Quantum phase transition and correlations in the multi-spin-boson model

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We consider multiple noninteracting quantum mechanical two-level systems coupled to a common bosonic bath and study its quantum phase transition with Monte Carlo simulations using a continuous imaginary time cluster algorithm. The common bath induces an effective ferromagnetic interaction between the otherwise independent two-level systems, which can be quantified by an effective interaction strength. For degenerate energy levels above a critical value of the bath coupling strength α all two-level systems freeze into the same state and the critical value α_c decreases asymptotically as 1/N with increasing N. For a finite number N of two-level systems the quantum phase transition (at zero temperature) is in the same universality class as the single spin-boson model; in the limit $N \rightarrow \infty$ the system shows mean-field critical behavior independent of the power of the spectral function of the bosonic bath. We also study the influence of a spatial separation of the spins in a bath of bosonic modes with linear dispersion relation on the location and characteristics of the phase transition as well as on correlations between the two-level systems.

DOI: 10.1103/PhysRevB.90.224401

PACS number(s): 05.10.Ln, 05.30.Jp, 03.65.Yz, 75.40.Mg

I. INTRODUCTION

The single spin-boson model [1,2] describes a two-level system, such as a spin-1/2 or a q-bit, which is linearly coupled to a bath of bosonic modes. In spite of its simplicity the system shows a multifaceted behavior in statics, dynamics, and quantum criticality, for which reason it became a paradigmatic model in the field of quantum dissipation. For Ohmic and sub-Ohmic dissipation the two-level system without energy bias shows a transition at zero temperature from a delocalized state (tunneling between the two states) at weak coupling to the bath to a localized, classical behavior (frozen in one state). The characteristics, in particular the universality class, of this quantum phase transition are identical to those of the thermal transition in the classical Ising chain with long-range interactions [3–5].

An interesting question is, what happens to this transition if one couples several two-level systems, such as impurities or qbits, to a bosonic dissipative bath? The simplest generalization of this kind is a pair of two-level systems, which was studied first in [6] and subsequently in [7-14]. In many studies of the two-spin-boson model an additional ferromagnetic coupling between the two-level systems of spins has been taken into account. Arranging N two-level systems in a chain with uniform ferromagnetic nearest-neighbor interactions and coupling each two-level system to its own bosonic bath yields the transverse Ising chain with dissipation that has been studied for an Ohmic bath in [15]. There it was shown that in the limit $N \to \infty$ a new quantum phase transition triggered by the bath coupling strength emerges, which is different from the quantum phase transition in the transverse Ising chain without dissipation and different from the single spin-boson model. If the ferromagnetic couplings between the two-level systems are not uniform but random the random transverse Ising chain with dissipation is obtained, which has been studied in [16,17]. Here not a sharp phase transition occurs but a smeared transition in which connected clusters of two-level systems undergo separate quantum phase transitions at different coupling strengths [17,18].

These earlier investigations for arbitrary two-level system number N assume an independent bath for each two-level system. This might not always represent correctly the physical situation, in which a single bath for all two-level systems might be more appropriate, as for instance in cold atom or trapped ion setups. In addition a direct coupling between the two-level systems might be absent. This is the situation which we address in this paper by studying the multi-spin-boson model (MSBM) with a single bosonic bath. One expects that in spite of the absence of a direct interaction between the two-level systems the common bath will mediate an effective interaction between the two-level systems that enhances the tendency towards localization and thus decreases the critical bath coupling strength. A related model is the large-spinboson model, where a single large spin (J > 1/2) is linearly coupled to a bosonic environment [19-21]. The large spin can be considered as an ensemble of 2J noninteracting spin-1/2. In [21] it was shown that the model resembles the spin-boson model phase diagram by rescaling the coupling as $\alpha'_c = (2J)^2 \alpha_c$. Since this model acts only on a 2J + 1dimensional subspace in contrast to a 2^{2J} dimensional Hilbert space in the multi-spin-boson model one might expect a slightly modified behavior. In Sec. III A we show that the universality class of the transition in the large-spin-boson model remains the same as the one of the multi-spin-boson model but the critical coupling strength α_c is shifted.

