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Identification of a separation wave number between weak and strong turbulence spectra for a vibrating plate

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A weakly nonlinear spectrum and a strongly nonlinear spectrum coexist in a statistically steady state of elastic wave turbulence. The analytical representation of the nonlinear frequency is obtained by evaluating the extended self-nonlinear interactions. The *critical* wave numbers at which the nonlinear frequencies are comparable with the linear frequencies agree with the *separation* wave numbers between the weak and strong turbulence spectra. We also confirm the validity of our analytical representation of the separation wave numbers through comparison with the results of direct numerical simulations by changing the material parameters of a vibrating plate.

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I. INTRODUCTION

The coexistence of weakly nonlinear and strongly nonlinear spectra, which is often called critical balance, has been predicted in several *anisotropic* wave turbulence systems [1]. The characteristic time of the linear regime is the wave period: e.g., the Alfvén wave period in magnetohydrodynamics, axial component of the rotation frequency in rotating turbulence, and horizontal component of the buoyancy frequency in stratified turbulence. The nonlinear time scale is the horizontal eddy turnover time in these three types of turbulence. The critical balance is a plausible conjecture that illustrates the energy flux in an anisotropic wave turbulence, where the nonlinear time is comparable with the linear dispersion time across a wide range of wave numbers. Thus, it is of great importance to understand the physics in the wave numbers at which the linear and nonlinear time scales are comparable. Even if a system is *isotropic*, the linear time scale may be comparable with the nonlinear time scale. In this paper, the clarification of the coexistence in an isotropic wave turbulence system is addressed from the viewpoint of the time scales.

Various energy spectra have been observed for the wave turbulence in a thin elastic plate. The possibility of the Kolmogorov-Zakharov spectrum, which is predicted by the weak turbulence theory [2], provides the motivation for the studies in this wave turbulence system. In a numerical study [3], the weak turbulence spectrum $\mathcal{E}(k) \propto k[\log(k_*/k)]^{1/3}$ is obtained as a statistically steady spectrum, where k is the magnitude of the two-dimensional wave-number vector \mathbf{k} , i.e., $k = |\mathbf{k}|$. Contrarily, energy spectra such as $\mathcal{E}(k) \propto k^{-0.2}$ and k^0 are experimentally observed [4–6]. Moreover, the existence of the spectra with the energy cascade $\mathcal{E}(k) \propto k^{-1}$ and with the wave action cascade $\mathcal{E}(k) \propto k^{-1/3}$ is predicted by the dimensional analysis [7]. The energy spectrum obtained by the numerical simulation is the steady solution of the kinetic equation of the weak turbulence theory, while the energy spectra obtained in the experiments and by the dimensional analysis are outside the scope of the weak turbulence theory.

In our earlier study [8], we provided a unified perspective on the variability of the spectra by performing direct numerical simulations using the Föppl-von Kármán equation and found that the strength of the nonlinearity causes the variability.

Recently, it was shown that the experimentally observed spectrum may be explained by the strong dissipation in the small frequencies [9]. The boundary conditions and homogeneity of the plate may also affect the energy spectra, as pointed out in Ref. [10]. We are interested in the universal dynamics in the inertial subrange, in which the external force and dissipation do not affect the dynamics. The subject concerned in this manuscript is an extension of the research in Ref. [3], whose attractive subtitle is “Can one hear a Kolmogorov spectrum?”

In Ref. [8], the weak turbulence spectrum was obtained at low-energy levels, while the strong turbulence spectrum, which is similar to the spectrum of the wave action cascade, was obtained at high-energy levels. Moreover, the coexistence of the spectra, in which the weak and strong turbulence spectra, respectively, appear in small and large wave numbers, was obtained at intermediate energy levels. The numerical simulations with larger mode numbers (Fig. 1) show that if the inertial subrange is large enough, the coexistence state is generally achieved. This generality was not clearly seen in Fig. 2 of Ref. [8]. In Fig. 1, the *separation* wave numbers between the two types of the spectra lie on a straight line, which suggests that there exists a simple law.

A similar bending structure of the energy spectra is numerically observed using another type of external force and dissipation [11]. This fact suggests that the bending structure is universal. The weak turbulence spectra in large wave numbers are commonly observed in the numerical simulations [3,8,11]. It should be possible to estimate the separation wave numbers by balancing the linear and nonlinear time scales owing to the following reason: the weak turbulence theory is not applicable in the wave-number range in which the nonlinear time scales of the energy transfer are comparable with or shorter than the linear time scales [12,13]. Note that the linear time scale is comparable with the nonlinear time scale at a certain scale in the statistically isotropic system of the elastic waves; this is distinct from the comparability of the time scales over a wide

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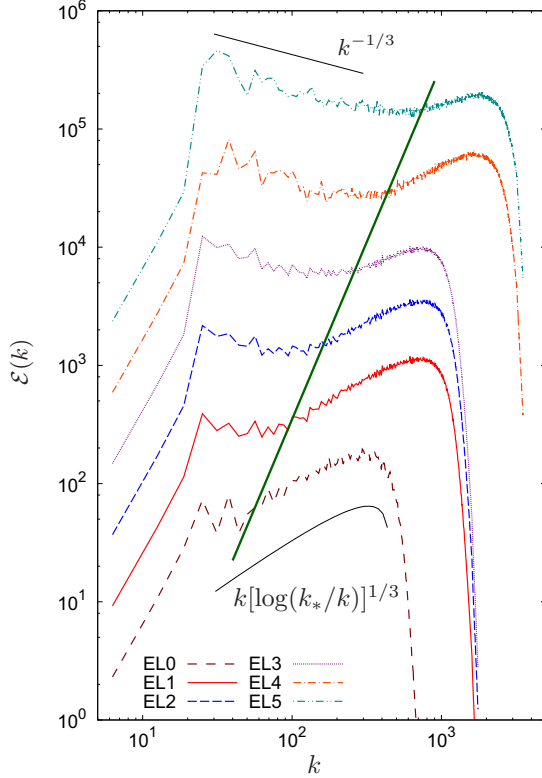


FIG. 1. (Color online) Energy spectra at various energy levels, EL0 to EL5, in statistically steady states. The solid line indicates the separation between the weak and strong turbulence spectra. The weak turbulence spectrum $k[\log(k_*/k)]^{1/3}$, where $k_* = 144\pi$, and a power-law spectrum $k^{-1/3}$ are shown for reference.

range of wave numbers in the systems reported as the critical balance.

