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Equality Connecting Energy Dissipation with a Violation of the Fluctuation-Response Relation

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In systems driven away from equilibrium, the velocity correlation function and the linear-response function to a small perturbation force do not satisfy the fluctuation-response relation (FRR) due to the lack of detailed balance in contrast to equilibrium systems. In this Letter, an equality between an extent of the FRR violation and the rate of energy dissipation is proved for Langevin systems under nonequilibrium conditions. This equality enables us to calculate the rate of energy dissipation by quantifying the extent of the FRR violation, which can be measured experimentally.

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Macromolecules and colloidal particles of the order of nanometers to submicrometers suspended in an aqueous solution provide an ideal ground to study the foundation of nonequilibrium statistical mechanics. Recent advances in experimental techniques for the observation and manipulation of such small systems have generated the possibility of directly investigating the nonequilibrium nature of fluctuations in the system. In particular, such techniques have been designed to verify several universal relations such as the fluctuation theorem [1–3], the Jarzynski equality [4,5], and the Hatano-Sasa identity [6,7].

Through the investigation of nonequilibrium systems, it has been recognized that the quantification of the violation of the fluctuation-response relation (FRR) [8] provides new information for systems driven far from equilibrium [9– 11]. On the other hand, the rate of energy dissipation is the most fundamental quantity that characterizes nonequilibrium steady states. Thus, it is naturally expected that the FRR violation is related to the amount of energy dissipation. Toward this end, in this Letter, we present an equality between the rate of energy dissipation and an extent of the FRR violation for a class of nonequilibrium stochastic systems.

In this Letter, for simplicity, we mainly study a system described by the Langevin equation

$$\gamma \dot{x}(t) = F(x(t), t) + \xi(t) + \varepsilon f^{p}(t), \qquad (1)$$

where γ is a friction coefficient and $\xi(t)$ is zero-mean white Gaussian noise with variance $2\gamma T$. In particular, we consider two models of the force F(x, t). In the first model, which is termed Model A, $F(x, t) \equiv$ $-\delta_{\sigma(t)0}\partial_x U_0(x) - \delta_{\sigma(t)1}\partial_x U_1(x)$, where $U_i(x)$ for i = 0, 1are periodic potentials with period l, and $\sigma(t)$ is a Poisson process on {0, 1} with a constant transition rate α for the transitions from both 0 to 1 and 1 to 0. This model was originally studied as a model of motor proteins, and is termed a *flashing ratchet model* [12,13]. In the second model, termed Model B, we address a time-independent force $F(x, t) = F(x) \equiv f - \partial_x U(x)$, where f is a constant driving force and U(x) is a periodic potential with period *l*. See Ref. [14] for the physical significance of this model. The last term in Eq. (1) with a sufficiently small ε represents a "probe" force used to investigate the linearresponse property of the system. The initial condition is given at $t = t_{init}$ and we consider the limit $t_{init} \rightarrow -\infty$ below.

First, we define several measurable quantities for this system. Assuming v_s to be the steady-state velocity with $\varepsilon = 0$, the response of the velocity to the probe force is characterized as

$$\langle \dot{x}(t) \rangle_{\varepsilon} = v_{\rm s} + \varepsilon \int_{-\infty}^{t} R(t-s) f^{\rm p}(s) ds + O(\varepsilon^2),$$
 (2)

where $\langle \cdot \cdot \cdot \rangle_{\varepsilon}$ denotes the ensemble average in the presence of the probe force with ε . The response function R(t) has the causality property R(t) = 0 for t < 0. Another important quantity is the time-correlation function of velocity fluctuations in the absence of perturbation, which is defined as

$$C(t) \equiv \langle [\dot{x}(t) - \boldsymbol{v}_{s}]] \dot{x}(0) - \boldsymbol{v}_{s}] \rangle_{0}.$$
(3)

Next, we quantify the energy dissipation in the Langevin description. According to the definition in Ref. [15], the rate of energy dissipation J(t) at time t for each trajectory is expressed as

$$J(t)dt \equiv [\gamma \dot{x}(t) - \xi(t)] \circ dx(t), \tag{4}$$

where • denotes the Stratonovich multiplication [16]. This definition of the rate of energy dissipation complies with both the conservation of energy and the second law of thermodynamics [17].

