Quantum Trajectories for Brownian Motion

Walter T. Strunz$^1$, Lajos Diósi$^2$, Nicolas Gisin$^3$, and Ting Yu$^3$
$^1$Fachbereich Physik, Universität GH Essen, 45117 Essen, Germany
$^2$Research Institute for Particle and Nuclear Physics, 1525 Budapest 114, POB 49, Hungary
$^3$Group of Applied Physics, University of Geneva, 1211 Geneva 4, Switzerland
(August 2, 1999)

We present the stochastic Schrödinger equation for the dynamics of a quantum particle coupled to a high temperature environment and apply it to the dynamics of a driven, damped, nonlinear quantum oscillator. Apart from an initial slip on the environmental memory time scale, in the mean, our result recovers the solution of the known non-Lindblad quantum Brownian motion master equation. A remarkable feature of our powerful stochastic approach is its localization property: individual quantum trajectories remain localized wave packets for all times, even for the classically chaotic system considered here, the localization being stronger as $\hbar \to 0$.

03.65.Bz, 42.50.Lc, 05.40.+j

The understanding of the dynamics of open or dissipative quantum systems is of fundamental importance both from a practical and conceptual point of view. The archetype of such a system is the standard quantum Brownian motion model [1] which describes a particle with Hamiltonian $H(q,p)$, coupled to an environment of harmonic oscillators $(q_\lambda,p_\lambda)$ via its position $q$, such that the total Hamiltonian of system and environment reads

$$H_{\text{tot}}(q,p,q_\lambda,p_\lambda) = H(q,p) + \sum_\lambda \left\{ \frac{p_\lambda^2}{2m_\lambda} + \frac{1}{2}m_\lambda\omega_\lambda^2(q_\lambda - \frac{g_\lambda}{m_\lambda\omega_\lambda^2}q)^2 \right\}.$$ (1)

Up to now, in order to determine the time dependent dynamics of the open ‘system’, the standard procedure was the derivation of a master equation for the reduced density operator, which, for the high temperature case considered below, is widely accepted to read

$$\hbar \dot{\rho}_t = -i[H,\rho_t] - \frac{\gamma}{2}[q,\{p,\rho_t\}] - \frac{m_\gamma kT}{\hbar} [q, [q, \rho_t]],$$ (2)

where $\gamma$ is the damping rate. This master equation is a Markov master equation not, however, of Lindblad form [2] and indeed it turns out that it may violate the positivity of $\rho_t$ on very short time scales, which has led to an ongoing debate about its range of applicability [3]. We will briefly address this issue later on in this Letter.

Our new approach to quantum Brownian motion is very different and circumvents the derivation of a master equation for $\rho_t$ altogether. Instead, we use a stochastic Schrödinger equation, derived straight from the microscopic model (1), for pure states $\psi_t(z)$ (quantum trajectories). Our construction recovers the reduced density operator as the ensemble mean $M[\ldots]$ over many of these quantum trajectories, in principle without any approximation:

$$\rho_t = M [\psi_t(z)\langle \psi_t(z) \rangle].$$ (3)

The mean $M[\ldots]$ is taken over the process $z_t$ which drives the stochastic Schrödinger equation. We are thus able to determine $\rho_t$ in a Monte-Carlo sense without an explicit master equation for its time evolution.

Quantum trajectory methods have been used extensively in recent years, mainly in the quantum optics community, due to their numerical efficiency, their intimate connection to (continuous) measurement, and their illustrative power helping to gain physical insight. The master equations encountered in quantum optics are of standard Lindblad type, for which Markov quantum trajectory methods are known for some time now: there are jump processes [4] and diffusive processes [5] recovering the reduced density operator. Despite being maybe the best known of all master equations, the Quantum Brownian motion master equation (2), being not of Lindblad form, has so far been excluded from a treatment with these powerful methods.

Only recently the authors managed to extend the quantum trajectory concept to non-Markovian situations [6], more precisely, we were able to determine a stochastic Schrödinger equation for the dynamics of a quantum system coupled to a bath of harmonic oscillators as in (1), without using the concept of a master equation for $\rho_t$. An alternative approach to non-Markovian quantum trajectories, more emphasizing the continuous measurement point of view, has now also been established [7].

