

226 This function can be determined by the next
 Lebesgue integral:

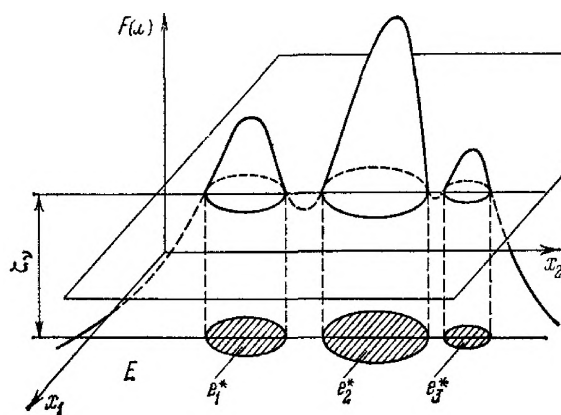


Fig. 1 Partition of the set E

$$m(E) = \int_{I_1} \int_{I_2} \dots \int_{I_s} \theta_V(x_1, \dots, x_n) dx_1 \dots dx_n \quad (2.2)$$

where $\theta_Z(x_1, \dots, x_n)$ is the characteristic function of the form

$$\theta_V(x_i) = \begin{cases} 1, & x_i \in E \setminus Z \\ 0, & x_i \in Z \end{cases} \quad (2.3)$$

One performs now the Lebesgue's partition of the function $F(x)$. An interval of the $F(x)$ range is divided for a set of subintervals,

$$0 = Z_0 < Z_1 < \dots < Z_v < \dots < Z^* = \max F(x). \quad (2.4)$$

Note that a finding of the minimum is principally the same.

It can be shown [34] that the transformed function \hat{Z} has the following principal properties: the function, independently of the prototype $F(x)$ dimension, is a function of the single variable; the function is monotone and nonincreasing one for all $F(x) \in L_p(E)$; a value Z^* , when $\hat{Z} = 0$, is equal to a value of the function $F(x) \in L_p(E)$ global extremum, $m\{x : F(x) = F_{\max}\} = 0$.

Analytical dependence between the functions $F(x)$ and \hat{Z} can be obtained in explicit form only in simplest cases. Other possibility of the function \hat{Z} construction consists in its determination in different points, next approximation and extrapolation of the function, and in a determination of the function \hat{Z} zero, which corresponds to a value of the function $F(x)$ global extremum.

To determine the function \hat{Z} in some point Z_v , statistical tests can be used. It is conditionally assumed that a measure of the set E , on which the function $F(x)$ is defined, is equal to the unit, that is, $m(E) = 1$. Let us define later a value of the measure $m\{x : F(x) > Z_v\}$ with respect to $m(E)$. So, it depends on the determination of some probability, which can be presented in the form of the integral (2.2). The probability P_v can be calculated by the next simple way. In fact, after some number s of statistical tests, it is possible to calculate the number of events $\%_v$, for which $F(x) > Z_v$, and to calculate a frequency or, a statistical probability

$$P_v = H_v/S. \quad (2.5)$$

According to the law of big numbers, if the number s increases, the frequency tends (on probability) to the desired value $P_v = m(E^*)$.

So, the determination of the desired measure of the set E^* consists in conducting of statistical tests for which the variables $x_1, x_2, \dots, x_i, \dots, x_n$ take equiprobable random values. Each test consists in determination of the function $F(x_1, x_2, \dots, x_n)$ value for given values of independent variables, and in comparison of the value with the given level Z_v . Depending on results of this comparison one has these points x , for which in the v -th test the inequality $F(x) > Z_v$ is satisfied. Then, the approximate value of the measure of the set E^* is determined.

As a result, by using the statistical tests method, one obtains the random function \hat{Z} instead of the function \hat{Z} . Construction of this function permits determination of a scalar value of the global extremum, but coordinates of it remain unknown. But, is it possible to determine all coordinates of the function $F(x_1, x_2, \dots, x_n)$ global extremum from the data obtained in statistical tests, which are used to construct the function \hat{Z} . For this, the function $x(Z)$ must be determined, where x_i is a mean value of such x_{ij} ($i = 1, \dots, n; j = 1, \dots, s$), for which $F(x_1, x_2, \dots, x_n) > Z$. Calculation of the mean values x_i for each level of the i th coordinate can be made by the formula A3 (see Appendix). As stated earlier, by substitution a value ζ^* , which is the function $F(x)$ global extremum, to $\bar{x}_i(\zeta)$, one obtains the i th coordinate of the global extremum.

So, to solve the complex many-dimensional optimization problems by using the \hat{Z} -transformation method, it is necessary to determine values of the functions \hat{Z} and $x \setminus (Z)$ in different points Z_v , then, approximating the obtained data, to solve the formulated problem by method of extrapolation.

Algorithm of the \hat{Z} -transformation method is described in the Appendix.

3 Energy transfer in some mechanical models 302

3.1 Energy transfer in a spring pendulum 303

One considers the transfer of energy in a spring pendulum (Fig. 2). This problem was investigated in [9, 16]. 304 305



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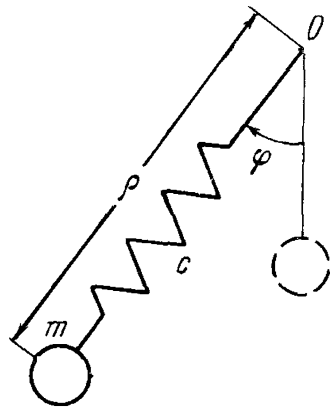


