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Supplemental Material: Coupled Nosé-Hoover Equations of Motions

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Abstract

Details of mathematical aspects and trajectories of control variables are described for a text entitled "Coupled Nosé-Hoover Equations of Motions." Note: An equation such as "Eq. (1)" refers to the equation that appears in the text, while an equation such as "Eq. (S1)" refers to the equation that appears in this document.

I. EVALUATION OF ϵ

The value of $\epsilon_i(\beta_0) \in \mathbb{R}$ (i = 1, ..., m), defined by Eq. (45), can be evaluated, under suitable mathematical conditions, such that

$$\epsilon_i(\beta_0) = -\int_{\mathbb{R}^{2n}} D_{\beta_i} \rho_{\mathrm{E}}(E(x,p),\beta_0) dx dp \left/ \int_{\mathbb{R}^{2n}} \rho_{\mathrm{E}}(E(x,p),\beta_0) dx dp \right|$$
(S1a)

$$= -\int_{\mathbb{R}^{2n}} \rho_{\mathrm{E}}(E(x,p),\beta_0) D_{\beta_i} \ln \rho_{\mathrm{E}}(E(x,p),\beta_0) dx dp \left/ \int_{\mathbb{R}^{2n}} \rho_{\mathrm{E}}(E(x,p),\beta_0) dx dp \right|$$
(S1b)

$$= \left\langle Y_i^{\beta_0} \right\rangle_{\rho_{\rm E};\beta_0},\tag{S1c}$$

where $Y_i^{\beta_0}(x,p) \equiv -D_{\beta_i} \ln \rho_{\rm E}(E(x,p),\beta_0)$ for $(x,p) \in \mathbb{R}^{2n}$. Namely, $\epsilon_i(\beta_0)$ is the average of the function $Y_i^{\beta_0}$ in the distribution whose density, $\rho_0(x,p,\ldots)$, is proportional to $\rho_{\rm E}(E(x,p),\beta_0)$. Note that β_0 is not a dynamical variable here, but just an arbitrarily fixed parameter value. By using e.g., the density dynamics [Phys. Rev. E **65**, 026105 (2002); *ibid.* **71**, 046708 (2005)], such a density ρ_0 can be created, so that $\epsilon_i(\beta_0)$ is obtained as the long-time average, generated by the density dynamics, of $Y_i^{\beta_0}$ in principle. Note that if m = 1 and $\rho_{\rm E}$ is the BG density, then $\epsilon(\beta_0)$ is the energy average under the BG ensemble at temperature $1/k_{\rm B}\beta_0$.

II. THE SOLUTION OF THE INVERSE PROBLEM

Consider s approximate equalities,

$$\langle h_{\nu}(\beta) \rangle \simeq \gamma_{\nu} \in \mathbb{R}, \quad \nu = 1, \dots, s,$$
 (S2)

for given values, $(\gamma_1 \ldots, \gamma_s) = \gamma$, in general [s = 1 is used in Eq. (65)]. Here the LHS is defined using a given function f that contains $l \geq s$ parameters $(\alpha_1 \ldots, \alpha_l) = \alpha \in \mathbb{R}^l$; viz., $f = f(\cdot; \alpha) : \sigma(\mathbb{R}^m) \to \mathbb{R}, \beta \mapsto f(\beta; \alpha)$. We solve Eq. (S2) with respect to α for a given γ . To do this, we assume three conditions: (i) for any given $\beta_0 \in \sigma(\mathbb{R}^m)$, there exists $U_0 \equiv \mathcal{E}(\beta_0) \in \mathbb{R}^M \ (M \in \mathbb{N})$ such that

$$\epsilon(\beta_0; U_0) \simeq 0, \tag{S3}$$

where $\epsilon(\beta_0; U_0)$ is defined in Eq. (47); (ii) we can represent such that $\beta_0 = \lambda(\alpha)$, where the mapping $\lambda : \mathbb{R}^l \to \sigma(\mathbb{R}^m)$ is specifically computed; (iii) for a map $\Phi = (\Phi_1 \dots, \Phi_s)$ defined by $\Phi_{\nu}(\alpha) \equiv \langle h_{\nu} \rangle_{f(\cdot;\alpha)}$ ($\nu = 1, \dots, s$), there exists $\alpha \in \mathbb{R}^l$ such that

$$\Phi(\alpha) = \gamma \tag{S4}$$

for a given $\gamma \in \mathbb{R}^s$.

