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Moments of work in the two-point measurement protocol for a driven open quantum system

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We study the distribution of work induced by the two-point measurement protocol for a driven open quantum system. We first derive a general form for the generating function of work for the total system, bearing in mind that the Hamiltonian does not necessarily commute with its time derivative. Using this result, we then study the first few moments of work by using the master equation of the reduced system, invoking approximations similar to the ones made in the microscopic derivation of the reduced density matrix. Our results show that already in the third moment of work, correction terms appear that involve commutators between the Hamiltonian and its time derivative. To demonstrate the importance of these terms, we consider a sinusoidally, weakly driven and weakly coupled open two-level quantum system, and indeed find that already in the third moment of work, the correction terms are significant. We also compare our results to those obtained with the quantum jump method and find a good agreement.

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I. INTRODUCTION

For microscopic systems driven out of equilibrium, the fluctuation theorems, e.g., Refs. [1–4], provide a powerful tool to analyze the thermodynamic nature of nonequilibrium processes beyond the linear response regime. When the microscopic system can be described in terms of classical mechanics, the fluctuation theorems have been examined for several systems [5–11]. However, when described in terms of quantum mechanics, the situation is more problematic. In quantum systems, it is far from obvious how to treat certain thermodynamic quantities such as work \( W \) that relate to the physical path of the system rather than to the state (wave function).

Work appears in the classical Jarzynski equation (JE) \( \langle e^{-W/k_B T} \rangle = e^{-\Delta F/k_B T} \), where \( \Delta F \) is the free-energy difference between the initial equilibrium and the final states, and the brackets \( \langle \cdot \rangle \) denote averaging over an infinite number of repetitions. Trying to generalize the JE to the quantum regime has caused much debate about how to define \( W \) in a physically meaningful way. Earlier quantum treatments of the JE were based on defining a genuine work operator [12–15]. Yet since work is not a traditional quantum observable [16], the use of a quantum work operator leads to corrections to the JE. It can be recovered by another approach, known as the two-point measurement protocol [17–20], in which the energy of the closed system is measured at the beginning and at the end of the process and there is no dissipated heat. The work of a single trajectory is then defined as the energy difference of the final and initial measurement outcomes. In the case of open systems, assuming that the interaction Hamiltonian is negligible, the energy measurement of the total closed system can be approximated by measuring the energy of the reduced system and the environment separately.

In a recent paper [21], the quantum jump method, also known as the Monte Carlo wave function method (MCWF), was proposed as an efficient way to discuss the problem of determining the statistics of work in driven quantum systems with dissipation. By interpreting a jump between the eigenstates of the Hamiltonian as an emission and absorption of a photon to the heat bath, the total energy exchanged between the system and the heat bath due to the jumps is then interpreted as heat. The work can then be defined as the energy difference between the initial and final states of the system plus the heat released to the heat bath. It should be noted that with this definition, a possible energetic contribution from the interaction between the system and the heat bath was not taken into account in work [22,23].

In this paper, we analyze in detail the first few moments of work by using the master-equation approach for an open quantum system. To characterize the stochastic nature of \( W \) and its distribution, it is natural to consider the moments of work instead of directly trying to calculate exponential averages such as that in the JE, which is a formidable task for open quantum systems in general. The first moment gives the mean work done, the second moment gives the variance, and the third moment gives the skewness of the work distribution for non-Gaussian distributions. As the first step, we derive the two-point measurement protocol generating function without making the implicit assumption in Ref. [24] that the total system Hamiltonian commutes with its time derivative. This result allows us to derive general expressions for the first three moments of work, which we compare with results obtained using the generating function of Ref. [24] (Eqs. (17), (18), (22), and (23) in Ref. [24]). Our results show that only the first two moments of work are identical in the two approaches above, and nontrivial correction terms appear to the third and higher moments when the Hamiltonian does not commute with its time derivative. To study this issue in a specific case, we consider the weakly coupled and weakly driven open two-level
quantum system of Ref. [21], where we invoke approximations similar to those used in the microscopic derivation of the Lindblad equation of the reduced system. The test system describes, for instance, a Cooper box coupled capacitively or a dc superconducting quantum interference device (dc-SQUID) coupled inductively to a calorimeter [25]. When calculating the dynamics of the test system, we neglect the interaction Hamiltonian in the energy measurements. We indeed find that our results for the first three moments are in agreement with the quantum jump results. When comparing the two different generating functions, we find a significant difference in the values of the third moment.

The general results derived here are not restricted to a Cooper box and a dc-SQUID, but can be used for various kinds of superconducting qubits [26] and quantum dot circuits [27–29].

II. GENERATING FUNCTION AND MOMENTS FOR WORK

In the two-measurement protocol for a closed quantum system, the probability to measure energy \(E_t\) at time \(t = 0\) and \(E_r\) at \(t = \tau\) is of the form [30]

\[
P[E_t, E_0] = \text{Tr}\{\hat{P}_E(t)\hat{U}(0, \tau)\hat{P}_E\hat{U}^\dagger(0, \tau)\hat{P}_E\},
\]

where \(\hat{U}(0, \tau) = T_{\text{ev}}\exp[-i\int_0^\tau dt \hat{H}(t)]\) is the unitary time evolution operator, \(T_{\text{ev}}\) describes the chronological time ordering, and the projection operators are given by \(\hat{P}_E = \{E_r\}\{E_t\}\), where \(\{E_t\}\) is the state corresponding to the measurement result \(E_t\) at time \(t\). The corresponding generating function is the Fourier transform of \(P[E_t, E_0]\) [24] (the calculation is also given in Appendix A):

\[
G(u) = \sum_{E_t, E_0} e^{iu(E_t - E_0)} P[E_t, E_0] = \text{Tr}[\hat{U}_{u/2}(0)\hat{P}_0\hat{U}_{-u/2}(0)],
\]

where

\[
\hat{U}_u(\tau, 0) = e^{iu\hat{H}(\tau)}\hat{U}(0, \tau)e^{-iu\hat{H}(0)},
\]

\[
\hat{P}_0 = \sum_{E_0} \hat{P}_E\hat{P}_0\hat{P}_E,
\]

and \(\hat{P}_0\) is the initial density matrix. If the initial density matrix is diagonal in the first measurement basis, then \(\hat{P}_0 = \hat{P}_0\).