For N = 2 and Ohmic bath a recent variational calculation [13] obtained a critical coupling of $\alpha_c \approx 1/2$, whereas a numerical renormalization group (NRG) calculation [12] predicted (for vanishing direct coupling between the two-level systems) $\alpha_c \approx 1/4$. Regarding such a large deviation between the naive mean-field prediction of $\alpha_c(N) \propto 1/N$ and the NRG result for N = 2 a close look at larger values for N using exact methods such as quantum Monte Carlo seems worthwhile. Therefore in this paper we analyze quantitatively in which way the location and characteristics of the quantum phase transition varies with the number N and the separation of two-level systems and also shed light on the correlations between the two-level systems mediated by the bath in the different phases.

The paper is organized as follows: In Sec. II we introduce the MSBM including its path-integral representation and discuss the zero-tunneling limit as well as the mean-field limit. The quantum phase transition in the MSBM model is analyzed in Sec. III for general number of two-level systems N, first theoretically and then with the help of extensive quantum Monte Carlo simulations. In Sec. IV the spin-spin correlations in the MSBM in the different phases are discussed and in Sec. V the effect of spatial separations of the two-level systems in the MSBM is analyzed. Section VI contains a summary and a discussion of open questions. Several details of the calculations as well as a detailed description of the quantum Monte Carlo algorithm are deferred to three Appendices.

II. THE MODEL AND MEAN-FIELD THEORY

For completeness and for further reference we first summarize the characteristics of the quantum phase transition in the single spin-boson model. Its Hamiltonian is given by ($\hbar = 1$)

$$H = \frac{\Delta}{2}\sigma^{x} - \frac{\epsilon}{2}\sigma^{z} + \sum_{k}\omega_{k}a_{k}^{\dagger}a_{k} + H_{I}, \qquad (1)$$

where $\sigma^{x,z}$ are the Pauli spin-1/2 operators, $a_k^{\mathsf{T}}(a_k)$ are the bosonic creation (annihilation) operators, Δ the bare tunneling amplitude, ω_k the frequency of the *k*th bath mode, and ϵ an energy bias of one of the two states, which we set to zero ($\epsilon = 0$). The interaction term

$$H_I = \frac{1}{2} \sum_k (\lambda_k a_k^{\dagger} + \lambda_k^* a_k) \sigma^z \tag{2}$$

describes the linear coupling between the spin and its environment. The bath spectrum $J(\omega) = \pi \sum_k |\lambda_k|^2 \delta(\omega - \omega_k)$ has a power-law form $J(\omega) = 2\pi \alpha \omega^s \omega_c^{1-s} \theta(\omega_c - \omega)$, with the bath exponent s > 0, the coupling strength α , and a sharp cutoff at the frequency ω_c .

The most prominent case is the Ohmic spectrum, s = 1, which has a phase transition in the Kosterlitz-Thouless universality class between a delocalized phase with finite effective renormalized tunneling amplitude Δ_{eff} and a localized phase, where tunneling is completely suppressed, at $\alpha_c = 1 + O(\frac{\Delta}{\omega_c})$ [1,2]. In the sub-Ohmic regime (0 < s < 1), the phase transition is of second order, which is described by a Gaussian fixed point with "classical" exponents for $0 < s \leq \frac{1}{2}$ and has nonclassical, *s*-dependent critical exponents for $\frac{1}{2} < s < 1$ [3,4].

In this paper we consider the generalization of (1) to N two-level systems coupled to a common bath. The Hamiltonian of this multi-spin-boson model (MSBM) is given by

$$H = \sum_{m=1}^{N} \left(\frac{\Delta}{2} \sigma_m^x - \frac{\epsilon}{2} \sigma_m^z \right) + \sum_k \omega_k a_k^{\dagger} a_k + H_I \qquad (3)$$

with the interaction term

$$H_I = \sum_{m=1}^N \frac{1}{2} \sum_{k=-\infty}^\infty (\lambda_k a_k^{\dagger} + \lambda_k^* a_k) \sigma_m^z.$$
(4)

Note that this setup is different from the dissipative spin chain [15,17,18], where every spin is embedded in its own dissipative bath. The main effect of the common bath is that each spin is simultaneously influenced by a polarization of the bosonic degrees of freedom. From this picture one gets the intuition that the localization of the multiple spin

system is enhanced in comparison to the single spin case. The delocalized-to-localized transition point α_c is hence expected to become lowered as the number of contributing spins is increased.