Now, take a solution (which is not necessarily unique) of Eq. (S4), $\alpha_{\gamma} \in \mathbb{R}^{l}$ [assumption (iii)]. Then we have $\beta_{0} = \beta_{0}^{\gamma} \equiv \lambda(\alpha_{\gamma})$ [assumption (ii)], and set $U_{0} = U_{0}^{\gamma} \equiv \mathcal{E}(\beta_{0}^{\gamma})$ that yields $\epsilon(\beta_{0}^{\gamma}; U_{0}^{\gamma}) \simeq 0$ [assumption (i)]. Remember that we have $\alpha_{\gamma}, \beta_{0}^{\gamma}$, and U_{0}^{γ} for the given γ in advance. We thereby see that α_{γ} also becomes a solution of Eq. (S2), because

$$\frac{\int_{\sigma(\mathbb{R}^m)} h_{\nu}(\beta) f(\beta; \alpha) Z(\beta; U_0) d\beta}{\int_{\sigma(\mathbb{R}^m)} f(\beta; \alpha) Z(\beta; U_0) d\beta} \bigg|_{\substack{\alpha = \alpha_{\gamma}, \\ U_0 = U_0^{\gamma}}} \equiv \langle h_{\nu}(\beta) \rangle \big|_{\substack{\alpha = \alpha_{\gamma}, \\ U_0 = U_0^{\gamma}}}$$
(S5a)

$$\simeq \frac{\int_{\sigma(\mathbb{R}^m)} h_{\nu}(\beta) f(\beta; \alpha) \mathrm{e}^{-(\epsilon(\beta_0; U_0)|\beta)} d\beta}{\int_{\sigma(\mathbb{R}^m)} f(\beta; \alpha) \mathrm{e}^{-(\epsilon(\beta_0; U_0)|\beta)} d\beta} \bigg|_{\substack{\alpha = \alpha_{\gamma}, \beta_0 = \beta_0^{\gamma} \\ U_0 = U_0^{\gamma} = \mathcal{E}(\beta_0^{\gamma})}} \equiv \langle \langle h_{\nu}(\beta) \rangle \rangle \bigg|_{\substack{\alpha = \alpha_{\gamma}, \beta_0 = \beta_0^{\gamma} \\ U_0 = U_0^{\gamma}}}$$
(S5b)

$$\simeq \left. \frac{\int_{\sigma(\mathbb{R}^m)} h_{\nu}(\beta) f(\beta; \alpha) d\beta}{\int_{\sigma(\mathbb{R}^m)} f(\beta; \alpha) d\beta} \right|_{\alpha = \alpha_{\gamma}} \equiv \langle h_{\nu} \rangle_{f(\cdot; \alpha_{\gamma})}$$
(S5c)

$$=\Phi_{\nu}(\alpha_{\gamma})=\gamma_{\nu} \tag{S5d}$$

holds for any ν .

Note that, while assumptions (ii) and (iii) may be loosened for small l, m, and s, assumption (i) should be critical to solve the problem. To specifically see this, suppose an ideal situation where the integration, Eq. (46b), is explicitly performed so that $\langle \langle h(\beta) \rangle \rangle$ is given by a known function, Ψ , of both $\epsilon(\beta_0)$ and α ; viz., $\langle \langle h(\beta) \rangle \rangle = \Psi(\epsilon(\lambda(\alpha)), \alpha)$. Nonetheless, we cannot easily solve $\Psi(\epsilon(\lambda(\alpha)), \alpha) = \gamma$ with respect to α , since the function ϵ is unknown [much effort is required to know the entire functional form; cf. Eq. (S1)], although the functions Ψ and λ are known (by the assumptions).