In its linear version [8], our non-Markovian quantum state diffusion (QSD) stochastic Schrödinger equation for the quantum Brownian motion model (1) takes the form

$$\hbar \dot{\psi}_t(z) = -iH'\psi_t(z) + qz_\lambda\psi_t(z) - q \int_0^t ds \alpha(t,s) \frac{\delta \psi_t(z)}{\delta z_s},$$ (4)

where we assumed a factorized total initial density operator $\rho_{\text{tot}} = |\psi_0\rangle\langle \psi_0 | \otimes \rho_T$ with a pure system state $|\psi_0\rangle$ and an environmental thermal density operator $\rho_T$. The

\[\text{Page 1 of 2}\]
influence of the environment on the system is encoded in the bath correlation function $\alpha(t,s) = \langle F(t)F(s) \rangle_{\rho_T}$ where $F(t) = \sum_{\lambda} g_\lambda q_\lambda(t)$ is the quantum force in (1) and $z_s$ is thus a complex Gaussian stochastic c-number force with correlation $M[z_s^* z_s] = \alpha(t,s)$. In the usual high temperature limit $K T \gg h \Lambda \gg \hbar \omega, h \gamma$, where $\Lambda$ is an environmental cutoff frequency and $\omega, \gamma$ are the typical system frequency and damping rate, respectively, one finds [1]

$$\alpha(t,s) = 2m\gamma k T \Delta(t-s) + i\hbar m\gamma \Delta(t-s), \quad (5)$$

where $\Delta(t)$ is a delta-like function decaying on the environmental ‘memory’ time scale $\Lambda^{-1}$ (here we use $\Delta(t) = 2^3 e^{-\Lambda |t|}$). In (4), the Hamiltonian $H' = H(q,p) + \frac{1}{2} m\gamma \Lambda q^2$ contains an additional potential term that turns out to be counterbalanced by a similar term arising from the memory integral.

Eq. (4) is exact, i.e. it provides a quantum trajectory method for Brownian motion for any temperature and any distribution of environmental oscillators in the model (1), i.e. for any $\alpha(t,s)$. In order to compute numbers, however, we have to express the functional derivative under the memory integral in (4) in terms of elementary operators. In the high temperature limit considered here, we simply need to expand in terms of the time delay $(t-s)$

$$\frac{\delta \psi_t(z)}{\delta z_s} = \frac{1}{\hbar} \left( q - \frac{p}{m} (t-s) + \ldots \right) \psi_t(z), \quad (6)$$

where the dots denote terms of the order $(t-s)^2$ and higher, leading to corrections of the order $\omega/\Lambda, \gamma/\Lambda$ and can therefore be neglected (see [9] for a general theory of such ‘post-Markov’ open systems). With (6), the memory integral in (4) takes the form

$$\int_0^t ds \alpha(t,s) \frac{\delta \psi_t(z)}{\delta z_s} = (g_0(t)q - g_1(t)p) \psi_t(z), \quad (7)$$

where we introduce time dependent coefficients $g_0(t) = \frac{1}{\hbar} \int_0^t ds \alpha(t,s)$ and $g_1(t) = \frac{i}{\hbar m} \int_0^t ds (t-s) \alpha(t,s)$. The imaginary part of $g_0(t)$ will be compensated by the additional potential term in $H'$. The imaginary part of $g_1(t)$ gives rise to damping. The real part of $g_0(t)$ describes diffusion and as the real part of $g_1(t)$ also gives rise to diffusion, yet smaller by a factor $\omega/\Lambda$, the latter can be neglected compared to the former in the regime we are interested in.

In order to get an efficient Monte Carlo method (importance sampling [10]), we go over to the nonlinear version of (4), which keeps the trajectories $\psi_t(z)$ normalized at all times while retaining the correct ensemble mean (3), see [6]. Using (7), the relevant stochastic Schrödinger equation for Brownian motion reads

$$\hbar \dot{\psi}_t(z) = -iH \psi_t(z) - i \left( \frac{1}{2} m\gamma \Lambda + \text{Im} \{g_0(t)\} \right) q^2 \psi_t(z) \quad (8)$$

Normalized quantum trajectories $\psi_t(z)$ whose ensemble mean gives the desired reduced density operator according to (3) can now be propagated using (8), where $\langle q \rangle = \frac{\langle q \rangle}{\hbar}$, a quantity which has to be determined numerically along with $\psi_t(z)$ (very often the replacement $m\langle q \rangle \approx \langle p \rangle$ turns out to be a good approximation).