Fig. 2 The spring pendulum

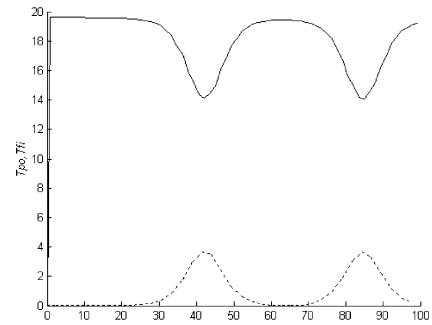
306 Differential equations of the system motion are the
 307 following:

$$\begin{aligned}
 p \ddot{\theta} + 2p \dot{\theta} \dot{\theta} &= -g p \sin \theta \\
 p \ddot{p} - p \dot{\theta}^2 &= -\frac{c}{m}(p - l) + g \cos \theta.
 \end{aligned}
 \tag{3.1}$$

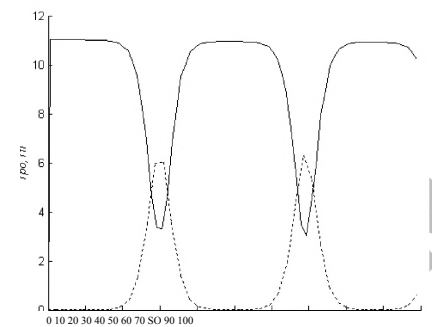
308 It was shown in [16] that for some values of the sys-
 309 tem parameters a disruption of the vertical vibrations
 310 takes place as a result of arbitrary small transversal
 311 perturbations.

312 One develops the numerical investigation of the
 313 Equations (3.1) in the point of the energy transfer and
 314 in some vicinity of the point. In this system the point
 315 is determined by the parameter $\gamma = mg$, and the en-
 316 ergy transfer takes place if $\gamma = 0.3333$. Results of thi
 317 numerical calculations are presented in Fig. 2. It is pre-
 318 sented as a dependence of the kinetic energies T_p and
 319 T_θ of time. Here $T_p = \frac{1}{2} m \dot{p}^2$ is a kinetic energy of the
 320 vertical vibration mode, and $T_\theta = \frac{1}{2} m l^2 \dot{\theta}^2$ is a kinetic
 321 energy of the angular vibration mode. The kinetic en-
 322 ergy maximum points are joined. We call correspond-
 323 ing lines as the kinetic energy envelopes. In Fig. 3, an
 324 envelop of the vertical vibration mode kinetic energy
 325 is depicted by solid line, and an envelop of the angular
 326 vibration mode kinetic energy is depicted by dot-and-
 327 dash line.

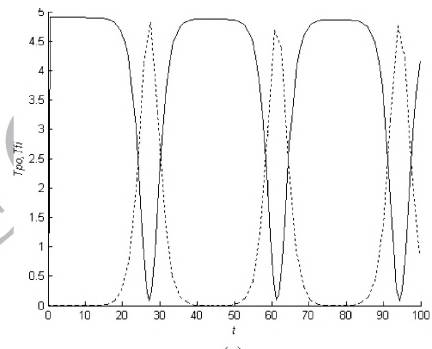
328 We can see from Fig. 3 that in some vicinity of
 329 the critical value of the parameter $\gamma = 0.333$, there
 330 is a decrease in the kinetic energy T_p amplitude val-
 331 ues, and the appearance of a kinetic energy T_θ peak.
 332 It is clear that all four diagrams illustrate a process
 333 of the energy transfer from one vibration mode to the



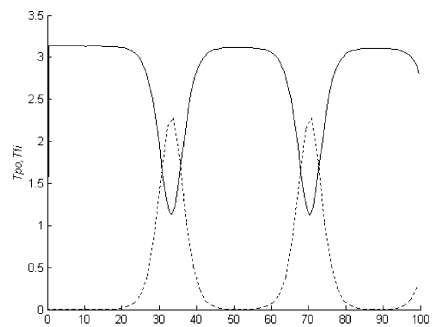
(a)



(b)



(c)



(d)

Fig. 3 Diagrams of envelopes of kinetic energies of the vertical and angular vibration modes, T_p and T_θ , for different values of the parameter γ : (a) $\gamma = 0.667$, (b) $\gamma = 0.5$, (c) $\gamma = 0.333$, (d) $\gamma = 0.266$. Here - T_p , - - - - T_θ

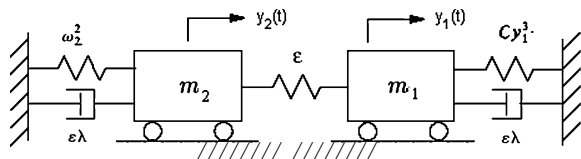


Fig. 4 System of weakly coupled linear and nonlinear oscillators

334 other. In addition, the full transfer of energy happens
 335 for $Y = 0.3333$.

336 3.2 Energy transfer in 2-DOF system containing an
 337 essentially nonlinear oscillator

338 One considers the energy transfer in a 2-DOF system
 339 (Fig. 4), containing the weakly coupled linear and non-
 340 linear oscillators.

341 One has the following differential equations of
 342 motion:

$$\begin{aligned} m_1 \ddot{y}_1 + ek y_1 + cy_1^3 + e (y_1 - y_2) &= 0 \\ m_2 \ddot{y}_2 + ek y_2 + C_2 y_2^2 + S (y_2 - y_1) &= 0. \end{aligned} \quad (3.2)$$

343 Here $\ell \ll 1$, other parameters have an order $O(1)$. Such
 344 models were considered in some works, for example,
 345 in works by Vakakis, Manevitch et al. Analysis of sys-
 346 tems with such essentially nonlinear absorbers permits
 347 to select principal advantages of the nonlinear absorp-
 348 tion, in particular, the effect of the energy localization
 349 In engineering practice, there is a possibility to obtain
 350 nonlinear absorbers with very small linear components
 351 in elastic characteristic.