III. THE EFFECT OF PARAMETER c

A special case.—We investigate the effect of parameter c introduced in Eq. (55). First we consider the function setting in Sec. II I in the text, where $\rho_{\rm E}$ is given by Eq. (53), or its generalization, Eq. (59), and $f \equiv f_{\rm G}$ is given by Eq. (56). Then the ODE is given by Eq. (5) with T(x, p, Q) represented by Eq. (54) and $\tilde{U}_{\epsilon}(Q)$ represented by Eq. (58) with $A_{\epsilon} = c(\epsilon + \alpha_2 + U_0)$. Let $\varphi : t \mapsto (x(t), p(t), \zeta(t), Q(t), \mathcal{P}(t), \eta(t))$ be any solution of ODE (5), $\dot{\omega} = X'(\omega)$. Suppose that we change the parameter value from c > 0 to c' > 0. We then have a number γ that satisfies $c' = e^{-\gamma/l}c =: c_{\gamma}$. Define a phase-space map $\nu_{\gamma} : \Omega \to \Omega, (x, p, \zeta, Q, \mathcal{P}, \eta) \mapsto (x, p, \zeta, Q + \gamma, \mathcal{P}, \eta)$, viz., the translation of Q. Then $\varphi_{\gamma} := \nu_{\gamma} \circ \varphi : t \mapsto (x(t), p(t), \zeta(t), Q(t) + \gamma, \mathcal{P}(t), \eta(t))$ is a solution of an ODE, $\dot{\omega} = X'_{\gamma}(\omega)$, where X'_{γ} is X' using c_{γ} instead of c. This is shown as follows: For any $\omega, X'_{\gamma}(\nu_{\gamma}(\omega)) = X'_{\gamma}(x, p, \zeta, Q + \gamma, \mathcal{P}, \eta)$ is given by Eq. (10) [RHS of Eq. (5)] with $T_{\gamma}(x, p, Q + \gamma)$ and $\nabla \tilde{U}^{\gamma}_{E(x,p)}(Q + \gamma)$, instead of T(x, p, Q) and $\nabla \tilde{U}_{E(x,p)}(Q)$, respectively. Here the index γ in T_{γ} and $\nabla \tilde{U}^{\gamma}_{E(x,p)}$ indicate the use of c_{γ} instead of c. We find that $T_{\gamma}(x, p, Q + \gamma) = 1/(k_{\rm B}c_{\gamma})e^{-(Q+\gamma)/l} = T(x, p, Q)$ and $\nabla \tilde{U}^{\gamma}_{E(x,p)}(Q + \gamma) = (c_{\gamma}/l)(E(x, p) + \alpha_2 + U_0)e^{(Q+\gamma)/l} - (\alpha_1 + 1)/l = \nabla \tilde{U}_{E(x,p)}(Q)$. Thus

$$X'_{\gamma}(\nu_{\gamma}(\omega)) = X'(\omega) \text{ for all } \omega.$$
(S6)

Therefore we have $D\varphi_{\gamma}(t) = D\nu_{\gamma}(\varphi(t)) \circ D\varphi(t) = D\varphi(t) = X'(\varphi(t)) = X'_{\gamma}(\nu_{\gamma}(\varphi(t))) = X'_{\gamma}(\varphi_{\gamma}(t)))$ for all t, so the proof is completed.

Namely, the translation of Q in any solution of the original ODE just yields a solution of the parameter-changed $(c \to c_{\gamma})$ ODE. Conversely, for any solution, φ_{γ} , of $\dot{\omega} = X'_{\gamma}(\omega)$, we similarly see that $\nu_{-\gamma} \circ \varphi_{\gamma}$ becomes a solution of $\dot{\omega} = X'(\omega)$. Thus, we have a mapping between the solution spaces, $S := \{$ solution of $\dot{\omega} = X'(\omega) \} \to S_{\gamma} = \{$ solution of $\dot{\omega} = X'_{\gamma}(\omega) \},$ $\varphi \stackrel{d}{\mapsto} \nu_{\gamma} \circ \varphi$, and find it is bijective, due to $\nu_{\gamma} \circ \nu_{-\gamma} = id_{\Omega} (\forall \gamma)$. This means that we completely know the solutions of the parameter-changed ODE from the solutions of the original ODE, just by the translation.

A general case.—The above results also hold in a more general function setting. Let the map $\sigma : \mathbb{R}^m \to \mathbb{R}^m$ be parameterized by $c \in A$ (a certain set), denoted as σ_c . We only assume two conditions. First, σ possesses the following relationship between a parameter change and a variable translation: for $c_1, c_2 \in A$, there exists $\gamma \in \mathbb{R}^m$ such that $\sigma_{c_2}(Q) = \sigma_{c_1}(Q - \gamma)$ for all Q (this σ_c is a generalization of the map $Q \mapsto ce^{Q/l}$ with m = 1). Second, the vector field, X_c , is smooth and takes the form: $X_c(\omega) = Y(x, p, \zeta, \mathcal{P}, \eta, \sigma_c(Q), D\sigma_c(Q), D^2\sigma_c(Q))$; viz., the dependence on Q is only through σ_c and its first and second differentials, $D\sigma_c$ and $D^2\sigma_c$ [Eq. (5) satisfies this condition, since T(x, p, Q) given by Eq. (6) involves only $\sigma(Q)$ and since $\tilde{U}_{E(x,p)}$ given by Eq. (8) involves only $\sigma(Q)$ and $D\sigma(Q)$ [see also Eq. (3)] so that $\nabla \tilde{U}_{E(x,p)}(Q)$ involves only $D^k \sigma_c(Q)$ (k = 0, 1, 2)]. We can show that if φ is a solution of $\dot{\omega} = X_{c_1}(\omega)$ then $\nu_{\gamma} \circ \varphi$ (ν_{γ} is the one given above) becomes a solution of $\dot{\omega} = X_{c_2}(\omega)$: It follows from $D^k \sigma_{c_2}(Q) = D^k \sigma_{c_1}(Q - \gamma)$ for every Q and k that $X_{c_2}(\nu_{\gamma}(\omega)) = X_{c_1}(\omega)$ for all ω , which corresponds to Eq. (S6). Hence the proof can be completed in the same manner as the special case. Similarly, a bijective mapping between the solution space of $\dot{\omega} = X_{c_1}(\omega)$ and that of $\dot{\omega} = X_{c_2}(\omega)$ is obtained.