In the experiment of a thin elastic plate [14], the nonlinear dispersion relation obtained from the space-time Fourier spectra indicates a scaling law that is a function of the energy flux. However, the coexistence of the weak and strong turbulence spectra as well as the balance between the linear and nonlinear time scales was not investigated. The nonlinear time scales are usually estimated by the nonlinear-interaction times according to the weak turbulence theory in Refs. [12,15,16]. However, because the unsteadiness of a system is required for the finiteness of the nonlinear-interaction times, the nonlinear-interaction times cannot be used in statistically steady states. Therefore, in the present analysis, the nonlinear time scales are evaluated using the nonlinear frequency shifts, which are the difference between the linear and nonlinear frequencies, instead of the nonlinear-interaction times. The nonlinear frequencies are derived by considering self-nonlinear interactions in Refs. [12,15,16]. The present system, however, requires an extension of this derivation since it has the $1 \leftrightarrow 3$ and $3 \leftrightarrow 1$ interactions as well as the $2 \leftrightarrow 2$ interactions. The *companion* elementary wave, which has the wave number with the opposite sign, should be taken into account for the weakly nonlinear expansion to be consistent in our derivation. The interactions are referred to as “extended self-nonlinear interactions” in this paper.

In this paper, we propose the analytical representation to evaluate the separation wave numbers from the viewpoint of

the time scales. First, we show that the nonlinear frequencies estimated by the extended self-nonlinear interactions agree well with the representative nonlinear frequencies obtained from the time series in our simulations. Then, we show that the *critical* wave numbers at which the frequency shift in the nonlinear frequency is comparable with the linear frequency agree well with the separation wave numbers. The validity of our analytical representation is confirmed by other series of numerical simulations in which the material parameters of elastic plates are changed.

II. FORMULATION OF ELASTIC WAVES

A. Fourier representation

The dynamics of the elastic waves propagating in a thin plate is described by the Föppl-von Kármán (FvK) equation [17]. Under the periodic boundary condition, the FvK equation is written using the Fourier coefficient of the displacement $\zeta_k(t)$ and that of the momentum $p_k(t)$ as follows:

$$\frac{dp_k}{dt} = -\frac{Eh^2}{12(1-\sigma^2)}k^4\zeta_k - \frac{E}{2} \sum_{k=k_1+k_2+k_3} \frac{|\mathbf{k} \times \mathbf{k}_1|^2 |\mathbf{k}_2 \times \mathbf{k}_3|^2}{|\mathbf{k}_2 + \mathbf{k}_3|^4} \zeta_{k_1} \zeta_{k_2} \zeta_{k_3}, \quad (1a)$$

$$\frac{d\zeta_k}{dt} = \frac{p_k}{\rho}, \quad (1b)$$

where E , σ , ρ , and h are the Young’s modulus, Poisson ratio, density, and thickness of an elastic plate, respectively.

The FvK equation (1) can be rewritten for the complex amplitude

$$a_k(t) = \frac{\rho\omega_k\zeta_k(t) + ip_k(t)}{\sqrt{2\rho\omega_k}}, \quad (2)$$

as

$$\frac{da_k}{dt} = -i\omega_k a_k - \frac{iE}{8\rho^2} \sum_{k=k_1+k_2+k_3} \frac{|\mathbf{k} \times \mathbf{k}_1|^2 |\mathbf{k}_2 \times \mathbf{k}_3|^2}{|\mathbf{k}_2 + \mathbf{k}_3|^4} \times \frac{(a_{k_1} + a_{-k_1}^*)(a_{k_2} + a_{-k_2}^*)(a_{k_3} + a_{-k_3}^*)}{\sqrt{\omega_k\omega_{k_1}\omega_{k_2}\omega_{k_3}}}. \quad (3)$$

If the nonlinear terms were absent, a_k would rotate clockwise in the phase space with the linear frequency

$$\omega_k = \sqrt{\frac{Eh^2}{12(1-\sigma^2)\rho}} k^2. \quad (4)$$

Thus, the complex amplitude a_k represents the elementary wave with the wavevector \mathbf{k} and plays a central role in the weak turbulence theory. The azimuthally integrated energy spectrum is defined as $\mathcal{E}(k) = (\Delta k)^{-1} \sum_{k-\Delta k/2 \leq |k| < k+\Delta k/2} \omega_k |a_k|^2$, where Δk is the bin width used to obtain the spectrum.

B. Formulation of simulation

Direct numerical simulations (DNS) were performed using the Fourier-spectral method according to Eq. (3). To obtain

statistically steady states, we added the external force \mathcal{F}_k , which is effective at small wave numbers, and dissipation \mathcal{D}_k , which is effective at large wave numbers, and the governing equation can be rewritten as

$$\frac{da_k}{dt} = -i\omega_k a_k + \mathcal{N}_k + \mathcal{F}_k + \mathcal{D}_k. \quad (5)$$

Here, \mathcal{N}_k symbolically expresses the four-wave nonlinear interactions in Eq. (3).

The external force \mathcal{F}_k is provided so that the absolute values of the complex amplitudes in the forced wave numbers $|k| \leq 8\pi$ are kept constant. The dissipation \mathcal{D}_k is provided as the eighth-order hyperviscosity to achieve larger inertial subranges. The pseudospectral method with the aliasing removal by the 4/2 law is used to obtain \mathcal{N}_k under the periodic boundary condition. The numbers of the Fourier modes are $N_{\text{mode}}^2 = 512^2$ (256² alias-free) for EL0, $N_{\text{mode}}^2 = 1024^2$ (512² alias-free) for EL1–EL3, and $N_{\text{mode}}^2 = 2048^2$ (1024² alias-free) for EL4 and EL5. The fourth-order Runge-Kutta method with a time step of 10^{-7} s is used for the time integration. The time step should be small enough to resolve the fast dynamics of a_k at the large wave numbers, which is governed by the linear dispersion relation [Eq. (4)], to reproduce the energy transfer due to the resonant interactions. The parameters of the steel plate which are shown in Ref. [4] were adopted as the standard values. The details of the numerical method are explained in Ref. [8].