352 A behavior of the system by using the envelopes
 353 kinetic energies (these envelopes were introduced in
 354 Section 3.1) of two subsystems is considered. Later,
 355 the adduced values of the kinetic energies of two par-
 356 tial oscillators are presented:

$$T_1 = m_1 \dot{y}_1^2, \quad T_2 = m_2 \dot{y}_2^2 \quad (3.3)$$

357 The kinetic energy envelopes obtained for the sys-
 358 tem (3.2) are shown in Fig. 5 for $m_1 = m_2 = 1$,
 359 $X = 0.5$, $c_2 = 0.9$, $c = 5.0$, $\ell = 0.1$ and for the next
 360 initial values: $y_1(0) = y_2(0) = 0$, $\dot{y}_1(0) = 0$; $\dot{y}_2(0) =$
 361 $\sqrt{2}h$, where h is the system energy at the moment;
 362 $t = 0$ (here $h = 0.8$). Here and later, a dotted line rep-
 363 resents an envelop of the adduced kinetic energy of ini-
 364 tially perturbed linear oscillator. Solid line represents

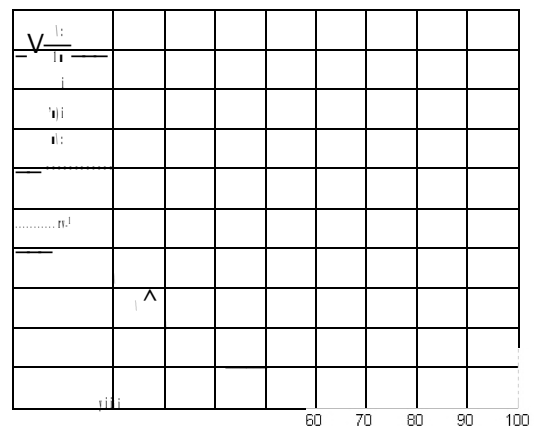


Fig. 5 Diagram of the kinetic energy envelopes for the system (3.2)

the same for the nonlinear absorber. Dot-a and-dash line
 365 represents the same for the linear oscil lator without the
 366 attached nonlinear subsystem. 367

By considering this diagram the next problem can
 368 be formulated: are there some parameters of the system
 when the kinetic energy envelop of the linear initially
 369 perturbed subsystem turns into zero, and the peak of
 370 such envelop of the perturbed nonlinear system simult-
 371 aneously appears. Such case will be assumed as a case
 372 of the full energy transfer. 373 374

One selects in Fig. 6 principal parameters which
 375 characterize the energy transfer. The point T_{2min} corre-
 376 sponds to the first minimum of the kinetic energy en-
 377 velop of the initially perturbed linear subsystem. The
 378

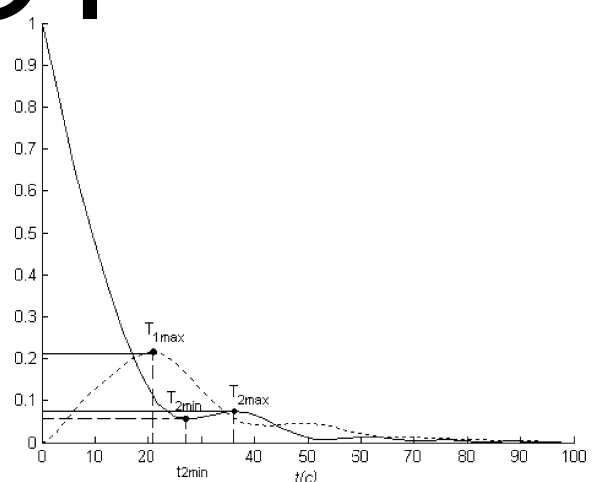


Fig. 6 Principal parameters of a diagram of the kinetic energy envelopes

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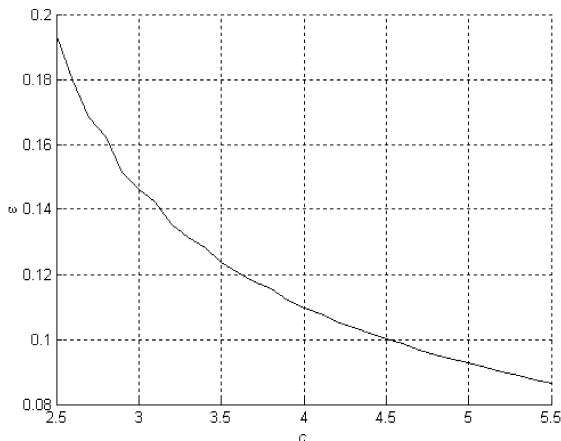


Fig. 7 Curve of the full energy transfer in plane (s, c)

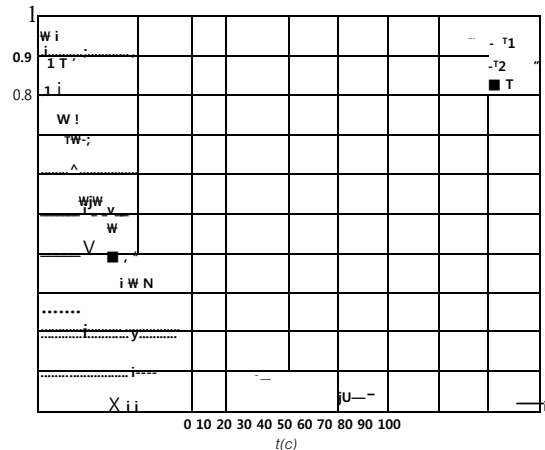


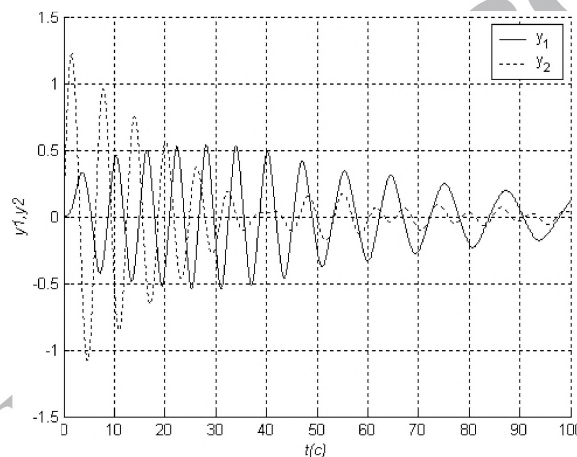
Fig. 8 Envelopes of the kinetic energies in point of the full energy transfer ($s = 0.09268, c = 5.0$)

379 point $T_{1 \max}$ represents a maximum of this envelop of
 380 the nonlinear subsystem. The point $T_{2 \max}$ represents
 381 the second maximum of this envelop for the linear sub-
 382 system, and the point $t_{2 \min}$ is a time of achievement
 383 of the first minimum of this envelop for the perturbed
 384 linear subsystem.