Based on the above, we present a *theorem* that constitutes the main claim of this Letter:

$$\langle J \rangle_0 = \gamma \left\{ v_s^2 + \int_{-\infty}^{\infty} \left[\tilde{C}(\omega) - 2T\tilde{R}'(\omega) \right] \frac{d\omega}{2\pi} \right\}.$$
 (5)

We express the Fourier transform of an arbitrary function A(t) as $\tilde{A}(\omega) \equiv \int_{-\infty}^{\infty} A(t) \exp(i\omega t) dt$. The prime denotes



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the real part of the function. It is widely acknowledged that in equilibrium, i.e., for cases wherein $\alpha = 0$ for Model A and f = 0 for Model B, the correlation function C(t) is connected to the response function R(t) as C(t) = TR(t)for t > 0, which is the FRR in model (1) [8,14]. Thus, from C(t) = C(-t) and R(t) = 0 for t < 0, we obtain the Fourier transform of this relation as

$$\tilde{C}(\omega) = 2T\tilde{R}'(\omega). \tag{6}$$

Here, we stress that Eq. (6) does not hold in nonequilibrium steady states; thus, it is observed that the right-hand side of Eq. (5) represents the extent of the FRR violation.

It is noteworthy that the equality (5) holds for variety of Langevin systems regardless of the magnitude of external driving as well as the manner in which the system is driven away from equilibrium. We will provide a detailed explanation of this equality in another paper. In the present Letter, we demonstrate the validity of Eq. (5) by the numerical verification for Model A and provide a mathematical proof for Model B.

First, we investigate Model A. Statistical quantities for Model A can be calculated by analyzing the following Fokker-Planck equation that corresponds to the Langevin equation (1) [12,13]:

$$\frac{\partial}{\partial t} \begin{pmatrix} P_0(x,t) \\ P_1(x,t) \end{pmatrix} = \begin{pmatrix} \mathcal{L}_0 - \alpha & \alpha \\ \alpha & \mathcal{L}_1 - \alpha \end{pmatrix} \begin{pmatrix} P_0(x,t) \\ P_1(x,t) \end{pmatrix}, \quad (7)$$

where $\mathcal{L}_i \equiv -\partial_x [F_i(x) + \varepsilon f^p(t) - T\partial_x]/\gamma$ and $F_i(x) \equiv -\partial_x U_i(x)$ for i = 0, 1. In particular, the potentials were selected as $U_0(x) = D \cos(2\pi x/l)$ and $U_1(x) = \text{const.}$, and all the quantities were converted into dimensionless forms by normalizing γ , l, and T to unity. The functions, $\tilde{C}(\omega)$ and $\tilde{R}(\omega)$, were calculated accurately by the matrix-continued-fraction method [14]. In Fig. 1, $\tilde{C}(\omega)$ and $2T\tilde{R}'(\omega)$ are represented as functions of ω when D = 5 and $\alpha = 10$. It exemplifies the violation of the FRR in a nonequilibrium

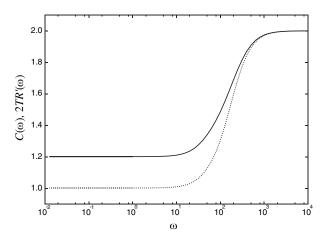


FIG. 1. $\tilde{C}(\omega)$ and $2T\tilde{R}'(\omega)$ as functions of ω when D = 5 and $\alpha = 10$. All the quantities are dimensionless under the normalization as $\gamma = l = T = 1$. The solid and dotted lines represent $\tilde{C}(\omega)$ and $2T\tilde{R}'(\omega)$, respectively.