IV. THE TOTAL ERROR

Figure 1 shows the total error measured by $\Delta(t)$, with respect to time t, for the distribution of x in the 1-dimensional double-well (1DW) potential system. See text for the details.

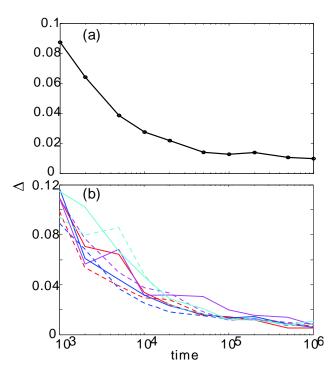


FIG. 1: (a) Difference in the 1D marginal distribution of x between simulation and theory as a function of the simulation time for the 1DW model. (b) The same as (a), but $c_{\rm Z} = 0.1$ (solid, blue) and 10 (dashed, blue), $c_{\rm Y} = 0.1$ (solid, cyan) and 10 (dashed, cyan), $\mathbf{M}_{\rm T} = 0.1$ (solid, red) and 10 (dashed, red), and l = 0.224 (solid, magenta) and 22.4 (dashed, magenta). $c_{\rm Z} = c_{\rm Y} = 1$, $\mathbf{M}_{\rm T} = 1$, and l = 2.24 are used unless otherwise mentioned. All quantities are dimensionless.

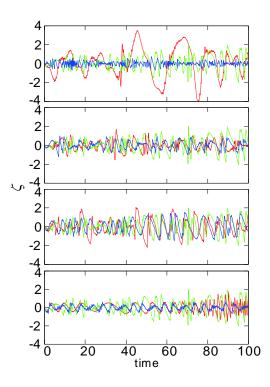


FIG. 2: Time courses of ζ for 1DW. Color schemes are the same as in Fig. 4 of the text.

V. TRAJECTORIES OF ζ , η , AND \mathcal{P} FOR THE 1DW MODEL SYSTEM

The oscillation, as well as the distribution, of ζ depends on c_Z , where a smaller c_Z tends to yield a larger period (Fig. 2). This is reasonable as an analogue that we observe in the single NH equation, because c_Z essentially corresponds to the inverse of Nose's mass $Q_{\rm NH}$ [recall the discussion in Sec. IIB in the text, such that the setting of our $\rho_Z(\zeta)$ leads to the friction term of the NH equation via $c_Z = 1/2k_{\rm B}T_{\rm ex}Q_{\rm NH}$]. In contrast, ζ , which directly controls the physical system, depends minimally on $c_{\rm Y}$ and $\mathbf{M}_{\rm T}$, as these are temperature-system parameters. The effects of l on the dynamics of ζ are not straightforward. ζ is governed by $\dot{\zeta} = T_{\rm P} - T_{\rm D}$ ($nk_{\rm B} = 1$ in 1DW), as seen in the typical examples shown in Fig. 3. When l = 22.4 [Fig. 3(a)], around $T_{\rm D}$, which reduces slowly, $T_{\rm P}$ oscillates fast, so that ζ has two oscillation modes: one is the slow (~3 periods for time 20) mode around zero with relatively large amplitudes and the other is the fast mode, inherited from $T_{\rm P}$, with relatively small amplitudes. While for l = 0.224 [Fig. 3(b)], $T_{\rm D}$ and $T_{\rm P}$ oscillate with larger amplitudes than those in Fig. 3(a). However, since the periods of $T_{\rm D}$ and $T_{\rm P}$ are similar, cancellations occur so that $T_{\rm P} - T_{\rm D}$ maintains small values and the amplitude of ζ is comparable to that in

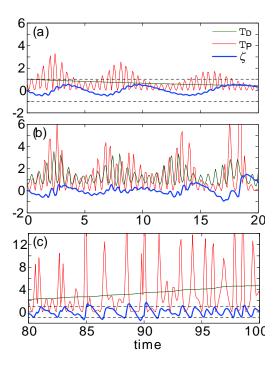


FIG. 3: Time courses of T_D (green) and T_P (red), and ζ (blue) for 1DW. (a) $l = 22.4, 0 \le t \le 20$; (b) $l = 0.224, 0 \le t \le 20$; and (c) $l = 22.4, 80 \le t \le 100$. Dotted lines in (a) and (c) indicate ± 1 .