The differentiation of the evolution operator \(\hat{U}_u(\tau, 0)\) [Eq. (3)] with respect to \(\tau\) yields the following equation of motion:

\[
\frac{d\hat{U}_u(\tau, 0)}{d\tau} = \left[\frac{-i}{\hbar}\hat{H}(\tau) + \sum_{n=1}^{\infty} \frac{(iu)^n}{n!}\hat{C}_n(\tau)\right]\hat{U}_u(\tau, 0),
\]

where \(\hat{C}_1(\tau) = \hat{\delta}_1\hat{H}(\tau)\), \(\hat{C}_2(\tau) = \{\hat{H}(\tau), \hat{\delta}_1\hat{H}(\tau)\}\), \(\hat{C}_3(\tau) = \{\hat{H}(\tau), [\hat{H}(\tau), \hat{\delta}_1\hat{H}(\tau)]\}\), etc. The generating function can be then written as (see Appendix A)

\[
G(u) = \text{Tr}\left\{T_{\text{ev}}\exp\left[\int_0^\tau dt \sum_{n=1}^{\infty} \frac{(-1)^{n+1}(iu)^n}{n!2^n}\hat{C}_n(H(t))\right]\times T_{\text{ev}}\exp\left[\int_0^\tau dt \sum_{n=1}^{\infty} \frac{(iu)^n}{n!2^n}\hat{C}_n(H(t))\right]\hat{P}_0\right\},
\]

where the superscript \(H\) indicates the Heisenberg picture, i.e., \(\hat{C}_n(H(t)) = \hat{U}(t, 0)^\dagger\hat{C}_n(0)\hat{U}(t, 0)\). The moments of work are then obtained by differentiating \(G(u)\) with respect to \(u\) at \(u = 0\):

\[
\langle W^n \rangle = \left(-i\right)^n \left.\frac{\partial^n G(u)}{\partial u^n}\right|_{u=0}.
\]

With the implicit assumption that \(\{\hat{H}(t), \hat{\delta}_1\hat{H}(t)\} = 0\) (Ref. [24]), \(\hat{C}_n = 0\) for \(n > 1\), and the generating function becomes

\[
G_0(u) = \text{Tr}\left\{T_{\text{ev}}\exp\left[i\int_0^\tau dt \hat{\Phi}(H(t))\right]\times T_{\text{ev}}\exp\left[i\int_0^\tau dt \hat{\Phi}(H(t))\right]\hat{P}_0\right\},
\]

where the power operator \(\hat{\Phi}\) (Ref. [22]) is the time derivative of the total Hamiltonian, i.e., \(\hat{\Phi}(H(t)) = \frac{d}{dt}\hat{H}(t)\hat{U}(t, 0)\).

The generating functions of Eqs. (6) and (8) are equivalent to the first order of \(u\). Thus, both generating functions trivially give the same expression for the first moment of work as

\[
\langle W \rangle = \int_0^\tau dt_1 \langle \hat{\Phi}(H(t_1))\rangle.
\]

Although the generating functions of Eqs. (6) and (8) differ already to second order in \(u\), the expressions for the second moment turn out to be equal as the corrections given by Eq. (6) cancel out,

\[
\langle W^2 \rangle = 2\int_0^\tau dt_1 \int_0^\tau dt_2 \text{Re}\{\langle \hat{\Phi}(H(t_1))\hat{\Phi}(H(t_2))\rangle\},
\]

where we have used the Hermiticity of \(\hat{\Phi}\) to further simplify the expression. The same expressions for the first and second moment are also obtained by using the work operator with and without the commutator of the Hamiltonian at different times [15]. However, for the third moment, the two generating functions give different results as

\[
\langle W^3 \rangle_0 = 3\int_0^\tau dt_1 \int_0^{t_1} dt_2 \int_0^{t_2} dt_3 \text{Re}\{\langle \hat{\Phi}(H(t_1))\hat{\Phi}(H(t_2))\hat{\Phi}(H(t_3))\rangle + \langle \hat{\Phi}(H(t_2))\hat{\Phi}(H(t_1))\hat{\Phi}(H(t_3))\rangle\},
\]

\[
\langle W^3 \rangle = \langle W^3 \rangle_0 + \frac{1}{4}\int_0^\tau dt \langle \hat{C}_1^3(H(t))\rangle + \frac{3}{2}\int_0^\tau dt_1 \int_0^{t_1} dt_2 \text{Re}\{\langle \hat{C}_1^2(H(t_1))\hat{C}_2(H(t_2))\rangle\},
\]

where \(\langle W^3 \rangle_0\) denotes the third moment given by Eq. (8) and \(\langle W^3 \rangle\) denotes the one given by our general expression of Eq. (6). The moments given by Eq. (8) consist of third-order correlation functions of the power operator. In our result here, there are additional correction terms that involve commutators.
between the Hamiltonian and its time derivative, as expected. Such correction terms appear also in the higher moments of work.