steady state. In this case, it should be noted that $v_s = 0$ because the selected potentials are symmetric with respect to the reflection of the variable *x*. We then calculated the right-hand side of Eq. (5), which amounted to 16.891, by integrating the difference between $\tilde{C}(\omega)$ and $2T\tilde{R}'(\omega)$ over the entire frequency domain. On the other hand, the rate of energy input for Model A has been calculated as [15,18]

$$\langle J_{\rm in} \rangle_0 = \alpha \int_0^\ell [U_1(x) - U_0(x)] [P_0^{\rm st}(x) - P_1^{\rm st}(x)] dx,$$
 (8)

where $(P_0^{\text{st}}(x), P_1^{\text{st}}(x))$ is the stationary solution of Eq. (7). As a result of the energy balance, the rate of energy input $\langle J_{\text{in}} \rangle_0$ should coincide with the rate of energy dissipation $\langle J \rangle_0$ irrespective of the validity of Eq. (5). Subsequently, if $\langle J_{\text{in}} \rangle_0$ coincides with the right-hand side of Eq. (5), it indicates the validity of Eq. (5). Using the above parameters, we obtained $\langle J_{\text{in}} \rangle_0 = 16.891$. Thus, in this case, Eq. (5) has been confirmed.

In order to demonstrate that this result is not accidental, both the quantities defined in Eqs. (5) and (8) are represented as functions of α for two values of *D* in Fig. 2. This figure clearly shows a good agreement between $\langle J \rangle_0$ and $\langle J_{in} \rangle_0$, independent of the values of the model parameters. In this manner, Eq. (5) has been numerically verified for Model A.

Next, we provide a mathematical proof of the theorem. For simplicity, we restrict our argument to Model B in this proof. In the case of Model B, it has recently been demonstrated that the response Eq. (2) can be derived from the Langevin equation (1) by transforming the force F(x(t)) as [19]

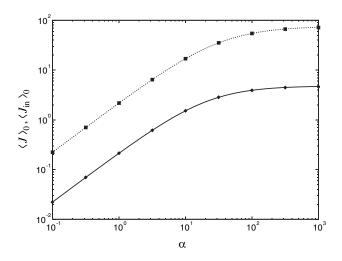


FIG. 2. Energy fluxes $\langle J \rangle_0$ and $\langle J_{\rm in} \rangle_0$ calculated on the basis of Eqs. (5) and (8), respectively, as functions of the transition rate α . The solid line and the closed circles represent $\langle J \rangle_0$ and $\langle J_{\rm in} \rangle_0$, respectively, when D = 1. The dotted line and the closed squares represent $\langle J \rangle_0$ and $\langle J_{\rm in} \rangle_0$, respectively, when D = 5. As in Fig. 1, all the quantities are dimensionless.



$$F(x(t)) = \gamma v_{s} + \int_{-\infty}^{t} K_{0}(t-s) \cdot [\xi(s) + \varepsilon f^{p}(s)] ds$$
$$+ \int_{-\infty}^{t} K_{\perp}(t-s;x(s)) \cdot [\xi(s) + \varepsilon f^{p}(s)] ds, \quad (9)$$

where \cdot denotes the Itô multiplication [16]. We define $K_0(t) = 0$ and $K_{\perp}(t; x(s)) = 0$ for t < 0. This decomposition of force is determined by defining $K_{\perp}(t; x(s))$ such that it satisfies the property

$$\langle K_{\perp}(t;x(s))\rangle_0 = 0 \tag{10}$$

for arbitrary t > 0 and s. By taking the average of Eq. (1) after substitution of Eq. (9), we obtain Eq. (2) by the use of Eq. (10). Then, it is established that

$$\gamma \tilde{R}(\omega) = \tilde{K}_0(\omega) + 1. \tag{11}$$

See Ref. [19] for further details.

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We now set $\varepsilon = 0$. Hereafter, in order to distinctly treat the coexistence of two types of multiplication \circ and \cdot , we discretize time as $t_n \equiv n\Delta t$. Furthermore, we introduce a Wiener process [16], W(t), in place of the noise, $\xi(t)$. Accordingly, the following notations are employed: $x_n \equiv x(t_n)$, $\Delta x_n \equiv x_{n+1} - x_n$, and $\Delta W_n \equiv W(t_{n+1}) - W(t_n)$. By the definition of the Wiener process, the relations, $\langle \Delta W_n \rangle_0 = 0$ and $\langle \Delta W_n \Delta W_m \rangle_0 = \delta_{nm} \Delta t$, hold.