385 One studies in detail the energy transfer in the
 386 system (3.2). One takes the system parameters and
 387 initial values which were used earlier. By using the
 388 proposed criterion of the full energy transfer and the
 389 -transformation method, the curve of the full energy
 390 transfer in a plane of the system parameters (s, c) was
 391 obtained. This curve is shown in Fig. 7.

392 One selects some point on the curve, and observes a
 393 change of the system vibration energy. As previously,
 394 the dotted line represents on diagram the envelop of the
 395 added kinetic energy of the initially perturbed linear
 396 oscillator, the solid line represents the same for the non-
 397 linear absorber, and the dot-and-dash line represents
 398 the same for the linear oscillator without the attached
 399 nonlinear oscillator. Envelops are shown in Fig. 8 for
 400 the point of the full energy transfer ($T_{2 \min} < 0.0001$),
 401 namely, $s = 0.09268, c = 5.0$. We can see here that the
 402 full energy transfer from the initially perturbed linear
 403 oscillator to nonlinear one takes place on the time inter-
 404 val from zero to 40. At the moment $t = 40$ all energy
 405 is concentrated in the nonlinear absorber.

406 The transient in the system (3.2) in point of the full
 407 energy transfer is shown on the Fig. 9. We can see that
 408 the vibration amplitude of the initially perturbed sys-
 409 tem decreases up to zero on the time interval from zero
 410 to 40. At that time the vibration amplitude of the non-
 411 linear attachment increase appreciably. Beginning with



9 Transient in the system (3.2) in point of the full energy transfer

some moment close to $t = 30$ the vibration amplitudes
 of the linear subsystem are smaller than the vibration
 amplitudes of the attached nonlinear subsystem in three
 and more times. Then, the back energy transfer to the
 linear subsystem happens. Vibration amplitudes of the
 main linear oscillator increase. Then the energy anew
 leaves the linear subsystem. So, one observes the phe-
 nomenon similar to that which was observed in the sys-
 tem (3.1). The sequential transfer from one vibration
 mode to another one takes place.

Let us see a behavior of the system under consider-
 ation, if the parameter s , representing the stiffness of
 the connected spring, is relatively small. One can see
 from the Figs. 10 and 11 that in this case the energy
 transfer from the initially perturbed linear oscillator to

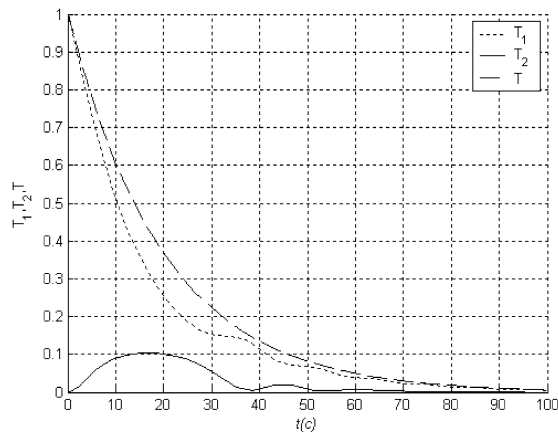


Fig. 10 Kinetic energy envelops for $\ell = 0.09$

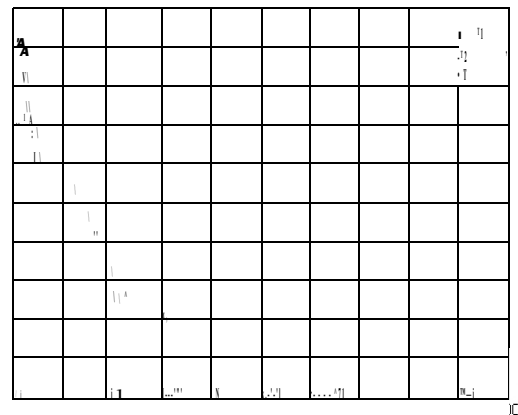


Fig. 12 Kinetic energies envelops for $\ell = 0.2$

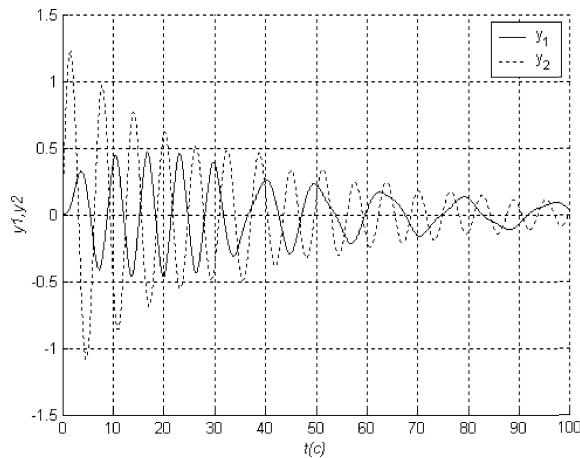


Fig. 11 Transient for $\epsilon = 0.09$

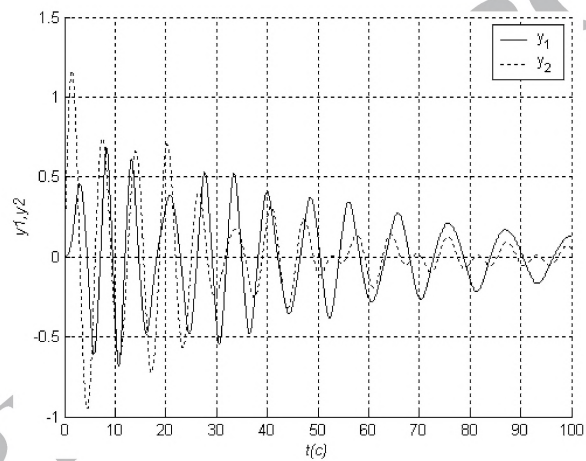


Fig. 13 Transient for $\epsilon = 0.2$

427 the nonlinear one is made/becomes weaker. Evidently,
 428 this is caused by the weak connection between oscilla-
 429 tors, which hinders the energy transfer to the nonlinear
 430 attachment.