By changing the magnitude of the external force \mathcal{F}_k , the total energy in the system can be controlled. The energy spectra are shown in Fig. 1. Three energy levels, EL1, EL2, and EL3, were selected as the representatives of the simulations to support our analytical results because of the limited amount of computational time. The ratio of the largest time scale to the smallest one is of the order $N_{\text{mode}}^2 \sim 10^6$ because of the linear frequency [Eq. (4)].

III. STRATEGY FOR IDENTIFYING SEPARATION WAVE NUMBER

Let us find an analytical representation to identify the separation wave number between the weak and strong turbulence spectra. Our analysis begins with the fact that the spectra in large wave numbers represent the weak turbulence spectra, which are consistent with the weak turbulence theory. For the weak turbulence theory to be applicable, the linear time scale should be much smaller than the nonlinear time scale. Therefore, the separation wave number is expected to be the critical wave number at which the nonlinear time scale is comparable with the linear time scale.

In the studies of unsteady wave turbulence, the nonlinear time scale is often estimated using the nonlinear-interaction time $n_k/(dn_k/dt)$, where n_k is the wave action defined as $n_k \delta_{k,k'} = \langle a_k a_{k'}^* \rangle$. (See, for example, Refs. [12, 15].) It should be noted that $n_k/(dn_k/dt)$ has a finite value only in a statistically unsteady state, although it describes the time scale of the net energy transfer. One might think that the self-similarity and the dimensional analysis can be used for the applicability-limit scale for the weak turbulence theory as is conventionally done [e.g. 12, 15, 18]. This idea is based on the fact that $n_k/(dn_k/dt)$, whose power-law dependence on wave

numbers is different from that of the linear dispersion, must cross over the linear time scale. However, the crossover scale depends not only on the power-law exponents of the time scales with respect to the wave numbers but also on their coefficients. Although the power-law dependence of $n_k/(dn_k/dt)$ on the wave numbers can be obtained by using the self-similarity and/or the dimensional analysis, the coefficient cannot be found within this framework. Furthermore, it cannot remain finite in the statistically steady states.

The energy flux P is generally used by employing the relation $d\mathcal{E}(k)/dt = -dP/dk$ for steady states of single turbulence [18]. However, it is not appropriate to use P in our present purpose to estimate the separation scale between the two types of turbulence, although P is used in Ref. [5]. It is difficult to expect that a single value of P can represent the energy flux over the entire range of wave numbers in the present system since the turbulence shows different characteristics on both sides of the separation wave number. Furthermore, since the Kolmogorov-Zakharov spectrum has the same self-similarity as the Rayleigh-Jeans equilibrium spectrum, which corresponds to the energy equipartition, the energy flux is zero at the first-order analysis of the weak turbulence theory [3].

Our strategy adopted here is that the nonlinear time scales are evaluated not by the energy transfer but instead by the frequency shift due to the extended self-nonlinear interactions. This implies that it is independent of the weak turbulence theory, which is a statistical theory; however, it depends on the weakly nonlinear expansion. In this sense, our estimation for the nonlinear time scale is different from that in Refs. [12, 16, 18], which is estimated by the resonant interactions. Then, the frequency shift is balanced to the linear frequency to estimate the separation wave number. In fact, this balance corresponds to equating the magnitudes of the linear and nonlinear terms on the right-hand side of Eq. (3).

IV. NONLINEAR FREQUENCY

A. Estimation by extended self-nonlinear interactions

Let us estimate the nonlinear frequencies to evaluate the nonlinear time scales. Among the nonlinear interactions, the combination of the wave numbers, including the self-nonlinear interactions in which one of \mathbf{k}_1 , \mathbf{k}_2 , and \mathbf{k}_3 appearing in Eq. (3) is \mathbf{k} , is confined to $(\mathbf{k}, \mathbf{k}_1, \mathbf{k}_2, \mathbf{k}_3) = (\mathbf{k}, \mathbf{k}', \mathbf{k}, -\mathbf{k}')$ or $(\mathbf{k}, \mathbf{k}', -\mathbf{k}', \mathbf{k})$. These self-nonlinear interactions are dominant if the nonlinearity is weak. Then, the FvK equation (3) can be rewritten as

$$\frac{da_k}{dt} = -i\omega_k a_k - i\omega_k^s (a_k + a_{-k}^*) + \mathcal{N}_k^{\text{nn}}, \quad (6)$$

where

$$\begin{aligned} \omega_k^s &= \frac{E}{4\rho^2} \sum_{k'} \frac{|\mathbf{k} \times \mathbf{k}'|^4}{|\mathbf{k} - \mathbf{k}'|^4} \\ &\quad \times \frac{|a_{k'}|^2 + |a_{-k'}|^2 + a_k a_{-k'} + a_{k'}^* a_{-k'}^*}{\omega_k \omega_{k'}} \\ &= \frac{E}{2\rho \omega_k} \sum_{k'} \frac{|\mathbf{k} \times \mathbf{k}'|^4}{|\mathbf{k} - \mathbf{k}'|^4} |\zeta_{k'}|^2. \end{aligned} \quad (7)$$

Equation (7) represents the frequency increment from ω_k due to the self-nonlinear interactions. Note that ω_k^s is nonnegative by definition. The nonself-nonlinear interactive term $\mathcal{N}_k^{\text{nn}}$ is defined as

$$\mathcal{N}_k^{\text{nn}} = -\frac{iE}{8\rho^2} \sum_{k=k_1+k_2+k_3, k_1, k_2, k_3 \neq k} \frac{|\mathbf{k} \times \mathbf{k}_1|^2 |\mathbf{k}_2 \times \mathbf{k}_3|^2}{|\mathbf{k}_2 + \mathbf{k}_3|^4} \times \frac{(a_{k_1} + a_{-k_1}^*)(a_{k_2} + a_{-k_2}^*)(a_{k_3} + a_{-k_3}^*)}{\sqrt{\omega_k \omega_{k_1} \omega_{k_2} \omega_{k_3}}}.$$

It should be noted that the companion mode for a_k , which has the wave number with the opposite sign a_{-k}^* , in the self-nonlinear interaction term of Eq. (6) appears in the first order for consistency in the weakly nonlinear expansion. In the weak turbulence systems that contain only the $2 \leftrightarrow 2$ interactions, $\omega_k + \omega_k^s$ is the nonlinear frequency [12,15,16]. However, the companion mode must be taken into account owing to the $1 \leftrightarrow 3$ and $3 \leftrightarrow 1$ interactions in the nonlinear term in Eq. (3).