Combining Eqs. (1) and (4) and taking into account the definition of the multiplication \circ , we obtain

$$J(t_n)\Delta t = \bar{F}_n \Delta x_n + O(\Delta t^{3/2}), \qquad (12)$$

where $\bar{F}_n \equiv [F(x_n) + F(x_{n+1})]/2$. On the other hand, integration of Eq. (1) from t_n to t_{n+1} gives

$$\gamma \Delta x_n = \bar{F}_n \Delta t + \sqrt{2\gamma T} \Delta W_n + O(\Delta t^2).$$
(13)

A straightforward calculation then gives

$$\langle J(t_n) \rangle_0 = \gamma v_s^2 + \gamma \left\langle \left(\frac{\Delta x_n}{\Delta t} - v_s \right)^2 \right\rangle_0 - \frac{2T}{\Delta t} - \sqrt{\frac{2T}{\gamma}} \frac{\langle \bar{F}_n \Delta W_n \rangle_0}{\Delta t} + O(\Delta t^{1/2}).$$
(14)

In the limit of $\Delta t \rightarrow 0$, the second and third terms on the right-hand side of Eq. (14) can be transformed as

$$\lim_{\Delta t \to 0} \left[\gamma \left\langle \left(\frac{\Delta x_n}{\Delta t} - \upsilon_s \right)^2 \right\rangle_0 - \frac{2T}{\Delta t} \right] = \int_{-\infty}^{\infty} [\gamma \tilde{C}(\omega) - 2T] \frac{d\omega}{2\pi}.$$
(15)

Next, since the discrete expression of Eq. (9) becomes

$$F(x_n) = \gamma v_s + \sqrt{2\gamma T} \sum_{k=1}^{\infty} K_0(t_k) \Delta W_{n-k} + \sqrt{2\gamma T} \sum_{k=1}^{\infty} K_{\perp}(t_k, x_{n-k}) \Delta W_{n-k}, \quad (16)$$

it follows from Eq. (10) that

$$\langle \bar{F}_n \Delta W_n \rangle_0 = \sqrt{\frac{\gamma T}{2}} K_0(\Delta t) \Delta t,$$
 (17)

In addition, using Eq. (11), Fourier's integration theorem, $\lim_{\Delta t \to 0} [K_0(-\Delta t) + K_0(\Delta t)]/2 = \int_{-\infty}^{\infty} \tilde{K}'_0(\omega) d\omega/2\pi,$ leads to

$$\lim_{\Delta t \to 0^+} K_0(\Delta t) = 2 \int_{-\infty}^{\infty} [\gamma \tilde{R}'(\omega) - 1] \frac{d\omega}{2\pi}.$$
 (18)

Substituting Eqs. (15), (17), and (18), into Eq. (14), we obtain the theorem.

We now present several discussions on the generality of the equality proved here. First, although this proof has been restricted to Model B, we can prove Eq. (5) in almost the same manner for Model A, by determining the decomposition of the force as expressed in Eq. (9). Furthermore, our result represented in Eq. (5) can be generalized for a larger class of Brownian ratchet models [20], including models with an inertia effect, a time-periodic potential, and spatially inhomogeneous temperature profile. We will provide a detailed description of these derivations in another paper.

Second, we can also analyze systems with many degrees of freedom, such as colloidal dispersions, in a parallel way. As a simple example, we consider a system where *N* spherical colloidal particles in a three dimensional aqueous solution are driven by an constant external force fe_x . (See Ref. [21] for a related experimental system.) Let us denote the coordinates of the particles by $\Gamma = \{x_i\}$ (*i* = 1,..., 3*N*), where $\mathbf{r}_{\mu} \equiv (x_{3\mu-2}, x_{3\mu-1}, x_{3\mu})$ represents the position of the μ th particle ($\mu = 1, ..., N$). Then, a widely used model describing the motion of the particles is provided as [22]