III. OPEN QUANTUM SYSTEM

To illustrate the importance of the results we have derived here, let us consider the special case of a weakly driven system, which is also weakly coupled to a heat bath [21]. The total Hamiltonian is taken to be of the form

\[
\hat{H}(t) = \hat{H}_S(t) + \hat{H}_B + \hat{H}_C,
\]

where subscripts \(S, B, \) and \(C\) denote the system, bath, and bath-system interaction (coupling) Hamiltonians, respectively. Both the bath and the system-bath interaction (coupling) Hamiltonians are assumed to be time independent. The system Hamiltonian \(\hat{H}_S(t) = \hat{H}_0 + \hat{V}(t)\) consists of a time-independent part \(\hat{H}_0\) and a time-dependent perturbative part \(\hat{V}(t)\). Therefore, the time derivative of the total Hamiltonian is simply given by \(\dot{\hat{H}}(t) = \partial_t \hat{H}(t) = \partial_t \hat{V}(t)\). In principle, we can calculate the moments of work from Eq. (6). However, already all the correlation functions of the third moment cannot be calculated just using the reduced density matrix \(\hat{\rho}(t)\), as the correlation functions contain the total Hamiltonian that does not depend only on the system degrees of freedom but also on the bath degrees of freedom. To proceed, we consider a specific model, where a two-level system as in Ref. [21] is bilinearly coupled to a heat bath of bosonic modes. The system Hamiltonian has the form

\[
\hat{H}_S(t) = \hat{H}_0 + \hat{V}(t),
\]

\[
\hat{H}_0 = \hbar \omega_0 \hat{a}^\dagger \hat{a},
\]

\[
\hat{V}(t) = \lambda(t) \hat{a}^\dagger \hat{a} + \hat{a},
\]

where \(\hat{a} = |e\rangle\langle g|\) and \(\hat{a}^\dagger = |g\rangle\langle e|\) are the creation and annihilation operators, respectively, in the ground-state (|g\rangle) and excited-state (|e\rangle) basis of the undriven system, \(\hbar \omega_0\) is the energy separation of the two levels, and \(\lambda(t)\) is the time-dependent drive. Further, the interaction and bath Hamiltonians are assumed to be of the form

\[
\hat{H}_C = \sum_k (\hat{a}^\dagger + \hat{a}) \otimes (g_k \hat{b}_k^\dagger + g_k^* \hat{b}_k),
\]

\[
\hat{H}_B = \sum_k \hbar \omega_k \hat{b}_k^\dagger \hat{b}_k,
\]

where \(g_k\) is the coupling strength, and \(\hat{b}_k\) and \(\hat{b}_k^\dagger\) are the bath annihilation and creation operators associated with energy \(\hbar \omega_k\), respectively. For the total Hamiltonian \(\hat{H}(t)\), this implies \([\hat{H}(t), \partial_t \hat{H}(t)] \neq 0\). In the calculations, we approximate the initial density matrix \(\hat{\rho}_0\) with the tensor product of the system and bath density matrices, where both the system and the heat bath start in thermal equilibrium. That is, we neglect the interaction Hamiltonian in the energy measurements. Due to the weak driving and coupling to the heat bath, the evolution of the two-level system can be approximated with the following Lindblad equation by invoking the Born-Markov and secular approximations (see Appendix B):

\[
\frac{d\hat{\rho}}{dt} = -\frac{i}{\hbar} [\hat{H}(t), \hat{\rho}(t)] + \Gamma_1 \left( \rho_{ee}(t)|g\rangle\langle g| - \frac{1}{2} \hat{\rho}(t), |e\rangle\langle e| \right)
\]

\[
+ \Gamma_1^\dagger \left( \rho_{gg}(t)|e\rangle\langle e| - \frac{1}{2} \hat{\rho}(t), |g\rangle\langle g| \right),
\]

where \(\Gamma_1\) and \(\Gamma_1^\dagger = \Gamma_1 e^{\gamma(t)/\hbar}\) are the photon emission and absorption transition rates, respectively, \(\hat{\rho}(t)\) is the density matrix of the reduced system in the Schrödinger picture, and \(\rho_{kl}(t) = \langle k|\hat{\rho}(t)|l\rangle\).

As the secular approximation neglects the fast oscillating coupling terms, the same master equation could have been achieved by starting with the following form of the interaction Hamiltonian:

\[
\hat{H}_C^{RWA} = \sum_k g_k \hat{b}_k^\dagger \hat{b}_k + g_k^* \hat{b}_k^\dagger \otimes \hat{b}_k,
\]

where the rotating wave approximation (RWA) has been invoked. With this form of the interaction Hamiltonian [Eq. (20)], the jumps can be easily interpreted as photon emission and absorption to the bath. The usual quantum jump method [31–34] can then be used to calculate the work distribution by interpreting the jumps as photon exchange while neglecting the energetic contribution due to \(\hat{H}_C\).

The first two moments for the system can be calculated in the usual manner by using the master equation of the reduced density matrix, as the operators in the correlation functions depend only on the system degrees of freedom [35,36]. For the third moment \(\langle W^3 \rangle\), we can simplify the expression by using the fact that the power operator \(\hat{P}(t)\) and the interaction Hamiltonian \(\hat{H}_C\) [Eq. (17)] commute,

\[
\langle W^3 \rangle = \langle W_3^S \rangle + \frac{1}{4} \int_0^\tau dt \langle [\hat{H}_C(t), [\hat{H}_S(t), \hat{P}(t)]] \rangle
\]

\[
+ \frac{3}{4} \int_0^\tau dt_1 \int_0^{t_1} dt_2 \text{Re} \{ [\hat{P}(t_1), [\hat{H}_S(t_2), \hat{H}_S(t_2)]] \}
\]

\[
+ \frac{1}{4} \int_0^\tau dt \langle [\hat{H}_C(t), [\hat{H}_S(t), \hat{P}(t)]] \rangle
\]

\[
= \langle W_3^S \rangle + \langle W_3^{S+B} \rangle,
\]

where \(\langle W_3^S \rangle\) is given in the first two lines of the above equation and consists of the correlation functions that include only system operators. The interesting part is the second term \(\langle W_3^{S+B} \rangle\) that contains also operators that depend on the bath degrees of freedom,

\[
\langle W_3^{S+B} \rangle = \frac{1}{4} \int_0^\tau dt \langle [\hat{H}_C(t), [\hat{H}_S(t), \hat{P}(t)]] \rangle.
\]

We can estimate the term \(\langle W_3^{S+B} \rangle\) by invoking approximations similar to those used in the derivation of the corresponding master equation (see Appendix C), yielding

\[
\langle W_3^{S+B} \rangle \approx \frac{\hbar^2 \omega_0}{2} (\Gamma_1 + \Gamma_1^\dagger) \int_0^\tau dt \langle \hat{\lambda}(t) \text{Im} \{\rho_{ee}(t)\} \rangle.
\]
of work can be calculated numerically by using the master equation for a weak \( \lambda(t) \).