Equation (6) can be rewritten in the following simultaneous equations:

$$\frac{d}{dt} \begin{pmatrix} a_k \\ a_{-k}^* \end{pmatrix} = \begin{pmatrix} -i(\omega_k + \omega_k^s) & -i\omega_k^s \\ i\omega_{-k}^s & i(\omega_{-k} + \omega_{-k}^s) \end{pmatrix} \begin{pmatrix} a_k \\ a_{-k}^* \end{pmatrix} + \begin{pmatrix} \mathcal{N}_k^{\text{nn}} \\ \mathcal{N}_{-k}^{\text{nn}*} \end{pmatrix}. \quad (8)$$

Because the extended self-nonlinear interactions preserve the inversion symmetry of the system, the frequency increment satisfies the relation $\omega_k^s = \omega_{-k}^s$. The eigenvalues of the matrix in the right-hand side of Eq. (8) are

$$\pm \omega_k^{\text{NL}} = \pm \omega_k \sqrt{1 + \frac{2\omega_k^s}{\omega_k}}. \quad (9)$$

If the remnant interaction term ($\mathcal{N}_k^{\text{nn}} \mathcal{N}_{-k}^{\text{nn}*}$) in Eq. (8) can be neglected, the solution is

$$a_k(t) = a_k(0) \left(\cos \omega_k^{\text{NL}} t - i \frac{\omega_k + \omega_k^s}{\omega_k^{\text{NL}}} \sin \omega_k^{\text{NL}} t \right) - i a_{-k}^*(0) \frac{\omega_k^s}{\omega_k^{\text{NL}}} \sin \omega_k^{\text{NL}} t, \quad (10)$$

which results in

$$\zeta_k(t) = \zeta_k(0) \cos \omega_k^{\text{NL}} t. \quad (11)$$

In other words, although $\omega_k^{\text{NL}} \approx \omega_k + \omega_k^s$ for $\omega_k^s \ll \omega_k$, ω_k^{NL} is preferable to $\omega_k + \omega_k^s$, which is considered as the nonlinear frequency in Refs. [12,15,16]. The difference in the derivation of the nonlinear frequencies in these references and in this paper is whether the companion mode is taken into account.

It should be noted that the evaluation of ω_k^s using Eq. (7) requires each Fourier component of the displacement $|\zeta_k|^2$, which is time-dependent in general. However, because the frequencies must be constant in time to be meaningful, the spectrum of the displacement should be replaced by its ensemble- or time-average $\langle |\zeta_k|^2 \rangle$. In the statistically steady state, the averaged $\omega_k^{\text{NL}} (\geq \omega_k)$ is expected to be a good approximation of the nonlinear frequency.

B. Comparison with numerically obtained representative frequency

The frequency spectrum $|\tilde{a}_k(\Omega)|^2$ is obtained from the Fourier transforms of the time series of $a_k(t)$. Note that $|\tilde{a}_k(\Omega)|^2$ lacks the symmetry with respect to the sign of Ω , since $a_k(t)$ is a complex number. Figure 2 shows the contours of the frequency spectra $|\tilde{a}_k(\Omega)|^2$. The horizontal and vertical axes, which are scaled logarithmically, show the wave number

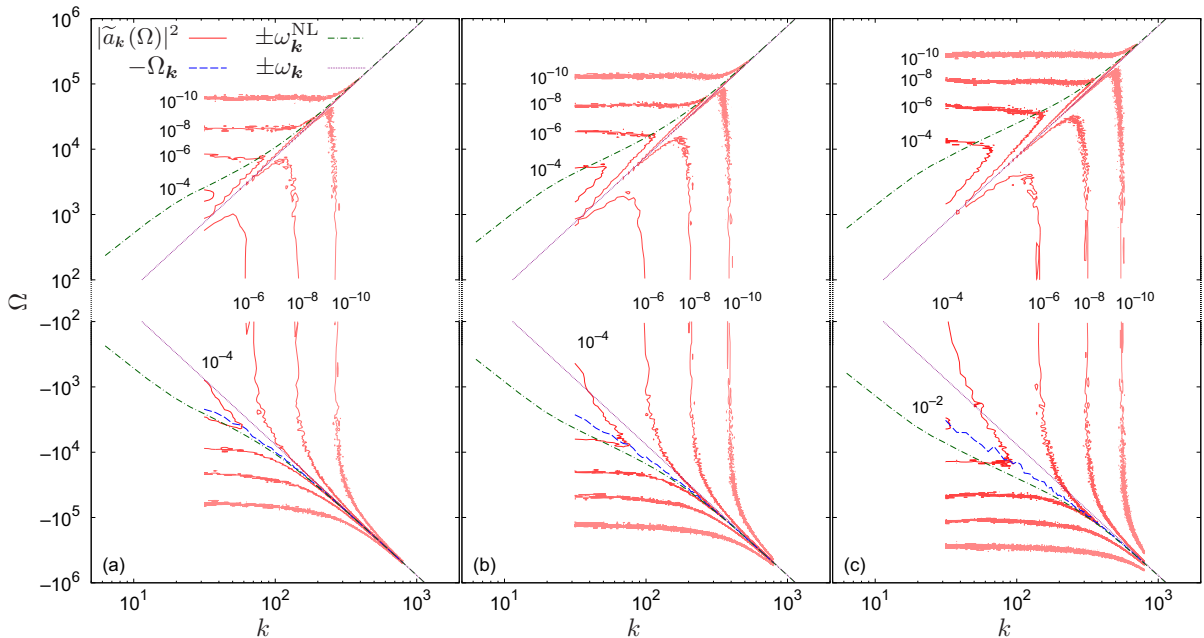


FIG. 2. (Color online) Contours of frequency spectra $|\tilde{a}_k(\Omega)|^2$ for three energy levels: (a) EL1, (b) EL2, and (c) EL3. The contours are drawn with the (red) solid curves. The representative frequencies Ω_k , nonlinear frequencies ω_k^{NL} , and linear frequencies ω_k are, respectively, indicated by the (blue) broken, (green) dashed-dotted, and (purple) dotted curves.

and frequency, respectively. Note that the negative frequency is also scaled logarithmically, and the figures for the positive and negative frequencies are arranged vertically with gaps. The contours (red curves) are drawn only in the inertial subrange $8\pi < k < 256\pi$. The cross-sections of these contours will exhibit a similar structure to those in Fig. 3 of Ref. [8]. (See Appendix A for the structure of the frequency spectrum.)