A. Path-integral representation

The partition function of the single spin-boson model can be exactly written as a path integral [3], which is straightforward to generalize to the multi-spin-boson model; see Appendix A. The partition function Z of the quantum system is expressed as a sum over all possible spin- $\frac{1}{2}$ world lines $s(\tau) \in \{-1,1\}$ with $0 \le \tau < \beta$:

$$Z = \operatorname{Tr}[\exp(-\beta H)] \propto \int \prod_{m=1}^{N} \mathcal{D}[s_m(\tau)] \, \exp(-S).$$
 (5)

The positions of the kinks, where $s(\tau)$ changes its sign, are Poissonian distributed and the effective action (for symmetric states, i.e., $\epsilon = 0$) is

$$S = -\frac{1}{2\beta} \sum_{m=1}^{N} \sum_{m'=1}^{N} \left[\int_{0}^{\beta} \int_{0}^{\tau} s_{m}(\tau) s_{m'}(\tau') K_{\beta}(\tau - \tau') d\tau' d\tau \right],$$
(6)

where the integral kernel is given by

$$K_{\beta}(\tau) = \int_{0}^{\omega_{c}} \frac{J(\omega)}{\pi\omega} \left\{ \frac{\beta\omega}{2} \frac{\cosh\left[\frac{\beta\omega}{2}\left(1 - \frac{2\tau}{\beta}\right)\right]}{\sinh\left(\frac{\beta\omega}{2}\right)} \right\} d\omega.$$
(7)

This function decays algebraically as $\propto \tau^{-1-s}$ for $\beta \to \infty$ and the finite cutoff frequency ω_c ensures the convergence at $\tau = 0$.

The path-integral representations of thermodynamic observables such as the order parameter (magnetization) m, the susceptibility χ , and the dimensionless moment ratio Q of the MSBM are then given by

$$m = \left\langle \sigma_i^z \right\rangle_{\text{MSBM}} = \left\langle |M_i| \right\rangle_{\text{PI}},\tag{8}$$

$$\chi = 2 \frac{\partial m}{\partial \epsilon} \bigg|_{\epsilon=0} = \beta \left(\left| M_i^2 \right|_{\rm PI} - m^2 \right), \tag{9}$$

$$Q = \langle M_i^2 \rangle_{\rm PI}^2 / \langle M_i^4 \rangle_{\rm PI}, \qquad (10)$$

where M_i is the magnetization of the *i*th world line,

$$M_i[s_i(\tau)] = \beta^{-1} \int_0^\beta d\tau \, s_i(\tau)$$
 (11)

(note that the observables m, χ , and Q are independent of the index i = 1, ..., N) and $\langle \cdots \rangle$ denotes expectation values with respect to the classical action,

$$\langle \mathcal{O} \rangle_{\text{PI}} = \int \prod_{m=1}^{N} \mathcal{D}[s_m(\tau)] \mathcal{O} \exp(-S).$$
 (12)

Our quantum Monte Carlo cluster algorithm [3] (see also Appendix B) simply samples stochastically world line configurations according to the probability measure and the classical action.

Let us consider the zero-tunneling limit $\Delta = 0$. This simplifies the terms in (6) and (7) drastically, because no kinks can occur and hence the spin world lines are only spin variables $s_m \in \{-1,1\}$ without any τ dependence. The path integral becomes a sum over these variables and the double integral over the curly brackets in Eq. (7) is just $\beta^2/2$. The remaining frequency integral over the spectral function can performed elementarily and the resulting partition function becomes

$$Z = \sum_{s_1,\dots,s_N=\pm 1} \exp\left(\frac{\beta}{2} \sum_{m=1}^N \sum_{m'=1}^N s_m s_{m'} \frac{\alpha \omega_c}{s}\right).$$
(13)

This partition function is identical to the one for a classical Ising model in which N Ising spins interact with each other with the ferromagnetic coupling strength

$$J = \omega_c \alpha / s \tag{14}$$

between all distinct spin pairs. This result can also easily be obtained by a polaron transformation $H' = U^{-1}HU$ of the form $U = \exp\{-\sum_{k} [\lambda_k/(2\omega_k)a_k^{\dagger}\sum_m \sigma_m^z - \text{H.c.}]\}$. The thermodynamic limit $(N \to \infty)$ is not yet well-defined, because the energy would grow quadratic with the system size. To ensure a proper limit, we introduce the scaled coupling strength

$$\tilde{\alpha} = \alpha N. \tag{15}$$

Thus, for zero tunneling strength the MSBM model is identical to the exactly solvable classical mean-field model [22], which provides a paramagnetic to ferromagnetic phase transition of mean-field type. Adapted to the present nomenclature in (13), the transition point is

$$\tilde{\alpha}_c = \frac{s}{\omega_c} T.$$
 (16)