431 Let us consider a behavior of the system under con-
 432 sideration, when the parameter ℓ , representing the stiff-
 433 ness of the connected spring, increases (Figs. 12-15).

434 One has from Figs. 12-15 that an increase of ℓ leads
 435 to more rapid energy transfer from the perturbed oscil-
 436 lator to unperturbed one. A peak of envelop of the un-
 437 perturbed mass kinetic energy increases together with
 438 increasing of ℓ , and it takes place for a smaller time
 439 interval. But, in this connection, the full energy trans-
 440 fer from the perturbed oscillator to unperturbed one is
 441 absent. There is a significant energy return back to the
 442 perturbed oscillator. Figure 14 shows that initially on
 443 the time interval from t_0 to t_{10} the rapid transfer of kinetic
 444 energy from the perturbed linear oscillator to nonlinear

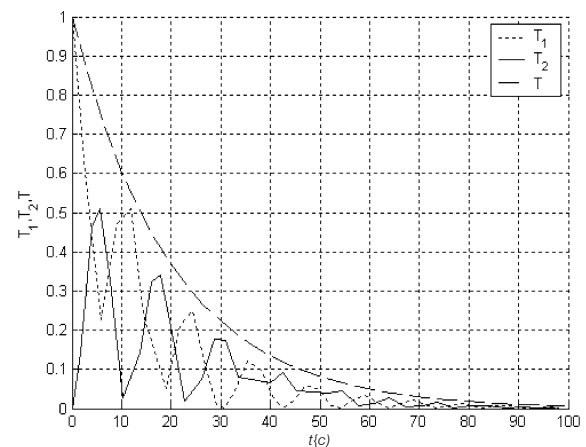


Fig. 14 Kinetic energy envelops for $\ell = 0.5$

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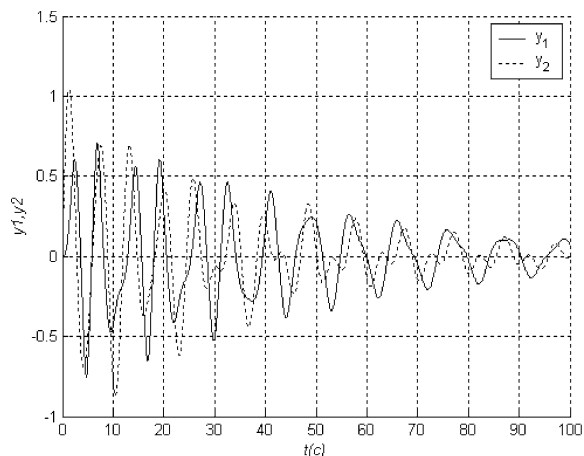


Fig. 15 Transient for $s = 0.5$

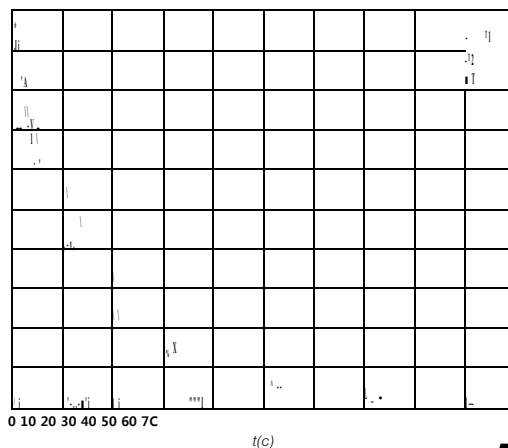


Fig. 16 Envelops of the kinetic energy in point of the full energy transfer, when $c = 2.5; s = 0.1933$

445 one takes place. But this is not the full energy transfer.
 446 Part of it remains in the perturbed oscillator. Then on
 447 the time interval from 10 to 19, this envelop of the ki-
 448 netic energy of the linear oscillator increases, while the
 449 kinetic energy of the nonlinear one sharply decreases.

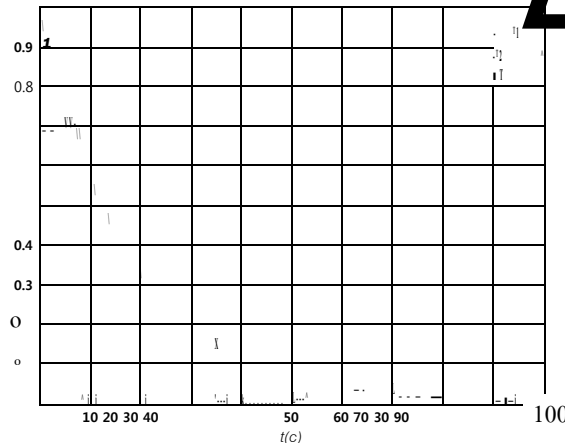
450 So, we conclude that an increase of the coefficient s
 451 leads to a larger return of the energy. The sequential
 452 transfer of the energy from one oscillator to another
 453 happens.

454 One considers anew a curve of the full energy trans-
 455 fer (Fig. 6). Let us analyze the energy envelops when
 456 we move to the left or to the right on the curve.