In (8), the time dependent coefficients quickly approach their asymptotic values $g_0(t) \to \frac{m\gamma k T}{\hbar} - \frac{i}{2} m\gamma \Lambda$ and $\text{Im} \{g_1(t)\} \to -\frac{i}{2}$ for times larger than the environmental memory time. After this initial slip $t \gg \Lambda^{-1}$, (8) becomes

$$\hbar \dot{\psi}_t(z) = -iH \psi_t(z) + (q - \langle q \rangle)z_t \psi_t(z)$$

$\quad - \text{Re} \{g_0(t)\} \left( (q - \langle q \rangle)^2 - (\langle q \rangle^2) \right) \psi_t(z)$

$\quad + \text{Im} \{g_1(t)\} \left( qp - \langle qp \rangle + m\langle q \rangle q - \langle p \rangle p \right) \psi_t(z). \quad (9)$

We now highlight the power of our stochastic Schrödinger equation for Brownian motion (8) by investigating the dynamics of a driven, damped, nonlinear, noisy system, the Duffing oscillator, where $H = \frac{1}{2} p^2 + \frac{1}{4} q^4 - \frac{1}{2} q^2 + gq \cos(t)$, here coupled to a heat bath at temperature $T$. This system has been studied before using the master equation (2) (see [11] and references therein), including a straight numerical solution which requires the propagation of a huge matrix. In our new approach, one propagates pure states $\psi_t(z)$ according to (8), a great reduction in resources, with the need, however, to solve (8) many times in order to evaluate the mean values. For Lindblad master equations, the power of quantum trajectory methods for investigating classically chaotic dissipative systems was shown in [12] (see also [14]).

We use parameters $g = 0.3$ with a damping rate $\gamma = 0.25$, thus the classical problem is chaotic [13]. The environment is furthermore characterized by $k T = 0.3$, and a cutoff frequency $\Lambda = 5$. With $h$ of the order $10^{-2}$ and smaller (see various choices of $h$ below), the parameters are in the required regime. As initial condition we choose a standard coherent state located at $\langle q \rangle = 0.1, \langle p \rangle = 0.1$.

In Fig.1 we show the ensemble mean $M[W_z(q,p,t)]$ over 1000, 5000, and 10000 Wigner functions of pure state trajectories $\psi_t(z)$ obtained solving (8) numerically up to a time $t = 4$. According to our construction, this quantity converges to the Wigner function of the reduced density operator for many realizations. Here we have chosen $h = 0.01$, a phase space area corresponding approximately to the extension of the wave packets shown in Fig.2.
FIG. 1. Contour plots of the Wigner function $W(q, p, t = 4)$ of the reduced density operator of the thermal Duffing oscillator with $\hbar = 0.01$ (for the phase space area corresponding to this $\hbar$ see Fig. 2). The contour plots show the ensemble mean over 1000, 5000, and 10000 Wigner functions $W_z(q, p, t = 4)$ of individual quantum trajectories obtained solving the quantum Brownian motion stochastic Schrödinger equation (8).

In Fig. 2 we show contour plots of Wigner functions $W_z(q, p, t = 4)$ of four realizations of (8), many of which add up to the Wigner function of the desired reduced density matrix shown in Fig. 1. One can see clearly that these individual Wigner functions are well localized in phase space compared to the phase space spread of the ensemble, even for this classically chaotic system.

FIG. 2. Contour plots of Wigner functions $W_z(q, p, t = 4)$ of four individual quantum trajectories obtained solving the quantum Brownian motion stochastic Schrödinger equation (8) for the thermal Duffing oscillator. Individual trajectories remain well localized in phase space with respect to the overall spread of the ensemble mean, even for this classically chaotic system. The chosen value of $\hbar = 0.01$ is slightly smaller than the phase space area covered by these states.