In the case of a simple sinusoidal resonance drive \( \lambda(t) = \lambda_0 \sin(\omega_0 t) \), \( \langle W^3 \rangle_{S+B} \) can be further approximated by changing to the interaction picture and neglecting the fast oscillating terms,

\[
\langle W^3 \rangle_{S+B} \approx \frac{\lambda_0^2 \omega_0^2}{4} (\Gamma_1 + \Gamma_1) \int_0^\tau dt \text{Im}\{\rho_{eg}(t)\}, \quad (24)
\]

where \( \hat{\rho}(t) \) is the density matrix of the reduced system in the interaction picture with respect to \( \hat{H}_0 + \hat{H}_B \) and \( \rho_{eg}(t) = \langle e | \hat{\rho}(t) | g \rangle \).

For the sinusoidal resonance drive, we can simplify the analytical calculations of the correlation functions of the work moments with an additional rotating wave approximation. By neglecting the fast oscillating terms, the power operator simplifies to the form \( \hat{P}(t) \approx \lambda_0 \omega_0 (\hat{a} + \hat{a}^\dagger) / 2 \) in the interaction picture. Using the regression theorem [37], we can then calculate analytical approximations for the moments of work.

The regression theorem results with the additional RWA were found to give an excellent agreement with the numerical master-equation results when the driving period \( \tau \) consists of full or half cycles. When the driving period is not \( n \pi \), where \( n \) is an integer, there can be a small difference between the regression theorem results and the numerical master-equation results. This difference is due to the oscillation caused by the fast oscillation terms of the drive for the latter and is illustrated in Fig. 1 for the second moment \( \langle W^2 \rangle \) with \( \lambda_0 = 0.05\hbar \omega_0 \). As the oscillation is caused by the fast oscillating terms of the drive, the deviation becomes larger when the value of \( \lambda_0 \) is increased.

![Fig. 1. (Color online) The numerical master-equation results for the second moment \( \langle W^2 \rangle \) as a function of time for two different coupling strengths. Inset: The numerical results are compared to the analytical approximation \( \langle W^2 \rangle_{\text{RWA}} \), achieved with the additional RWA. The driving is assumed to be in resonance with \( \hbar \omega_0 \), i.e., \( \omega = \omega_0 \), \( \beta\hbar\omega_0 = 2.0 \), and \( \lambda_0 = 0.05\hbar \omega_0 \). The oscillation in the numerical results is caused by the fast oscillating terms of the drive. These are neglected in the analytical results by invoking the additional RWA. Inset: The oscillation for both coupling strengths is almost identical up to \( \omega_0 \tau = 10\pi \).](image1)

![Fig. 2. (Color online) Comparison of the quantum jump method and master-equation results for the first three moments for different coupling amplitudes. The solid and dashed lines correspond to the analytical results with the additional RWA, the dots correspond to the numerical quantum jump results, and the crosses correspond to the numerical master-equation data. The driving is assumed to be in resonance with \( \hbar \omega_0 \), i.e., \( \omega = \omega_0 \), \( \beta\hbar\omega_0 = 2.0 \), \( \lambda_0 = 0.05\hbar \omega_0 \) and the drive lasts for 10 cycles, i.e., \( \omega \tau = 20\pi \). The numerical results are calculated with \( 10^4 \) time steps. The quantum jump results consist of \( 10^6 \) realizations. The numerical master-equation and quantum jump results give a good agreement within the numerical accuracy: The largest difference in \( \langle W^n \rangle/\langle \hbar \omega_0 \rangle^n \) is less than 0.0032.](image2)

![Fig. 3. (Color online) The standard fluctuation dissipation theorem is valid not only in the linear response regime \( \lambda_0 \rightarrow 0 \) but also in the limit of no coupling \( \Gamma_1 \rightarrow 0 \) with arbitrary drive amplitude within this model. In the case of no coupling, the probability to end up in the excited state when starting from the ground state, denoted as \( p_{eg} = |\langle e | \hat{U}(\tau,0) | g \rangle |^2 \), is exactly the same as the probability to end up in the ground state when starting from the excited state, \( p_{ge} \). Hence, \( \langle W^n \rangle = \langle \hbar \omega_0 \rangle^n p_{eg}(0)p_{ge} + (-\hbar \omega_0 \rangle^n p_{ge}(0)p_{eg} = \).](image3)
are needed to check fluctuation relations such as the JE. To properly evaluate the higher moments of work, which is the case of the third moment. This emphasizes the importance have shown that there is a significant difference already in the total Hamiltonian and its time derivative commute and 10 cycles, i.e., \( \alpha \tau = 20\pi \). As expected, significant deviations start appearing with increased coupling and drive.

\[
(h\omega_0)^n p_{\text{PE}}(0) + (-1)^n p_{\text{PE}}(0),
\]
which immediately gives the FDT when we start from thermal equilibrium.