C. Mean-field theory

The mean-field approximation consists of neglecting terms that are quadratic in the fluctuations of the order parameter $m = \langle \sigma^z \rangle$, which means one sets $(\sigma_i^z - m)(\sigma_j^z - m) \approx 0$. Since the integral in Eq. (7) can be computed for $\tau = \beta$ and yields $K^{(1)}(\beta) = 2\alpha\omega_c/s$, the action (6) is in mean-field approximation

$$S_{MF} = \sum_{n=1}^{N} \left[\frac{\beta \tilde{\alpha} \omega_c}{2s} m^2 - \frac{1}{2} \int_0^\beta \left(\epsilon + \frac{2 \tilde{\alpha} \omega_c}{s} m \right) s_n(\tau) \right].$$
(17)

Solving the self-consistent equation $\frac{m}{2} = \frac{1}{N\beta} \frac{\partial \ln(Z)}{\partial \epsilon}$, one obtains a critical coupling strength

$$\tilde{\alpha}_c = \frac{s\Delta}{2\tanh(\beta\Delta/2)\omega_c},\tag{18}$$

which is identical to the result in Ref. [23] for N = 1. Note that the mean-field approximation predicts a phase transition also at nonvanishing temperatures (T > 0, i.e., $\beta < \infty$), which is absent in the MSBM for any finite N (see next subsection). Within the mean-field approximation the dependence of the critical coupling α_c on the number of spins N is fully compensated by the rescaling of the coupling strength (15) and Eq. (18) predicts a phase transition at the temperature

$$T_c = \frac{\Delta}{2 \operatorname{artanh}[s\Delta/(2\tilde{\alpha}_c \omega_c)]}.$$
 (19)

At zero temperature one gets $\tilde{\alpha}_c(T \to 0) = s\Delta/(2\omega_c)$. For zero tunneling one obtains $T_c(\Delta \to 0) = \tilde{\alpha}\omega_c/s$, which is identical to the exact result (16) as expected, since for zero tunneling the spins do not fluctuate and therefore the meanfield approximation is exact.

Also in the limit $N \to \infty$ one expects mean-field theory to be exact (see Appendix C), which we checked with extensive QMC simulations by finite N scaling at fixed temperatures (see next subsection).

III. THE QUANTUM PHASE TRANSITION IN THE MSBM

In this section, we study the quantum phase transition (at T = 0, i.e., $\beta = \infty$, and vanishing bias $\epsilon = 0$) of the MSBM for different values of N. First we argue in the next subsection that the universality class is independent of N; then we validate this prediction with finite β scaling of QMC data.

A. Universality class: Theoretical considerations

The universality class of the transition of the MSBM for general N is not known and before we embark on a QMC study we argue in the following that for any finite N the transition is in the same universality class as the single spin-boson model (N = 1). Consider a single world line of a 1/2-spin in a transverse field $\Delta \sigma^x/2$. The field defines a characteristic length 2/ Δ between two kinks [26]. The mean number of segments of the world line is hence $L = \beta \Delta/2$. Therefore, the partition function (5), (6) can approximately be written for large β in a discretized form

$$Z = \sum_{\{s_{m,i}\}} \exp\left[\frac{1}{2\beta} \sum_{m,m'}^{N} \frac{1}{2} \sum_{i,j}^{L} s_{m,i} s_{m',j} K_{\beta} \left(|j-i|\frac{2}{\Delta}\right) \left(\frac{2}{\Delta}\right)^{2}\right],$$
(20)

where $s_{m,i} \in \{-1,1\}$ denotes the spin variables of the *m*th world line at the *i*th segment. Basically, the two integrals in (6) are replaced by sums. The spin variables can be summed up to a large spin variable $M_i = \sum_m s_{m,i} \in \{-N, -N+2, ..., N\}$ involving a combinatorial factor

$$Z = \sum_{\{M_i\}} \prod_{i=1}^{L} {N \choose \frac{N+M_i}{2}} \exp\left[\frac{1}{4\beta} \sum_{i,j}^{L} M_i M_j K_\beta \times \left(|j-i|\frac{2}{\Delta}\right) \left(\frac{2}{\Delta}\right)^2\right].$$
(21)