The *representative* nonlinear frequency Ω_k for each k is defined such that $|\tilde{a}_k(-\Omega_k)|^2$ is maximal. The negative sign is introduced in the argument because the maxima of $|\tilde{a}_k(\Omega)|^2$ always appear in the negative region of Ω . Therefore, $\Omega = -\Omega_k$ corresponds to the ridge curve of the contours of $|\tilde{a}_k(\Omega)|^2$. Note that $|\tilde{a}_k(\Omega)|^2 \propto \delta(\Omega + \omega_k)$ and $\Omega_k = \omega_k^{\text{NL}} = \omega_k$ if the nonlinear interactions were zero. In Fig. 2, the ridges of the frequency spectra are drawn only in the inertial subrange with the (blue) broken curves. The nonlinear frequencies ω_k^{NL} , which are obtained from Eqs. (7) and (9) by averaging ten sets of data of ζ_k , as well as the linear frequencies [Eq. (4)] are drawn over the entire range of wave numbers.

The structure of the contour curves drastically changes near the linear dispersion line $\Omega = \pm\omega_k$. The asymptotic behavior of the contour curves in $|\Omega| < \omega_k$ ($|\Omega| > \omega_k$) is vertical (horizontal), which means that the spectral amplitude $|\tilde{a}_k(\Omega)|^2$ is nearly constant along the constant wave number (frequency) lines. The hornlike extensions of the contour curves to the upper right along ω_k^{NL} and to the lower right along $-\omega_k^{\text{NL}}$ correspond to the humps of $|\tilde{a}_k(\Omega)|^2$, whereas the extension to the lower left along ω_k corresponds to the depressions. From the angles of the curves in Figs. 2(a)–2(c), one can see that the humps and depressions broaden for higher energy levels. The widths between the contour curves in Figs. 2(a)–2(c) are also larger as the energy levels increase, and it corresponds to a broadening of the frequency spectrum.

The negative frequency region is a primary interest to evaluate the nonlinear time scales. For the large wave numbers in the negative frequency region, where the nonlinearity is weak [8], the representative nonlinear frequencies Ω_k are close to the linear frequencies ω_k and the nonlinear frequencies ω_k^{NL} . Owing to the weak nonlinearity, the frequency spectra are sharp around the maximum frequency, and Ω_k and ω_k^{NL} are barely distinguishable from ω_k .

For the small wave numbers in the negative frequency region, where the nonlinearity is relatively strong, the frequency spectra are broad, and the nonlinear frequencies ω_k^{NL} deviate from the linear frequencies ω_k to the higher frequencies. The representative nonlinear frequencies Ω_k lie between ω_k^{NL} and ω_k . Comparing the three figures, Figs. 2(a)–2(c), we observe that the difference among these three frequencies, Ω_k , ω_k , and ω_k^{NL} , is larger for a wave number k as the energy levels increase.

In the positive frequency regions shown in Fig. 2, we found the secondary maxima of the frequency spectra near the nonlinear frequencies ω_k^{NL} (see Appendix A). Our results show the two eigenvalues [Eq. (9)] agree quite well with the primary maximum frequencies and the secondary maximum frequencies in the large wave numbers. The secondary maxima of the frequency spectra can be observed even in the weakly nonlinear regime, which can be accounted for by the coupling with the companion mode via the $1 \leftrightarrow 3$ and $3 \leftrightarrow 1$

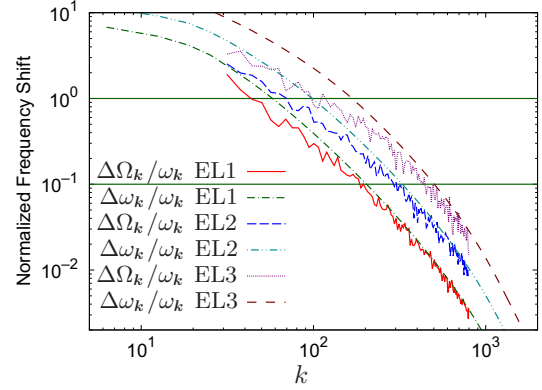


FIG. 3. (Color online) Nonlinear frequency shifts normalized by the linear frequencies for three energy levels. The two horizontal lines represent the normalized frequency shifts equal to 10^{-1} and 1.

interactions. It is a different mechanism from the strong nonlinearity [19] to make the secondary maxima.

To observe the nonlinear frequency shifts more clearly, the nonlinear frequency shifts normalized by the linear frequencies are illustrated in Fig. 3. The frequency shifts due to the extended self-nonlinear interactions and those of the representative frequency are, respectively, obtained as $\Delta\omega_k = \omega_k^{\text{NL}} - \omega_k$ and $\Delta\Omega_k = \Omega_k - \omega_k$. The frequency shifts $\Delta\omega_k$ agree quite well with $\Delta\Omega_k$ at low energies and large wave numbers, where the nonlinearity is weak. Conversely, at high energies and small wave numbers, the frequency shifts $\Delta\omega_k$ overestimate the actual frequency shifts since the remnant interactions absorb the frequency shift due to the extended self-nonlinear interactions. The nonlinear frequency $\omega_k^{\text{NL}} = \omega_k \sqrt{1 + 2\omega_k^s / \omega_k}$ agrees with Ω_k better than $\omega_k + \omega_k^s$ at small wave numbers, at which ω_k^s is large.

The deviation of $\Delta\Omega_k / \omega_k$ from $\Delta\omega_k / \omega_k$ appears in the wave numbers where $\Delta\Omega_k / \omega_k$ and $\Delta\omega_k / \omega_k$ are between 10^{-1} and 1. Since $\Delta\omega_k$ is derived from the extended self-nonlinear interactions, the deviation of $\Delta\Omega_k / \omega_k$ from $\Delta\omega_k / \omega_k$ is caused by the broad-mode interactions, including strongly nonlinear interactions. Namely, the deviation between 10^{-1} and 1 indicates that the strongly nonlinear effects are no longer negligible in this region. Therefore, the ratio $\Delta\omega_k / \omega_k$ can be understood as the measure of the nonlinearity for each wave number. In other words, the separation wave number between the weak and strong turbulence, which is expected to be the application limit of the weak turbulence theory, is determined by the critical wave number where the ratio of the frequencies is between 10^{-1} and 1.