Fig. 3(a). For l = 22.4 and $80 \le t \le 100$ [Fig. 3(c)], against the slow motion of $T_{\rm D}$ as in $0 \le t \le 20$ [Fig. 3(a)], $T_{\rm P}$ oscillates with very large amplitudes [cf. Fig. 4(b) of the text], which affect the motion of ζ . The faster mode of ζ seen in $0 \le t \le 20$ seems to grow and attain the large amplitudes in $80 \le t \le 100$. This is the reason why ζ for l = 22.4 suddenly has the large fluctuations seen in Fig. 2 (bottom panel).

The trajectories of η show that a larger period and amplitude in the oscillating-like motion of η appear as $c_{\rm Y}$ is decreased (Fig. 4). This dependence of $c_{\rm Y}$ on η is similar to the dependence of $c_{\rm Z}$ on ζ , as expected from the roles of these two parameters [exactly speaking, there is a difference in that the friction term $\tau_{\rm Z}(\zeta)$ described by ζ and $c_{\rm Z}$ in Eq. (5) is multiplied by T(x, p, Q), whereas the friction term $\tau_{\rm Y}(\eta)$ described by η and $c_{\rm Y}$ is not]. The temperature-system variable η does not strongly depend on the physical-system parameter $c_{\rm Z}$. We also see that η —a control variable in the temperature system—does not strongly depend on the temperature-system parameters, $\mathbf{M}_{\rm T}$, which is the mass parameter, and l, which is the potential parameter [see Eq. (58)]. This tendency seems to be reasonable, by considering the analogous tendency in the single NH equation that the control variable in the physical system does not depend strongly on the mass and potential parameters of the physical system.

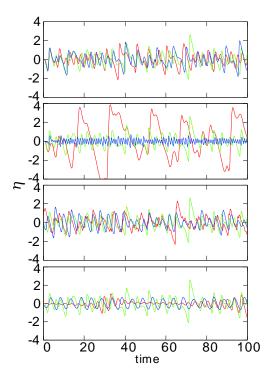


FIG. 4: Time courses of η for 1DW. Color schemes are the same as in Fig. 4 of the text.

The dependence on the parameters of the trajectory of \mathcal{P} (Fig. 5) is similar to that of the dynamical temperature $T_{\rm D}$. This is suggested by the relations, $\mathbf{M}_{\rm T}^{-1}\mathcal{P} = \dot{Q}$ and $T_{\rm D} \propto \exp\left[-Q/l\right]$. Even though the cases of $c_{\rm Y} = 10$ and $\mathbf{M}_{\rm T} = 10$ are seemingly exceptional, the periods of the two oscillating modes existing in $T_{\rm D}$ and in \mathcal{P} are actually similar, whereas the faster modes in \mathcal{P} have large amplitudes and those in $T_{\rm D}$ have small amplitudes. As expected from the relation, $\dot{\eta} = \mathbf{M}_{\rm T}^{-1}\mathcal{P}^2 - 1$, there are also correlations between \mathcal{P}^2 and η . This may not be clear for l = 0.224 in the behavior of \mathcal{P} shown in Fig. 5, but it is actually observed if we view them on a shorter time scale, as shown in Fig. 6.

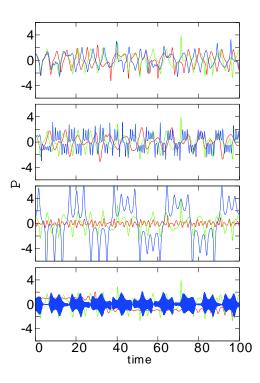


FIG. 5: Time courses of \mathcal{P} for 1DW. Color schemes are the same as in Fig. 4 of the text.

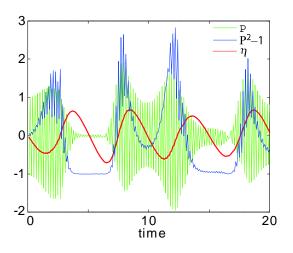


FIG. 6: Time courses of \mathcal{P} (green), $\mathcal{P}^2 - 1$ (blue) and η (red) for 1DW. l = 0.224 and $\mathbf{M}_{\mathrm{T}} = 1$.