For large N the binomial coefficient tends to a Gaussian and the normalized spin variable $m_i = M_i/N \in [1, -1]$ becomes a continuous variable

$$Z = \sum_{\{M_i\}} A \exp\left[\frac{N}{2} \sum_{i,j}^{L} J_{i,j} m_i m_j - N \sum_{i=1}^{L} \frac{m_i^2}{2}\right]$$
(22)

with

$$\mathbf{A} = \left(\frac{2^{2N+1}}{\pi N}\right)^{\frac{L}{2}}, \quad J_{i,j} = \frac{1}{2\beta} \tilde{K}_{\beta} \left(|j-i|\frac{2}{\Delta}\right) \left(\frac{2}{\Delta}\right)^{2}, \quad (23)$$

and $\tilde{K}_{\beta}(\tau) = NK_{\beta}(\tau)$. The form in (22) shows a onedimensional continuous spin model with long-range interaction, where the states $m_i \approx 0$ are favored due to the quadratic term. One can extract the N dependence from this model, even without having to solve it in detail. Without the last term in (22), the factor N in the exponential would simply scale the coupling. As a consequence, the critical coupling would scale as $\tilde{\alpha}_c \propto N^{-1}$. In the following, we will therefore use temporarily the replacement

$$\tilde{\tilde{\alpha}} = N\tilde{\alpha}.$$
 (24)

Effectively, this replacement is only introduced for bookkeeping the N dependence in the following argumentation.

The effect of the last term in (22) can understood by its influence on the corresponding Ginzburg-Landau functional (GLF) [25,27,28]

$$\mathcal{F}[m(\tau)] = \int d\tau \left[\frac{r}{2} m^2 - \frac{N}{2} m^2 + \frac{u}{4} m^4 + c \int d\tau' \frac{m(\tau)m(\tau')}{|\tau - \tau'|^{1+s}} \right],$$
(25)

where

$$Z = \int \mathcal{D}[m(\tau)] \exp\{\mathcal{F}[m(\tau)]\}.$$

The first and third terms are the common second and fourth order terms of the GLF and the last term represents the longrange interaction in an integral form. The parameters r, u, c are all functions of $s, \Delta, \tilde{\alpha}$, but do not depend on N in an explicit manner. The local term in (22) now enters the functional as another quadratic term $Nm^2/2$ that does not depend on the coupling $\tilde{\alpha}$ or on other system parameters, but depends linearly on the number of spins N. The parameter r can always be shifted in such a way that the phase transition occurs at r = 0. In the case where the $Nm^2/2$ term is neglected, this parameter must then have the form $r = b(s, \Delta)[\tilde{\alpha}_c - a(s, \Delta)]$, with unknown functions a and b. In the full model with included $Nm^2/2$ term, the equation

$$r - N = b(s, \Delta)[\tilde{\tilde{\alpha}}_c - a(s, \Delta)] - N = 0$$
(26)

determines the transition point $\tilde{\alpha}_c$.

Even the offset function $b(s, \Delta)$ can be determined. Since the spin variables in (22) are continuous and independent of N, a saddle point integration becomes exact in the limit $N \to \infty$ and the onset of the magnetization of the most probable state gives the phase transition point for $N \to \infty$:

$$\sum_{j} m_{j} J_{i,j} - m_{i} = 0 \stackrel{m \to 0^{+}}{\Rightarrow} \tilde{\alpha} = \frac{s \Delta}{2\omega_{c}} =: b(s, \Delta)^{-1}.$$
(27)

Instead of the sum on the left-hand side, the integral over the Kernel $\tilde{K}_{\beta}(\tau)/(\beta \Delta)$ is used (cf. Appendix C) and the homogeneity of the spin variables at equilibrium $m = m_k$ is assumed. With (24), (26), and (27), we arrive at the following prediction of the asymptotic N, s, Δ , and ω_c dependence of the critical coupling strength:

$$\alpha_c \simeq \frac{1}{N} \frac{s\Delta}{2\omega_c} + \frac{a(s,\Delta)}{N^2}.$$
(28)

In the large-spin-boson model, the combinatorial factor in (21) is absent but otherwise no change occurs. As a consequence the universality class of the transition remains unchanged but the critical coupling strength is $\alpha_c \propto 1/N^2$ as predicted in [21].