This remarkable feature of the quantum Brownian motion stochastic Schrödinger equation (8) is highlighted again in Fig. 3, where we show the mean position spread, $M[\Delta q] = M[\sqrt{\langle (q - \langle q \rangle)^2 \rangle}]$ and the mean uncertainty product $M[\Delta q \Delta p/\hbar]$ in units of $\hbar$ of individual trajectories as a function of time for three different choices of $\hbar$.

FIG. 3. Localization property of the QBM stochastic Schrödinger equation. Individual runs are well localized in phase space, the localization being stronger the smaller $\hbar$: (a) the average position spread $M[\Delta q] = M[\sqrt{\langle (q - \langle q \rangle)^2 \rangle}]$ of solutions of the QBM stochastic Schrödinger equation for the choices $\hbar = 0.01$ (solid line), $\hbar = 0.005$ (dashed line), and $\hbar = 0.001$ (dotted line). Fig. (c) shows the mean uncertainty product $M[\Delta q \Delta p/\hbar]$, which remains of the order one almost independently of $\hbar$. Thus the quantum trajectories remain almost minimum uncertainty wave packets for all times.

The quantities shown in Fig. 3 can only be given sense...
in the framework of quantum trajectories, they have no meaning from a density operator point of view as they are the ensemble mean over an expression non-quadratic in \( \psi_t(z) \). It is apparent from Fig.3 that individual trajectories are well localized in phase space for all times, the localization being stronger the smaller \( \hbar \). As can be seen, our quantum trajectories remain almost ‘classical’ states, yet recover the fully quantum master equation (2). Thus, the representation (3) expresses the reduced density operator of quantum Brownian motion explicitly as a mixture of almost ‘classical’ states.

The observed localization property of QSD is well known in the Markov case and has been studied for instance in [14]. Here we see that similar properties hold for the generalized non-Markovian QSD equation (8) which has now been applied to quantum dynamics beyond the class of Lindblad master equations. As in the Markov case, the localization property can be exploited to further reduce the numerical effort.

Finally, let us briefly address the connection between our approach and the widely used QBM master equation (2). Since a quantum trajectory approach strictly preserves positivity of the reduced density operator, our QBM stochastic Schrödinger equation (8) cannot be identical to (2) in the mean, as the latter is known to violate positivity on short time scales. Taking the ensemble mean \( M \ldots \) in (3) with (4) analytically, we were able to show in [9] that in the regime considered in this Letter, the evolution of the ensemble mean (3) is well described by the master equation

\[
\hbar \dot{\rho} = -i[H, \rho] - i \left( \frac{1}{2} \mu g \Lambda + \text{Im} \{g_0(t)\} \right) [q^2, \rho] + \text{Im} \{g_1(t)\} [g, \{p, \rho\}] - \text{Re} \{g_0(t)\} [\{g, g, \rho\}],
\]

which reduces to (2) for times larger than the environmental memory time, \( t \gg \Lambda^{-1} \) due to the asymptotics of the coefficients \( g_0(t), g_1(t) \). Thus, apart from an initial slip on the environmental memory time scale \( \Lambda^{-1} \), our approach recovers (2) in the mean. It is known in the case of the exact master equation for a damped harmonic oscillator [15] that such time dependent coefficients may ensure the positivity of the reduced density operator for non-Lindblad master equations, a result that is here supported for general system Hamiltonian \( H(q, p) \).

To conclude, we have presented the stochastic Schrödinger equation for Brownian motion. It is compatible with the standard QBM master equation yet allows to compute states rather than a matrix, a huge reduction in resources, which becomes even more relevant for QBM in more than one space dimension. Individual trajectories are well localized in phase space, the localization being stronger the smaller \( \hbar \). Thus, in (3), the reduced density operator for Brownian motion is explicitly represented as an ensemble of almost ‘classical’ states.

We thank F Haake and IC Percival for helpful comments. WTS would like to thank the Deutsche Forschungsgemeinschaft for support through the SFB 237 “Unordnung und große Fluktuationen”. NG and TY thank the Swiss National Science Foundation.