C. Identification of boundary wave number

The weak turbulence theory is not applicable when the nonlinear time scale is comparable with the linear time scale. Therefore, the critical wave number $k_c = |k_c|$ can be defined as

$$\Delta\omega_{k_c} = \epsilon\omega_{k_c}, \quad (12)$$

where ϵ is a constant between 10^{-1} and 1, as shown in Fig. 3.

To find the analytical representation of the critical wave number k_c for arbitrary energy levels, it is convenient to use

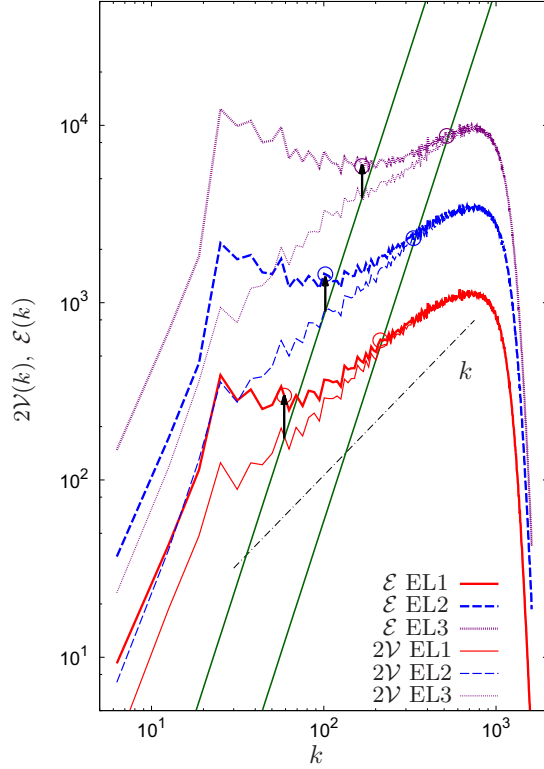


FIG. 4. (Color online) Energy spectra (thick curves) and doubled linear potential energy spectra (thin curves) for three energy levels. The left and right inclined parallel lines represent $2\check{\mathcal{V}}(k_c; 1)$ and $2\check{\mathcal{V}}(k_c; 10^{-1})$, respectively. The circles indicate the position of the energy spectra at the wave numbers at which the linear potential energy spectra intersect with $2\check{\mathcal{V}}(k_c; 1)$ and with $2\check{\mathcal{V}}(k_c; 10^{-1})$.

the linear potential energy $\mathcal{V}(k)$. As shown in the previous two subsections, a good estimation of ω_k^{NL} as well as ω_k^s is obtained by its ensemble- or time-average $\langle |\zeta_k|^2 \rangle$. Furthermore, as derived in Appendix B, $\langle |\zeta_k|^2 \rangle \approx 4\pi \mathcal{V}(k) / (\rho k \omega_k^2)$ in isotropic fields. Therefore, the averaged ω_k^{NL} is obtained from $\mathcal{V}(k)$.

It is assumed that the linear potential energy has the self-similar form $\mathcal{V}(k) = Ck$. The self-similarity of $\mathcal{V}(k)$ over the entire inertial subrange for all the energy levels will be seen in Fig. 4, which is a distinctive feature from the energy spectra $\mathcal{E}(k)$. Equations (7), (9), and (12) give the analytical representation of the critical wave number with the parameter C corresponding to the energy level as

$$k_c = \frac{6(1 - \sigma^2)}{h^2} \sqrt{\frac{3C}{\epsilon(\epsilon + 2)E}} \quad (13)$$

(see Appendix C for the derivation).

There exists a more convenient representation to graphically find the value of k_c . Equation (13) can be rewritten as

$$\mathcal{V}(k_c) = \frac{\epsilon(\epsilon + 2)Eh^4}{108(1 - \sigma^2)^2} k_c^3 \equiv \check{\mathcal{V}}(k_c; \epsilon). \quad (14)$$

To illustrate the validity of Eq. (14), the energy spectra $\mathcal{E}(k)$ and twice the linear potential energy spectra $2\mathcal{V}(k)$ are drawn in Fig. 4. The two inclined parallel lines in Fig. 4 represent $2\check{\mathcal{V}}(k_c; 10^{-1})$ and $2\check{\mathcal{V}}(k_c; 1)$. The critical wave number k_c for

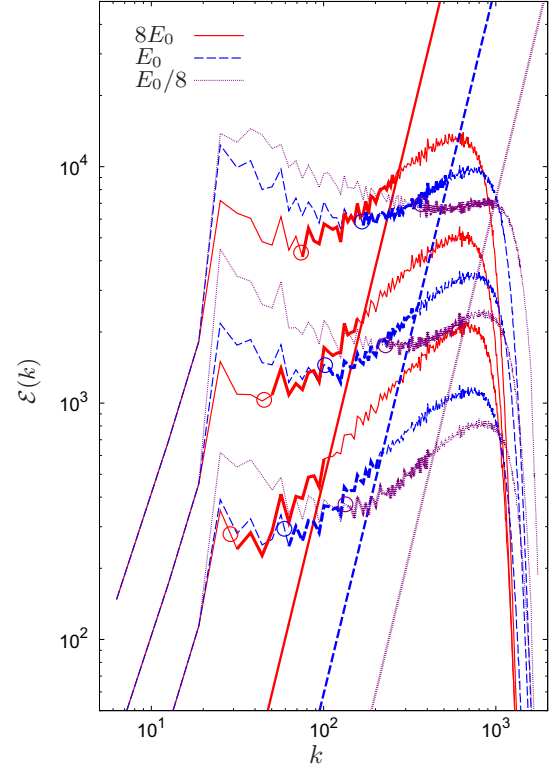


FIG. 5. (Color online) Energy spectra for different values of the Young's modulus. The straight lines show $2\check{\mathcal{V}}(k_c; 10^{-1})$ for each Young's modulus. The energy spectra between k_c for $\epsilon = 10^{-1}$ and that for $\epsilon = 1$ are indicated by the thick curves, and the circles indicate $\mathcal{E}(k_c)$, where k_c is determined by $\mathcal{V}(k_c) = \check{\mathcal{V}}(k_c; 1)$.