B. The critical point: QMC results

For a sub-Ohmic bath (0 < s < 1) one expects a secondorder phase transition characterized by the following scaling laws (N, s, Δ , and ω_c fixed):

$$m(\alpha = \alpha_c, \beta) \propto \beta^{-x/\nu},$$
 (29)

$$\chi(\alpha = \alpha_c, \beta) \propto \beta^{\gamma/\nu}, \tag{30}$$

$$Q(\alpha,\beta) \sim \tilde{Q}(\beta^{1/\nu}(\alpha-\alpha_c)/\alpha_c).$$
 (31)

At zero temperature (i.e., $\beta \to \infty$), then $m(\alpha > \alpha_c, T = 0) \propto (\alpha - \alpha_c)^x$ and $\chi(\alpha, T = 0) \propto (\alpha_c - \alpha)^{-\gamma}$ holds. The critical exponents *x* (usually denoted as β , the order parameter exponent, which we changed to avoid confusion with the inverse temperature $\beta = 1/T$), ν and γ obey the scaling relation $\gamma = \nu - 2x$ and are expected to be independent of the tunneling strength Δ and the cutoff frequency ω_c , for which reason we fix both to $\Delta = 0.1$ and $\omega_c = 1$ in most calculations. The dependence of γ and ν on *s* is known for N = 1 [3,4], but not for general *N*.

We performed large scale QMC simulations for N = 1, ..., 128 and $\beta = 2000, ..., 1024000$ using a continuous imaginary time algorithm based on the world line representation (see Sec. II A) and [3]). To determine the critical point α_c it is most convenient to use the relation (31): For fixed N, s, Δ , and ω_c the quantity Q is at the critical point $\alpha = \alpha_c$ asymptotically independent of β , which can be used to locate the critical point. This is demonstrated for s = 0.75 and s = 0.9 in Figs. 1 and 2 for N = 2 and N = 16, respectively.

As a check for accuracy we compare the estimates for α_c that we obtain in this way for N = 2 with the predictions of the numerical renormalization group (NRG) calculation of Ref. [12], which is shown in Fig. 3. The agreement is very good in the sub-Ohmic regime for s = 1/2 but differs slightly for an Ohmic spectrum (s = 1). The phase transition of the latter case is known to be notoriously difficult to investigate with Monte Carlo simulations, because it belongs to the Kosterlitz-Thouless universality class. Throughout this paper, we set $\Delta = 0.1$ and $\omega_c = 1$, for which the two independent methods coincide within 5% for the Ohmic case (which is compatible with the error bar) and much less in the sub-Ohmic regime.

Next we determined the critical bath coupling strength for different values of N and s; the result is shown in Fig. 4. For increasing N the rescaled critical bath coupling strength appears to approach an *s*-dependent constant, which means that asymptotically (for large N) α_c would decrease as 1/N. In the last subsection we derived a more precise prediction for

This prediction is checked in Fig. 5. The simulation results

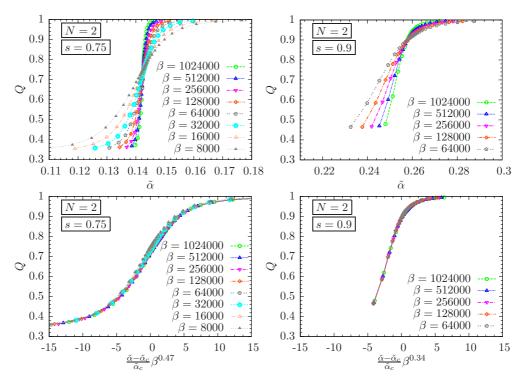


FIG. 1. (Color online) Top row: Fourth-order cumulant Q versus the rescaled bath coupling $\tilde{\alpha}$ for s = 0.75 (left) and s = 0.9 (right) for N = 2. Bottom row: Scaling plots of Q for the data in the top row for s = 0.75 (left) and s = 0.9 (right). The best data collapse is obtained for $1/\nu = 0.47$ for s = 0.75 and $1/\nu = 0.34$ for s = 0.9. The other parameters are $\Delta = 0.1$ and $\omega_c = 1$.

the asymptotic behavior of α_c in Eq. (28):