For small values of \( \lambda_0 \) and \( \Gamma_\tau \), the deviation from the FDT increases almost parabolically when the drive amplitude \( \lambda_0 \) increases and the transition rate \( \Gamma_\tau \) remains constant for small values of \( \lambda_0 \) and \( \Gamma_\tau \). This can be seen by Taylor expanding \( (W^2)_{\text{RWA}}/(W)_{\text{RWA}} \) around \( (\lambda_0, \Gamma_\tau) = (0,0) \),

\[
(W^2)_{\text{RWA}}/(W)_{\text{RWA}} = h\omega_0 \coth(\beta h\omega_0/2)
\]

\[
+ h\omega_0\Gamma_\tau^2 \frac{\lambda_0^3}{6\hbar^3} (1 - e^{-\beta h\omega_0})
\]

\[
\times \left[ 1 - \Gamma_\tau \frac{\tau}{6} (1 + e^{-\beta h\omega_0}) \right]
\]

\[
+ O((\Gamma_\tau \tau)^3) + O((\lambda_0 \tau/h)^3).
\]

This expression is valid up to \( \Gamma_\tau \lambda_0/\hbar \lesssim 0.01\lambda_0 \) in Fig. 3 as the higher-order terms become important already when \( \Gamma_\tau \lambda_0/\hbar = 0.01\lambda_0 \), due to the high number of drive cycles.

IV. SUMMARY AND CONCLUSIONS

In summary, we have examined in detail the distribution of work done when a two-measurement protocol is applied to a driven open quantum system. To this end, we have first derived a general form for the generating function of work and studied the first three moments of work by using the master equation of the reduced system and invoking approximations similar to the ones made in the microscopic derivation of the reduced density matrix. We have compared our results to the earlier derivations [24] that were carried out implicitly assuming that the total Hamiltonian and its time derivative commute and have shown that there is a significant difference already in the case of the third moment. This emphasizes the importance of properly evaluating the higher moments of work, which are needed to check fluctuation relations such as the JE. To make our results concrete, we have considered a weakly driven and weakly coupled two-level system by using a number of different techniques, including the quantum jump method. Our results demonstrate the influence of the correct choice of the generating function already in the results for the third moment of work distribution.

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APPENDIX A: GENERATING FUNCTION OF THE TWO-POINT MEASUREMENT PROTOCOL

In the two-measurement protocol for a closed quantum system, the probability to measure \( E_0 \) at time \( t = 0 \) and \( E_\tau \) at \( t = \tau \) is of the form

\[
P[E_\tau, E_0] = \text{Tr}\left\{ \tilde{P}_{E_\tau} \tilde{U}_\tau (\tau,0) \tilde{P}_{E_0} \tilde{P}_{E_\tau} \tilde{U}_\tau^{\dagger}(0,0) \tilde{P}_{E_\tau} \right\},
\]

where \( \tilde{U}_\tau (\tau,0) = T_\tau \exp \left\{-i \int_0^\tau dt \tilde{H}(t) \right\} \) is the unitary time evolution operator, \( T_\tau \) describes the chronological time ordering, and the projection operators are given by \( \tilde{P}_{E_\tau} = |E_\tau \rangle \langle E_\tau| \), where \( |E_\tau \rangle \) is the state corresponding to the measurement result \( E_\tau \) at time \( t \). The corresponding generating function is given by [24]

\[
G(u) = \sum_{E_0, E_\tau} e^{iuE_\tau - E_0} P[E_\tau, E_0] = \sum_{E_\tau} \text{Tr} \left\{ \tilde{U}_\tau (\tau,0) \sum_{E_0} e^{iuE_0} \tilde{P}_{E_0} \tilde{P}_{E_\tau} \tilde{U}_\tau^{\dagger}(0,0) \tilde{P}_{E_\tau} e^{iuE_\tau} \right\}
\]

\[
= \text{Tr} \left\{ \tilde{U}_\tau (\tau,0) e^{-i(u/2)E_\tau} \tilde{P}_{E_\tau} \tilde{P}_{E_0} e^{-i(u/2)E_0} \tilde{U}_\tau^{\dagger}(0,0) \tilde{P}_{E_\tau} e^{iuE_\tau} \right\}
\]

\[
= \text{Tr} \left\{ \tilde{U}_{\tau/2}(\tau,0) \tilde{P}_{E_\tau} \tilde{U}_{\tau/2}^{\dagger}(0,\tau) \right\}.
\]

where

\[
\tilde{U}_\mu (\tau,0) = e^{iuH(0)} \tilde{U}_\mu (\tau,0) e^{-iuH(0)},
\]

\[
\tilde{P}_{\text{BE}} = \sum_{E_\tau} \tilde{P}_{E_\tau} \tilde{P}_{E_\tau},
\]

and \( \tilde{P}_{\text{BE}} \) is the initial density matrix. If the initial density matrix is diagonal in the first measurement’s basis, then \( \tilde{P}_{\text{BE}} = \tilde{P}_{\text{BE}} \). In the case of energy measurement, this means that if the total system Hamiltonian \( \tilde{H}(0) \) commutes with \( \tilde{P}_{\text{BE}} \), e.g., the density matrix is diagonal in the eigenbasis of \( \tilde{H}(0) \), then \( \tilde{P}_{\text{BE}} = \tilde{P}_{\text{BE}} \).