457 Some part of calculation of envelops of the kinetic
 458 energy are shown in Figs. 16 and 17 for different values .
 459 of the parameters c and s . It is possible to conclude t_{tran}^*
 460 that a decrease of c and an increase of s leads to a
 461 more rapid energy transfer from the initially perturbed
 462 linear oscillator to the nonlinear one. But, in this case,
 463 the larger return of the energy to the linear subsystem
 464 takes place. Increase of c gives us a reduction of the
 465 energy return to the linear oscillator.

466 The conclusion stated earlier is confirmed by
 467 Figs. 18 and 19, wherein points of the full energy trans-
 468 fer for the relations of the return energy value (after the full
 469 energy transfer) T_{2max} , and a time of the passage time of
 470 the full energy transfer t_{min} , are presented, depending
 471 on the change of the parameter c .

472 One considers now a problem of the energy trans-
 473 fer in the system of two connected oscillators from the
 474 point of view of the absorption problem. In this prob-
 475 lem the fast energy transfer from the main system to
 476 absorber is principal.



17 Envelops of the kinetic energy in point of the full energy transfer, when $c = 6; s = 0.0818$

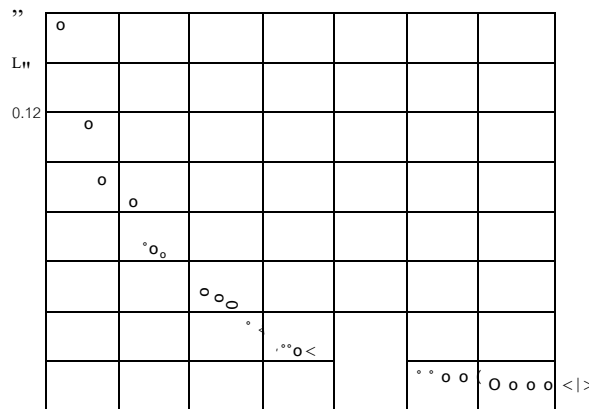
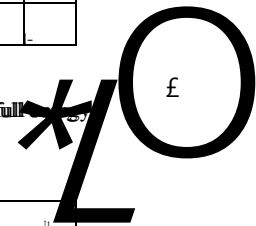


Fig. 18 Relation of T_{2max} depending on c



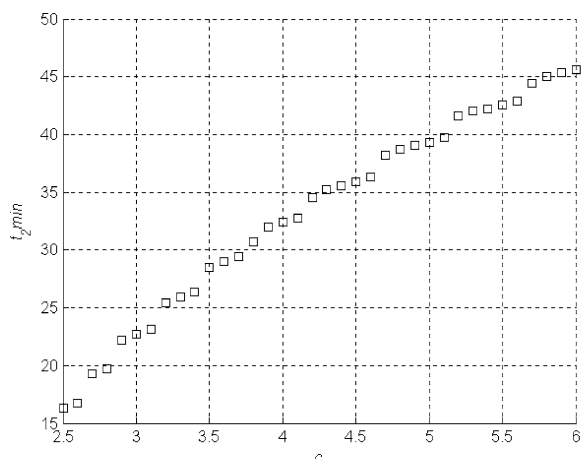


Fig. 19 Relation of t_{2min} depending on c

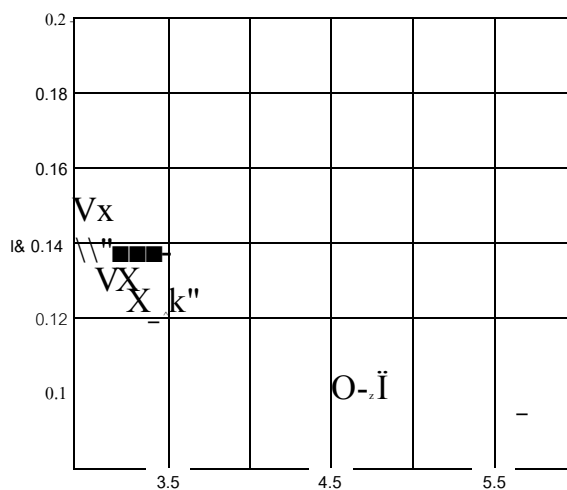
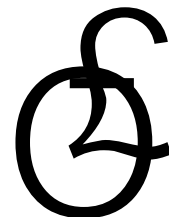


Fig. 21 Region of the parameters c and s , where the energy transfer takes place



0.15

0.08
2.5

5.5

Fig. 20 Classification of points of the full energy transfer

477 Additional limitation to the precedent criterion is in-
 478 troduced: is introduced it needs to minimize too T_{2max}
 479 (Fig. 6), that is the second maximum of the kinetic en-
 480 ergy envelop of the initially perturbed linear oscillator.
 481 It forms as a result of the energy return from the non-
 482 linear absorber to linear one. Introducing this criterion,
 483 we can to classified points of the full energy transfer
 484 (Fig. 7) on dependence of T_{2max} . Points on the Fig. 20
 485 represent values of parameters for which $T_{2max} < 0.5$,
 486 triangles represent a case when $0.5 < T_{2max} < 1$, and
 487 circles represent a case when $T_{2max} > 1$.
 488 Obtained results permit to formulate the next more
 489 general problem of the parametric investigation of the
 490 energy transfer: to find the values of the parameters
 491 c and s , when the envelop of the adduced kinetic en-
 492 ergy of the initially perturbed subsystem tends to zero.

In a neighborhood of this minimum the maximum of
 envelop of the nonlinear absorber exists.

Set of points, obtained by using this additional cri-
 terion, are shown in Fig. 21, where the solid line cor-
 responds to a maximal extinguishing of the initially
 perturbed oscillator energy ($r_{2min} < 0.0001$). Set of
 parameters, bounded by dotted lines, corresponds to
 the limitations $T_{2min} < 0.05$ and $T_{2max} < 0.05$. Re-
 gion of the parameter values where $T_{2min} < 0.1$ and
 $T_{2max} < 0.1$, are limited by dot-and-dash lines.