ϵ can be found at the intersection of $2\mathcal{V}(k)$ and $2\check{\mathcal{V}}(k_c; \epsilon)$. Then, our estimation predicts the separation wave number is located between the critical wave number for $\epsilon = 10^{-1}$ and that for $\epsilon = 1$. In other words, the wave numbers at which $\mathcal{E}(k)$'s bend are located between k_c for $\epsilon = 10^{-1}$ and that for $\epsilon = 1$ for all the energy levels. In the larger wave-number side of the line $2\check{\mathcal{V}}(k_c; 10^{-1})$, we find that $\mathcal{E}(k) \approx 2\mathcal{V}(k)$ and that the system is weakly nonlinear. Conversely, in the smaller wave-number side of the line $2\check{\mathcal{V}}(k_c; 1)$, we found that $\mathcal{E}(k)$ is far from $2\mathcal{V}(k)$ and that a_k cannot be approximated by the harmonic oscillation.

To verify Eq. (14), other series of DNS, in which the material parameters such as the Young's modulus E and thickness h of the elastic plates are artificially changed, were performed. The energy spectra for different values of the Young's modulus are shown in Fig. 5. Because we used $8E_0$ and $E_0/8$ for the Young's modulus, the critical wave numbers should be $2k_{c0}$ and $k_{c0}/2$ according to Eq. (14), where E_0 is the Young's modulus of the originally used steel plate and k_{c0} is the critical wave number for $E = E_0$. Clearly, Fig. 5 supports that the right-hand side in Eq. (14) is proportional to $E k_c^3$.

Similarly, the energy spectra for different values of the thickness h are shown in Fig. 6. Because we used $2^{3/4}h_0$ and $h_0/2^{3/4}$ for the thickness, the critical wave numbers should be $2k_{c0}$ and $k_{c0}/2$ again, where h_0 is the originally used thickness. This also supports that the coefficient in Eq. (14) is proportional to $h^4 k_c^3$. We also confirmed that the critical wave number does not depend on the density ρ , although the figures

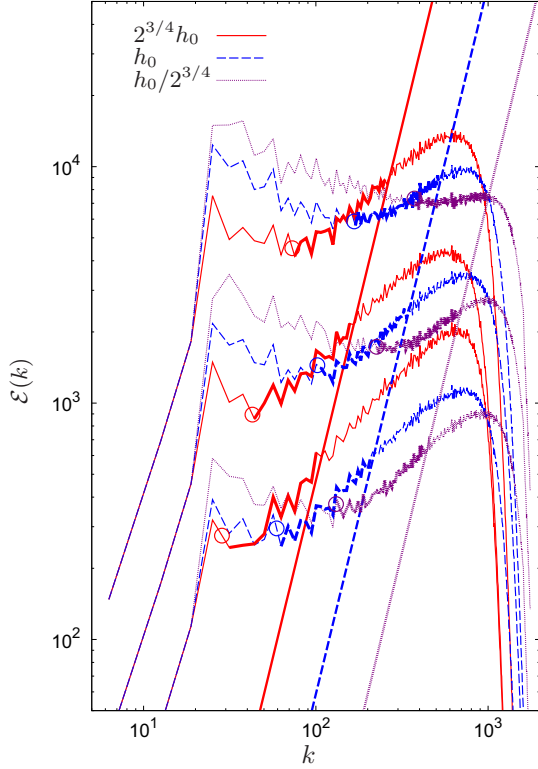


FIG. 6. (Color online) Energy spectra for different values of the thickness of the plate. The straight lines show $2\check{\mathcal{V}}(k_c; 10^{-1})$ for each thickness. The energy spectra between k_c for $\epsilon = 10^{-1}$ and that for $\epsilon = 1$ are indicated by the thick curves, and the circles indicate $\mathcal{E}(k_c)$, where k_c is determined by $\mathcal{V}(k_c) = \check{\mathcal{V}}(k_c; 1)$.

are omitted here. Note that we do not examine the dependence of the critical wave number on the Poisson ratio σ since the variation of the Poisson ratio σ among elastic media is small. In Figs. 4–6, we have demonstrated the validity of the estimation of the separation wave number according to Eq. (14) including its coefficients.

V. CONCLUSION

In this paper, we have obtained the analytical representation of the critical wave numbers at which the nonlinear frequency shifts are comparable with the linear frequencies. In the derivation of the nonlinear frequency, the extended self-nonlinear interactions, including the coupling with the companion elementary wave, are explicitly evaluated. The coupling with the companion elementary wave must be taken into account generally in the weak turbulence systems containing the $1 \leftrightarrow 3$ and $3 \leftrightarrow 1$ interactions as well as the $2 \leftrightarrow 2$ interactions. The nonlinear frequencies evaluated by the extended self-nonlinear interactions agree well with the representative nonlinear frequencies determined by the maxima of the frequency spectra obtained in the direct numerical simulations. The agreement is remarkable at large wave numbers, in which the weak turbulence spectrum is observed.

The analytically obtained critical wave number successfully reproduced the separation wave number between the weak and strong turbulence spectra. It is consistent with the view that the weak turbulence theory is not applicable for wave numbers at

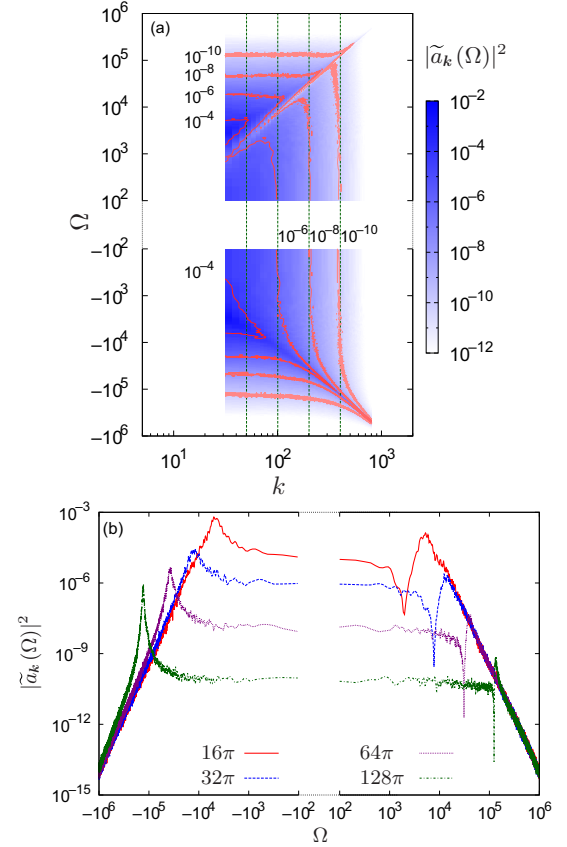


FIG. 7. (Color online) (a) Density plots and contours of the frequency spectra $|\tilde{a}_k(\Omega)|^2$ for EL2. The vertical straight lines show $k = 16\pi, 32\pi, 64\pi,$ and 128π . (b) Frequency spectra for $k = 16\pi, 32\pi, 64\pi,$ and 128π for EL2. The positive and negative frequencies are arranged horizontally with the gap.

which the nonlinear time scales are comparable with the linear time scales. The agreement between the critical wave numbers and the separation wave numbers as well as the dependence of the wave numbers on the material parameters were also confirmed.