With the assumption \( \{ \tilde{H}(t), \tilde{H}(t) \} = 0 \), the evolution operator \( \tilde{U}_\mu (\tau,0) \) satisfies the following equation of motion:

\[
\frac{d}{d\tau} \tilde{U}_\mu (\tau,0) = -i \hbar \left[ \tilde{H}(\tau) - \hbar \frac{\partial \tilde{H}(\tau)}{\partial \tau} \right] \tilde{U}_\mu (\tau,0).
\]
Let us denote the time derivative of the total Hamiltonian as
\[
\dot{U}_u(0,0) = \frac{\partial}{\partial \tau} \begin{array}{c} \nu \frac{\partial}{\partial \epsilon} \hat{H}(t) \end{array}.
\] (A6)

However, contrary to Ref. [24], this solution [41] is not the general one due to the implicit assumption that \([\hat{H}(t), \partial \hat{H}(t)] = 0\). With this form of \(\dot{U}_u(0,0)\), the generating function simplifies to
\[
G_0(u) = \text{Tr} \left[ \tau_u \exp \left\{ -\frac{i}{\hbar} \int_0^\tau dt \left[ \hat{H}(t) - \nu \frac{\partial}{\partial \epsilon} \hat{H}(t) \right] \right\} \right].
\]

Let us denote the time derivative of the total Hamiltonian as the power operator \(\hat{P}(t) = \partial \hat{H}(t)/\partial \epsilon\). In order to get an expression where the operators are expressed in the Heisenberg picture, we can use the unitarity of \(\hat{U}u(0,0)\) and calculate the equation of motion for the operators \(\hat{U}u(0,0)\) and \(\hat{U}u(0,0)^T\). Changing to this Heisenberg picture and using the periodicity of the trace then gives the final form,
\[
G_0(u) = \text{Tr} \left[ \tau_u \exp \left\{ -\frac{i}{\hbar} \int_0^\tau dt \left[ \hat{H}(t) - \nu \frac{\partial}{\partial \epsilon} \hat{H}(t) \right] \right\} \right].
\]

After changing to the Heisenberg picture described earlier, the exact generating function reads
\[
G(u) = \text{Tr} \left[ \tau_u \exp \left\{ \int_0^\tau dt \sum_{n=1}^{\infty} (-1)^{n+1}(iu)^n \hat{C}_n(t) \right\} \right].
\]

APPENDIX B: CALCULATION OF THE MASTER EQUATION

Let us denote the density matrix of the total system with \(\hat{P}(t)\). The density matrix of the reduced system \(\hat{\rho}(t)\) is obtained by tracing over the bath degrees of freedom,
\[
\hat{\rho}(t) = \text{Tr}_B \{ \hat{P}(t) \}. \quad (B1)
\]

Similarly, the density matrix of the bath \(\hat{\rho}_B(t)\) is obtained by tracing over the system degrees of freedom,
\[
\hat{\rho}_B(t) = \text{Tr}_S \{ \hat{P}(t) \}. \quad (B2)
\]

The Hamiltonian of the total closed system can be written as
\[
\hat{H}(t) = \hat{H}_0 + \hat{H}_B + \hat{V}(t) + \hat{H}_C. \quad (B3)
\]

We will approximate the initial density matrix after the first measurement with \(\hat{\rho}_B(0) = \hat{\rho}_B(0) \otimes \hat{\rho}_B(0)\), where both the system and the heat bath start in thermal equilibrium. This approximation corresponds to that of neglecting the interaction Hamiltonian in the energy measurements. A similar approximation is done also in the calculation of the moments. Tracing over the bath degrees of freedom, we get the following equation for the reduced density matrix:
\[
\frac{d\hat{\rho}_B(t)}{dt} = -\frac{i}{\hbar} \left[ \hat{V}(t), \hat{\rho}_B(t) \right] - \frac{1}{\hbar^2} \int_0^t dt' \left[ \hat{H}_C(t'), \hat{V}(t') + \hat{H}_C(t'), \hat{\rho}_B(t') \right]. \quad (B4)
\]

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\hat{\rho}_B(t) = \text{Tr}_S \{ \hat{P}(t) \}. \quad (B2)
\]

The Hamiltonian of the total closed system can be written as
\[
\hat{H}(t) = \hat{H}_0 + \hat{H}_B + \hat{V}(t) + \hat{H}_C. \quad (B3)
\]

Let us change to the interaction picture with respect to \((\hat{H}_0 + \hat{H}_B)\), denoted by the superscript \(I\). We can write the equation of motion for the total density matrix as
\[
\frac{d\hat{\rho}_I(t)}{dt} = -\frac{i}{\hbar} \left[ \hat{V}(t), \hat{\rho}_I(t) \right] - \frac{1}{\hbar^2} \int_0^t dt' \left[ \hat{H}_C(t'), \hat{V}(t') + \hat{H}_C(t'), \hat{\rho}_I(t') \right]. \quad (B4)
\]

We will approximate the initial density matrix after the first measurement with \(\hat{\rho}_B(0) = \hat{\rho}_I(0) \otimes \hat{\rho}_B(0)\), where both the system and the heat bath start in thermal equilibrium. This approximation corresponds to that of neglecting the interaction Hamiltonian in the energy measurements. A similar approximation is done also in the calculation of the moments. Tracing over the bath degrees of freedom, we get the following equation for the reduced density matrix:
\[
\frac{d\hat{\rho}_I(t)}{dt} = -\frac{i}{\hbar} \left[ \hat{V}(t), \hat{\rho}_I(t) \right] - \frac{1}{\hbar^2} \int_0^t dt' \text{Tr}_B \left[ \left[ \hat{H}_C(t'), \hat{V}(t') + \hat{H}_C(t'), \hat{\rho}_I(t') \right] \right]. \quad (B5)
\]

Let us denote the last term on the right-hand side of Eq. (B5) as \(\chi(t)\). Invoking the Born approximation \(\hat{\rho}_I(t) = \hat{\rho}_I(t) \otimes \hat{\rho}_B(0)\) and the Markov approximation, it changes
to
\begin{equation}
\chi(t) = -\frac{1}{\hbar^2} \int_0^\infty dt' \text{Tr}_B \left[ \left\{ \hat{H}_C(t), \left[ \hat{V}(t'), \hat{\rho}(t') \otimes \hat{\rho}_B(0) \right] \right\} \right. \\
+ \left[ \hat{H}_C(t), \left[ \hat{\rho}(t'), \hat{\rho}(t) \otimes \hat{\rho}_B(0) \right] \right].
\end{equation}
(B6)