4 Character of the energy transfer in 2-DOF system containing the essentially nonlinear oscillator with a small mass

One considers now a case, when the attached nonlinear
 absorber has a mass which is essentially smaller than
 that of the linear subsystem. It corresponds to the real
 engineering practice when a use of the absorbers with
 big mass is impossible. The damping coefficient X is
 equal here to 0.1.

The energy envelops in a point of the full energy
 transfer for $m_1 = 0.1$ is shown in Fig. 22, and the cor-
 responding transient is shown in Fig. 23. Note that in
 Fig. 23, showing the transient, the solid line represents
 vibrations of the initially perturbed linear oscillator
 with attached absorber, and the dotted line represents
 such vibrations without this absorber. We can see here
 the essential extinguishing of the main linear subsystem

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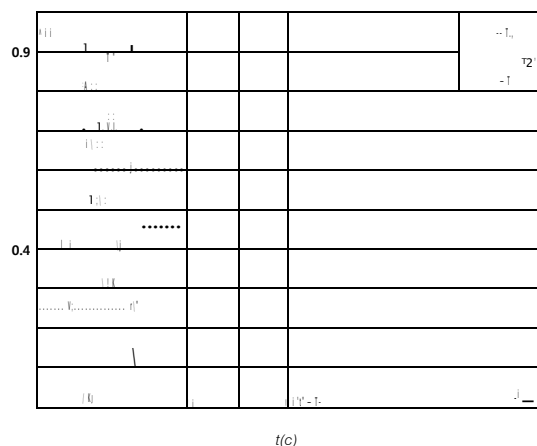


Fig. 22 Energy envelopes in a point of the full energy transfer for $m_1 = 0.1$



Fig. 24 Region of the effective energy transfer

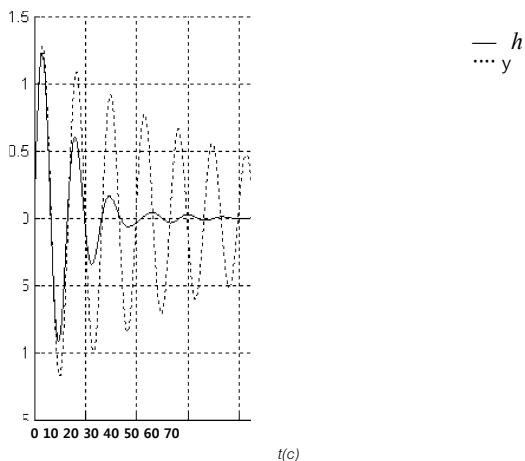


Fig. 23 Transient for linear subsystem with absorber (solid line) and without absorber (dotted line) when $m_1 = 0.1$

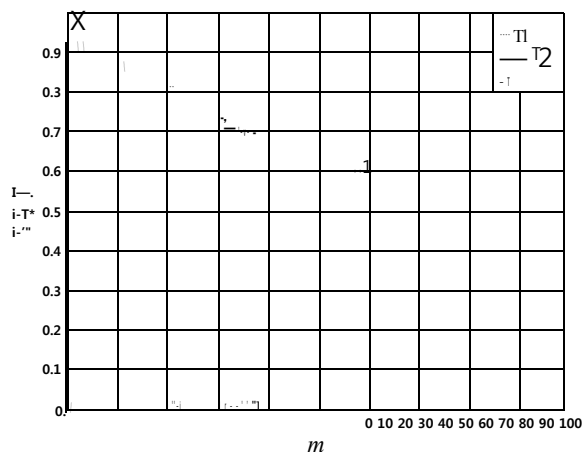
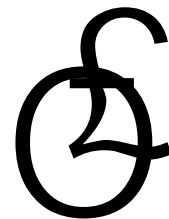


Fig. 25 Envelops of kinetic energies in a point of the full energy transfer ($c = 0.018, s = 0.04$)



520 vibrations if this nonlinear absorber exists. So, <
 521 ing the mass of absorber, the rapid one-way channeling
 522 of energy in the system under consideration is found.
 523 The next criterion of efficiency of the energy transfer
 A2524 (Fig. 24) is now introduced. Namely:
 525 The energy transfer must be effected for a fixed and
 526 not large time interval. (For example, here $t = 20$ is
 527 introduced). For a quantitative valuation of this trans-
 528 fer one estimates the following: if a loss of energy for
 529 the initially perturbed linear oscillator T_1 during this
 530 time interval is equal to 70% or more, the energy trans-
 531 fer is considered as effective (or optimal). Moreover,
 532 this condition has to be fulfilled in the point T_{2max} ,
 533 that is in a point of the second maximum of the kinetic
 534 energy T_1 envelop.

For the system (3.2), with the parameters $m_1 =$ 535
 0.1, $m_2 = 1$, $X = 0.1$, $c_2 = 0.9$, $s_1 = 0.1$ and initial 536
 conditions $y_1(0) = y_2(0) = 0$, $\dot{y}_1(0) = 0$; $\dot{y}_2(0) =$ 537
 $V_2 h$, where $h = 0.8$, the region of the effective energy 538
 transfer in a place of the parameters (c, s) was obtained. 539
 Boundary of this region is marked by boldface line in 540
 Fig. 26. Thin line represents a curve of the full energy 541
 transfer for this case. 542

One of the solutions from the region is presented in 543
 Figs. 25 and 26. 544

Investigation of influence of degree of nonlinearity 545
 and initial energy to the energy transfer was made too. 546
 Namely, three cases for the system (3.2) are considered: 547
 (1) the anchor spring is linear; (2) the anchor spring 548
 has a cubic nonlinearity; (3) the anchor spring has a 549
 fifth degree of nonlinearity. Corresponding points are 550