The coexistence of the weak and strong turbulence spectra implies that the large-scale strongly nonlinear coherent structures coexist with a large number of small-scale weakly nonlinear random waves in the real space. The characterization of each structure is of great importance. The relations of this real-space coexistence to the different properties between $\mathcal{E}(k)$ and $\mathcal{V}(k)$ appearing in Fig. 4 are now under consideration. We will also identify the energy transfer among waves in the coexistence.

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APPENDIX A : STRUCTURES OF FREQUENCY SPECTRA

The density plots and contours of the frequency spectra for EL2 are shown in Fig. 7(a). It corresponds to Fig. 2(b), where

only the contours are drawn for the spectra. The ridge of the frequency spectra in the negative frequencies corresponds to the cusps of the contours which extend to the lower right. Similarly, the ridge and trough in the positive frequencies correspond to the cusps that extend to the upper right and lower left, respectively. To see the ridges and trough more clearly, the frequency spectra for $k = 16\pi$, 32π , 64π , and 128π on the line parallel to the vertical axis at each k are shown in Fig. 7(b). Humps can be observed in the positive and negative frequency regions of Ω ; the hump in the negative frequency region is larger than that in the positive frequency region. The null amplitude $|\tilde{a}_k(\omega_k)|^2 = 0$ produces a slit in the right humps. This structure corresponds to the hornlike extension of the contour curves to the lower right along ω_k in Fig. 2. The secondary maximum frequencies are located near ω_k^{NL} .

The $|\tilde{a}_k(\omega_k)|^2 = 0$ property is a direct consequence of the definition of a_k . The frequency-space representation of Eq. (1b), $\tilde{p}_k(\Omega) = i\rho\Omega\tilde{\zeta}_k(\Omega)$, and the definition of the complex amplitude, Eq. (2), lead to the relation $|\tilde{a}_k(\Omega)|^2 = \rho(\omega_k - \Omega)^2|\tilde{\zeta}_k(\Omega)|^2/(2\omega_k)$. We have checked these relations in our simulations, though the figures are omitted here.

APPENDIX B : SPECTRA OF LINEAR POTENTIAL ENERGY AND OF DISPLACEMENT

In the present study, the periodic boundary condition with a period of $L = 1$ m is used. The Fourier series is defined as

$$F_k = \frac{1}{L^2} \int_0^L \int_0^L dx f(\mathbf{x}) \exp(-i\mathbf{k} \cdot \mathbf{x}),$$

$$f(\mathbf{x}) = \sum_k F_k \exp(i\mathbf{k} \cdot \mathbf{x}),$$

where $f(\mathbf{x})$ is a physical quantity such as the displacement $\zeta(\mathbf{x})$ and momentum $p(\mathbf{x})$ in the real space, and F_k is its Fourier coefficient. The azimuthally integrated linear potential energy spectrum is given as

$$\mathcal{V}(k)\Delta k = \sum_{k-\Delta k/2 \leq |k| < k+\Delta k/2} \frac{\rho\omega_k^2}{2} |\zeta_k|^2,$$

where Δk is the bin width used to obtain the spectrum; Δk is set to $2\pi/L$ in the present study. The summation in the two-dimensional wave-number space can be replaced by the integral as

$$\sum_k \approx \int_0^\infty dk \int_0^{2\pi} d\theta \frac{k}{\Delta k_x \Delta k_y}.$$

In a statistically isotropic system, the ensemble- and/or time-averaging $\langle \dots \rangle$ is performed to evaluate the spectra:

$$\mathcal{V}(k)\Delta k \approx \frac{2\pi k \Delta k}{\Delta k_x \Delta k_y} \frac{\rho\omega_k^2}{2} \langle |\zeta_k|^2 \rangle,$$

where $\Delta k_x = \Delta k_y = 2\pi/L$ is the grid spacing.

APPENDIX C : FREQUENCY DUE TO SELF-INTERACTIONS AND LINEAR POTENTIAL ENERGY SPECTRA

In an isotropic system, the frequency due to the self-interactions can be rewritten as

$$\omega_k^s = \frac{E}{2\pi\rho^2\omega_k} \int_0^\infty dk' \int_0^{2\pi} d\theta' \frac{|\mathbf{k} \times \mathbf{k}'|^4}{|\mathbf{k} - \mathbf{k}'|^4} \frac{\mathcal{V}(k')}{\omega_{k'}^2}$$

$$= \frac{3Ek^4}{8\rho^2\omega_k^3} \left[\int_0^k dk' \mathcal{V}(k') + k^4 \int_k^\infty dk' k'^{-4} \mathcal{V}(k') \right], \quad (\text{C1})$$

where the integral over θ' is performed using the following equation for a constant c :

$$\int_0^{2\pi} d\theta \frac{\sin^4 \theta}{(1 + c^2 + 2c \cos \theta)^2} = \begin{cases} \frac{3\pi}{4} & \text{for } 0 \leq c \leq 1 \\ \frac{3\pi}{4c^4} & \text{for } c > 1 \end{cases}.$$

The self-similarity of the linear potential energy is assumed as $\mathcal{V}(k) = Ck^\alpha$. Then, the two integrals in Eq. (C1) converge when $-1 < \alpha < 3$, and Eq. (C1) is reduced to

$$\omega_k^s = \frac{3}{2(\alpha+1)(3-\alpha)} \sqrt{\frac{12^3(1-\sigma^2)^3}{\rho E h^6}} \frac{\mathcal{V}(k)}{k}. \quad (\text{C2})$$

Finally, Eqs. (9), (12), and (C2) yield Eqs. (13) and (14) for $\alpha = 1$.

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