The interaction Hamiltonian can be expressed as \( \hat{H}_C(t) = \sum_j \hat{A}_j(t) \otimes \hat{B}_j(t) \), where \( \hat{A}_j(t) \) acts on the system degrees of freedom and \( \hat{B}_j(t) \) acts on the bath degrees of freedom. With this expression of \( \hat{H}_C(t) \) and assuming that \( \text{Tr}_B (\hat{B}_j(t) \hat{\rho}_B(0)) = 0 \), \( \chi \) changes to the form
\begin{equation}
\chi(t) = -\frac{1}{\hbar^2} \int_0^\infty dt' \sum_{j,k} \left[ \hat{A}_j(t) \hat{A}_j(t') \hat{\rho}(t) \right] \\
- \hat{\lambda}_j(t') \hat{\lambda}_j(t') \text{Tr}_B \left[ \hat{B}_j(t) \hat{B}_j(t') \hat{\rho}_B(0) \right] \\
+ \left[ \hat{\rho}(t') \hat{\lambda}_j(t') \hat{\lambda}_j(t) \right] \text{Tr}_B \left[ \hat{B}_j(t') \hat{B}_j(t) \hat{\rho}_B(0) \right].
\end{equation}
(B7)

For the system studied, the bath correlation functions are given by
\begin{equation}
\text{Tr}_B \left[ \hat{B}_j(t) \hat{B}_j(t') \hat{\rho}_B(0) \right] = \sum_k |g_k|^2 e^{i\omega_k(t-t')} n_k \\
+ e^{-i\omega_k(t-t')} (n_k + 1),
\end{equation}
where \( \hat{B}_j(t) = \sum_k e^{i\omega_k t} g_k \hat{a}_j \) and \( n_k \) is the average number of photons with frequency \( \omega_k \). The expression of \( \chi \) can be simplified by taking into account that \( \int_0^\infty \text{d}t e^{i\omega t} = \pi \delta(\omega) + i\mathcal{P} \left( \frac{\delta}{\omega} \right) \), where \( \mathcal{P} \) denotes the Cauchy principal value and the imaginary part only affects the Lamb shift. By neglecting the Lamb shift and invoking the secular approximation, i.e., neglecting the fast oscillating terms, we get
\begin{equation}
\chi(t) = \Gamma_1 \left( \rho_{\text{ss}}(t)|g\rangle\langle g| - \frac{1}{2}\{\hat{\rho}(t), |e\rangle\langle e|\} \right) \\
+ \Gamma_1 \left( \rho_{\text{ss}}(t)|e\rangle\langle e| - \frac{1}{2}\{\hat{\rho}(t), |g\rangle\langle g|\} \right),
\end{equation}
(B9)
where \( \rho_{\text{ss}}(t) = (|k\rangle\langle k|) \) and the transition rates are given by
\begin{equation}
\Gamma_1 = \frac{2\pi}{\hbar^2} \sum_k (\alpha_k + 1) |g_k|^2 \delta(\omega_k - 0),
\end{equation}
\begin{equation}
\Gamma_1 = \frac{2\pi}{\hbar^2} \sum_k n_k |g_k|^2 \delta(\omega_k - 0),
\end{equation}
and they satisfy the detailed balance \( \Gamma_1 = \Gamma_1 e^{-\beta \omega_0} \).

With the approximation \( \hat{\rho}_B(0) = \hat{\rho}'(0) \otimes \hat{\rho}_B(0) \), the second term on the right-hand side of Eq. (B5) goes to zero due to \( \text{Tr}_B (\hat{B}_j(t) \hat{\rho}_B(0)) = 0 \). Thus, switching back to the Schrödinger picture gives us the following master equation:
\begin{equation}
\frac{d\hat{\rho}}{dt} = -\frac{i}{\hbar} \left[ \hat{H}_S(t), \hat{\rho}(t) \right] \\
+ \Gamma_1 \left( \rho_{\text{ss}}(t)|g\rangle\langle g| - \frac{1}{2}\{\hat{\rho}(t), |e\rangle\langle e|\} \right) \\
+ \Gamma_1 \left( \rho_{\text{ss}}(t)|e\rangle\langle e| - \frac{1}{2}\{\hat{\rho}(t), |g\rangle\langle g|\} \right).
\end{equation}
(B12)

\section*{APPENDIX C: CALCULATION OF \( \langle W^3 \rangle_{S+B} \)}

Using the same notation as in the derivation of the master equation, we can write the total density matrix in the interaction picture with respect to \((\hat{H}_S + \hat{H}_B)\) as
\begin{equation}
\hat{\rho}'(t) = \hat{\rho}'(0) - \frac{i}{\hbar} \int_0^t dt' \left[ \hat{H}_S(t') + \hat{V}(t'), \hat{\rho}(t') \right].
\end{equation}
(C1)

With this form of \( \hat{\rho}'(t) \), the term inside the integral in Eq. (22) can be written as
\begin{align}
\Xi(t) &= \text{Tr}_{S+B} \left[ \left[ \hat{H}_S(t'), \hat{H}_S(t), \hat{\rho}'(t) \right] \right] \\
&= \text{Tr}_{S+B} \left[ \left[ \hat{H}_S(t'), \hat{H}_S(t), \hat{\rho}'(t) \right] \right] \\
&= -\frac{i}{\hbar} \int_0^t dt' \text{Tr}_{S+B} \left[ \left[ \hat{H}_S(t'), \hat{H}_S(t), \hat{\rho}'(t) \right] \right] \\
&\times \left[ \hat{H}_S(t') + \hat{V}(t'), \hat{\rho}(t') \otimes \hat{\rho}_B(0) \right].
\end{align}
(C2)