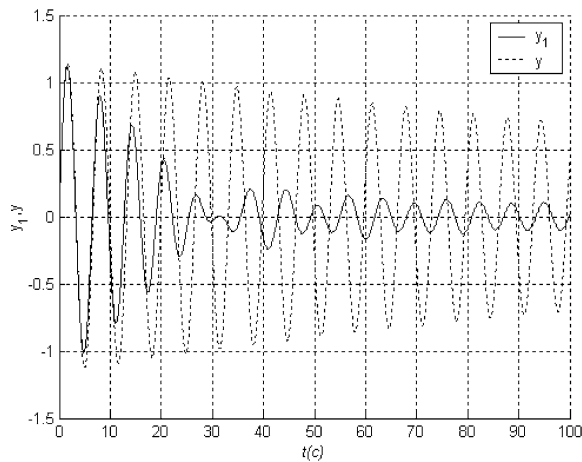


Fig. 26 Transient for $c = 0.018, s = 0.04$

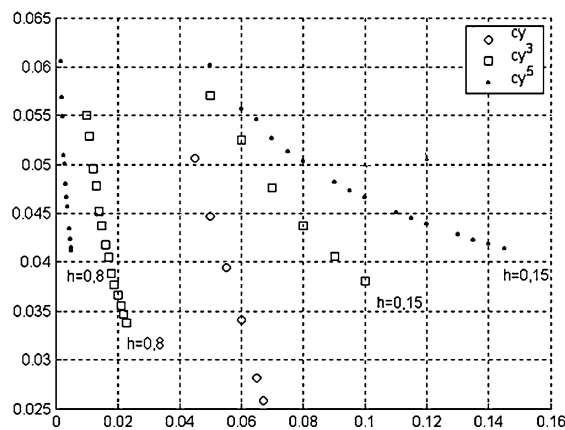


Fig. 27 Points of the full energy transfer in systems with different degree of nonlinearity and different initial energy

551 presented in Fig. 27 by using circles (linear spring),
 552 squares (cubic nonlinearity), and black points (fifth-
 553 degree nonlinearity) on a place of the parameters (c, s).
 554 It is very interesting that points of the full energy trans-
 555 fer for the nonlinear spring are different for different
 556 values of the initial energy.

557 5 Conclusions

558 In the present study, an analysis of the energy transfer in
 559 some 2-DOF nonlinear mechanical systems have been
 560 carried out. The effective method of global optimiza-
 561 tion, namely, the V -transformation method, is used
 562 here. Principal characteristics of the energy transfer,
 563 namely, envelopes of the subsystems kinetic energies,

are selected to use the numerical investigation of this
 process. Criterion of the full energy transfer is proposed
 and discussed. By using this criterion, the curves of the
 full energy transfer in a place of the system parameters,
 are obtained. This full energy transfer is illustrated by
 numerous numerical simulations. Additional criterion
 of optimization, implied with the energy, which returns
 to the linear subsystem, is discussed and used to obtain
 corresponding regions of the effective energy transfer
 in the system parameter place. Regions of the effective
 energy transfer in the parameter place are obtained. In-
 fluence of degree of nonlinearity and initial energy to
 the energy transfer is discussed too. It seems that the
 proposed approach can be used to investigate the trans-
 fer of energy in different nonlinear systems.

Appendix

The algorithm of the V -transformation method is pre-
 sented here.

Input data is the following: a, b are vectors of lower
 and high limitations of the parameter limitation; $f(\dots)$
 is a procedure of calculation of the minimized function,
 N is the number of the sketch points.

Output data is the following: x as a point of the
 minimum, $f(x)$.

1. The values $x_1, x_2, \dots, x_t, \dots, x_n$ are chosen by
 random low with uniform distribution.
2. A value of the function $F(x_1, x_2, \dots, x_n)$ is calcu-
 lated.
 The points 1 and 2 s times are repeated.
3. The $\sup F$ and $\inf F$ among s values of the function
 $F(x_1, x_2, \dots, x_n)$ are determined.
4. The interval $[(\sup F - \inf F)/2, \sup F]$ is divided
 into k equal components.
5. For all $k, Z_v (v = 1, 2, \dots, k)$ and $\hat{\cdot}$ are deter-
 mined according to the formulae:

$$Z_v = (\sup F - \inf F)/2 + (v - 1)AZ \quad (A.1)$$

$$V_v = \xi_v/s, AZ = Z_{v-1} - Z_{v-2} \quad (A.2)$$

where ξ_v is a number of events when, for the given
 ξ_v , one has $F(x) > Z_v$.

6. The obtained data are approximated so as to deter-
 mine the parabolic approximation.

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603 7. The formulated problem is solved by calculating
 604 the roots Z_1 and Z_2 of the parabola.
 605 8. The smaller values of the roots are found, cor-
 606 responding to a scalar value Z^* of the function
 607 $F(x_1, x_2, \dots, x_n)$ global extremum F_{\max} .
 608 9. For each Z_v ($v = 1, 2, \dots, k$) a mean value X_i ($i =$
 609 $1, 2, \dots, k$) is calculated by the formula:

$$X_i = \quad /H_v \quad (A.3)$$

610 10. The coefficients of parabolas which approximate
 611 the functions $x_i(Z)$ ($i = 1, \dots, k$) are deter-
 612 mined.
 613 11. The value Z^* , obtained in P. 8, is substituted to the
 614 expression, which was determined in P. 10, then
 615 one determines i th coordinate x_i^* of the global ex-
 616 tremum.
 617 12. The values x_i^* are substituted to the function
 618 $F(x_1, x_2, \dots, x_n)$ expression and one determines
 619 the required value F^* .
 620 13. The value F^* obtained in P. 12 is compared, with the
 621 scalar value of the global extremum Z^* , determined
 622 in P. 8.
 623 Equality of both values shows, with some error, that
 624 the problem is solved correctly.

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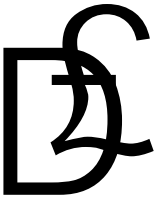
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Queries to Author

A1: Au: Please note references have been renumbered to provide for sequential arrangement. Please check.

A2: Au: Please check the citation of Fig. 24 for appropriateness



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