$$\gamma \dot{x}_i(t) = F_i(\Gamma(t)) + \xi_i(t) + \varepsilon f_i^{\rm p}(t), \qquad (19)$$

where $F_i(\Gamma) = \sum_{\mu=1}^N f \delta_{i,3\mu-2} - \partial_{x_i} \sum_{\mu=1}^N U(\mathbf{r}_{\mu}) - \partial_{x_i} \sum_{\mu,\nu=1}^N U_{\mu\nu}^{int} (|\mathbf{r}_{\mu} - \mathbf{r}_{\nu}|)/2$ represents single-body forces and two-body interactions, and the noise satisfies that $\langle \xi_i(t)\xi_j(s)\rangle_0 = 2\gamma T \delta_{ij}\delta(t-s)$. For this model, the energy dissipated into the solvent is expressed as $J(t)dt \equiv \sum_{i=1}^{3N} [\gamma \dot{x}_i(t) - \xi_i(t)] \circ dx_i(t)$ [23]. We can then prove

$$\langle J \rangle_0 = \sum_{i=1}^{3N} \gamma \left\{ \langle \dot{x}_i \rangle_0^2 + \int_{-\infty}^{\infty} [\tilde{C}_{ii}(\omega) - 2T\tilde{R}'_{ii}(\omega)] \frac{d\omega}{2\pi} \right\}, \quad (20)$$

where $C_{ij}(t) \equiv \langle [\dot{x}_i(t) - \langle \dot{x}_i \rangle_0] [\dot{x}_j(0) - \langle \dot{x}_j \rangle_0] \rangle_0$ are the cross correlations of velocity fluctuations, and $R_{ij}(t)$ are the cross response functions defined as

$$\langle \dot{x}_i(t) \rangle_{\varepsilon} = \langle \dot{x}_i \rangle_0 + \varepsilon \sum_{j=1}^{3N} \int_{-\infty}^t R_{ij}(t-s) f_j^{\mathrm{p}}(s) ds + O(\varepsilon^2).$$
(21)

The derivation of Eq. (20) will be presented in another paper, although it is a straightforward extension of the above argument for Model B.

In addition, we present another generalization of Eq. (5). For Model B, by a similar argument with the above proof.

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For Model B, by a similar argument with the above proof, we can derive an expression for the symmetrized time correlation between the velocity and force $I(t) \equiv \langle [\dot{x}(t) \circ F(x(0)) + \dot{x}(0) \circ F(x(t))] \rangle_0 / 2$ as

$$I(t) = \gamma \left\{ v_{\rm s}^2 + \int_{-\infty}^{\infty} [\tilde{C}(\omega) - 2TR'(\omega)] e^{i\omega t} \frac{d\omega}{2\pi} \right\}.$$
 (22)

In the case of equilibrium, this equality leads to the FRR [Eq. (6)] since we can derive I(t) = 0 and $v_s = 0$ from the detailed balance condition. Thus, in order to further investigate the physical meanings of the FRR violation, it might be important to study the function I(t).

Before we conclude, let us discuss the physical significance of the theorem. From an experimental point of view, it has been difficult to estimate the amount of the rate of energy dissipation in systems under investigation. The virtue of the expression in Eq. (5) is that it enables us to determine the energy dissipation from experimentally accessible quantities alone, without knowing every detail of the system such as the profile of the force, F(x, t). It is expected that the present result is also useful in the study of biological molecular motors.

In conclusion, we have presented an equality between the rate of energy dissipation and the extent of the FRR violation for a class of Langevin systems under nonequilibrium conditions. To the best of our knowledge, no previous study has addressed this equality, with the exception of a similar, but not precise, expression that was conjectured by one of the authors [24]. However, we should state that an inequality between the rate of entropy production and the FRR violation was demonstrated in Ref. [25].

In the present Letter, we restricted our argument to Langevin dynamics. To extend our result for more general nonequilibrium systems that are not described by Langevin dynamics is an important open problem. Related to this problem, one might be able to prove the present result on the basis of microscopic dynamics, e.g., the Liouville equation, at least with focusing on the vicinity of equilibrium. Experimental examinations on various nonequilibrium systems are also of great importance.

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