Again, we will approximate the initial density matrix with \( \hat{\rho}'(0) = \hat{\rho}'(0) \otimes \hat{\rho}_B(0) \), where both the system and the heat bath start in thermal equilibrium. Using the Born approximation \( \hat{\rho}'(t) = \hat{\rho}'(t) \otimes \hat{\rho}_B(0) \), we can approximate \( \Xi(t) \) with
\begin{align}
\Xi(t) &= \left[ \hat{H}_S(t'), \hat{H}_S(t), \hat{\rho}'(t) \right] \\
&= -\frac{i}{\hbar} \int_0^t dt' \text{Tr}_{S+B} \left[ \left[ \hat{H}_S(t'), \hat{H}_S(t), \hat{\rho}'(t) \right] \right] \\
&\times \left[ \hat{H}_S(t') + \hat{V}(t'), \hat{\rho}(t') \otimes \hat{\rho}_B(0) \right].
\end{align}
(C3)

The interaction Hamiltonian can be written as \( \hat{H}_C(t) = \sum_j \hat{A}_j(t) \otimes \hat{B}_j(t) \), where \( \hat{A}_j(t) \) acts on the system degrees of freedom and \( \hat{B}_j(t) \) acts on the bath degrees of freedom. Assuming \( \text{Tr}_B (\hat{B}_j(t) \hat{\rho}_B(0)) = 0 \), \( \Xi(t) \) reduces to
\begin{align}
\Xi(t) &= -\frac{i}{\hbar} \int_0^t dt' \text{Tr}_{S+B} \left[ \left[ \hat{H}_S(t'), \hat{H}_S(t), \hat{\rho}'(t) \right] \right] \\
&\times \left[ \hat{H}_S(t') + \hat{V}(t'), \hat{\rho}(t') \otimes \hat{\rho}_B(0) \right].
\end{align}
(C4)

Invoking the Markov approximation, the expression changes to
\begin{align}
\Xi(t) &= -\frac{i}{\hbar} \int_0^t dt' \text{Tr}_{S+B} \left[ \left[ \hat{H}_S(t'), \hat{H}_S(t), \hat{\rho}'(t) \right] \right] \\
&\times \left[ \hat{H}_S(t') + \hat{V}(t'), \hat{\rho}(t') \otimes \hat{\rho}_B(0) \right].
\end{align}
(C5)

Expressing the interaction Hamiltonian as \( \hat{H}_C(t) = \sum_j \hat{A}_j(t) \otimes \hat{B}_j(t) \) and denoting \( \hat{Q}_j(t) = [\hat{A}_j(t), \hat{H}_S(t), \hat{\rho}(t)] \), Eq. (C5) changes to the form
\begin{align}
\Xi(t) &= -\frac{i}{\hbar} \sum_{j,k} \int_0^t dt' \text{Tr}_S \left[ \hat{Q}_j(t') \hat{A}_j(t') \hat{\rho}(t') \right] \\
&\times \text{Tr}_B \left[ \hat{B}_j(t') \hat{B}_j(t') \hat{\rho}_B(0) \right] \\
&- \text{Tr}_S \left[ \hat{A}_j(t') \hat{Q}_j(t') \hat{\rho}(t') \right] \\
&\times \text{Tr}_B \left[ \hat{B}_j(t') \hat{B}_j(t') \hat{\rho}_B(0) \right].
\end{align}
(C6)
For the system studied, $\Xi(t)$ reduces to

$$\Xi(t) = -i2\omega_0\lambda(t)\left[\rho_{\varepsilon\varepsilon}(t)\int_0^\infty dt\, e^{-i\omega_0t} \xi(t,t') - \rho_{\varepsilon\varepsilon}(t)\int_0^\infty dt\, e^{i\omega_0t} \xi(t,t')\right], \quad (C7)$$

where $\rho_{\varepsilon\varepsilon}(t) = \langle \varepsilon | \hat{\rho}(t) | \varepsilon \rangle$ and the term $\xi(t,t') = \text{Tr}_B(\dot{\hat{B}}(t)\hat{B}(t')\hat{\rho}_B(0)) + \text{Tr}_B(\dot{\hat{B}}(t')\hat{B}(t)\hat{\rho}_B(0))$. Neglecting the Lamb shift, we get

$$\Xi(t) = -i\hbar^2\omega_0\lambda(t)\left[\rho_{\varepsilon\varepsilon}(t)e^{-i\omega_0t} - \rho_{\varepsilon\varepsilon}(t)e^{i\omega_0t}\right](\Gamma_1 + \Gamma^-)$$

$$= 2i\hbar^2\omega_0\lambda(t)\text{Im}[\rho_{\varepsilon\varepsilon}(t)](\Gamma_1 + \Gamma^-). \quad (C8)$$

With this form of $\Xi(t)$, $\langle W^3 \rangle_{S+B}$ reduces to

$$\langle W^3 \rangle_{S+B} \approx \frac{\hbar^2\omega_0}{2} (\Gamma_1 + \Gamma^-) \int_0^\infty dt\, \lambda(t)\text{Im}[\rho_{\varepsilon\varepsilon}(t)]. \quad (C9)$$

[30] $P[E_1,E_0]$ can also be written as $P[E_1,E_0] = \langle |E_1\rangle \langle \hat{U}(\tau,0)|E_0\rangle \rangle^2 |E_0\rangle \langle \rho_0 |E_0\rangle$. However, in the main text, we use the trace form as it is useful in the following calculations.
[41] Note that if the observable $\hat{A}(t) = \hat{H}(t)$, then $e^{i\lambda_0(t)}\hat{H}(t)e^{-i\lambda_0(t)} = \hat{H}(t)$ in Eq. (22) in Ref. [24].