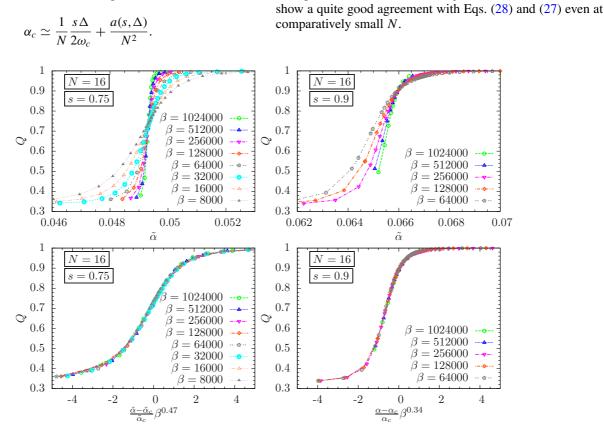


FIG. 2. (Color online) The same as Fig. 1, but for N = 16 instead of N = 2. The scaling plot in the lower row yields the best data collapse for a different critical rescaled bath coupling but the same values for the exponent $1/\nu = 0.47$ for s = 0.75 and $1/\nu = 0.34$ for s = 0.9.

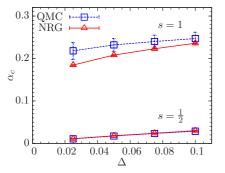


FIG. 3. (Color online) Comparison of QMC and NRG predictions for the critical bath coupling α_c as a function of the tunneling amplitude for N = 2, s = 1/2, and s = 1. Squares are our QMC estimates and triangles the prediction of the NRG calculation of Ref. [12].

According to the theoretical considerations of the last subsection, the universality classes of the MSBM model for general, finite $N < \infty$ should be identical with the universality class of the single spin-boson model. More precisely, one expects a Gaussian fixed point for s < 1/2 (i.e., x = 1/2, $\gamma = 1$, and $\nu = 1/s$; see [24,25]), nontrivial exponents for $1/2 < s \leq 1$, and no phase transition above s = 1. Since the quantum phase transition of the single spin-boson model is well described by the zero-temperature phase transition in the mean-field model (17) (see also [23]), one expects in the regime s < 1/2 mean-field exponents also for the MSBM with N > 1. We confirmed with our QMC simulation that the classical exponents x = 1/2, $\gamma = 1$, and $\nu = 1/s$ for s < 1/2are indeed independent of N (data not shown) and focus here on the more interesting, the nonclassical regime $1/2 < s \leq 1$, and show results for two explicit values: s = 0.75 and s = 0.9.

Figures 1 and 2 show scaling plots of Q according to the scaling relation (31). Since corrections to scaling increase noticeably if *s* approaches 1, we restricted the scaling plot to $\beta \ge 64\,000$ for s = 0.9, whereas for s = 0.75 the data for $\beta \ge 8000$ have been included for scaling. The data collapse is very good for the exponent values $1/\nu = 0.47$ and $1/\nu = 0.34$ for s = 0.75 and s = 0.9, respectively, for both N = 2 and N = 16. These estimates for ν also agree with those

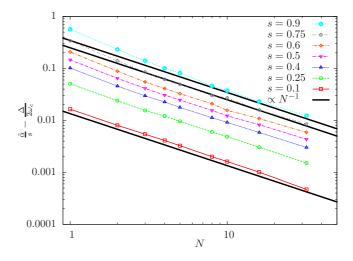


FIG. 5. (Color online) Check of the prediction (28) for the rescaled critical bath coupling $\tilde{\alpha}_c/s - \Delta/(2\omega_c) \propto 1/N$ for several bath exponents $s = 0.1, \ldots, 0.9$ (from bottom to top). The solid lines are a guide for the eyes.

that we obtained for N = 1 (data not shown; cf. [3]) and therefore support our prediction of the last subsection that the universality class of the MSBM is independent of N for finite N.

We obtain a second independent exponent, namely γ , from the behavior of the susceptibility χ (30) at the critical point. The data are shown in Fig. 6 in a log-log plot. For both bath exponents the asymptotic slope of the susceptibility $\chi \simeq \beta^{\gamma/\nu}$ stays unchanged as the number of spins is increased, which confirms that the critical exponent is $\gamma/\nu = 0.78$ for s = 0.75and $\gamma/\nu = 0.975$ for s = 0.9 is universal for any value of *N*.

The pre-asymptotic behavior ($\beta \ll 10^4$) displays small systematic deviations from the straight line. For small N the asymptotic straight line is approached from below indicating corrections to scaling. For large N it is approached from above indicating a crossover from the $N = \infty$ mean-field critical behavior to the asymptotic single spin-boson behavior N = 1.

Finally we study the $N \rightarrow \infty$ limit of the MSBM model, for which we expect the mean-field theory described in Sec. II C

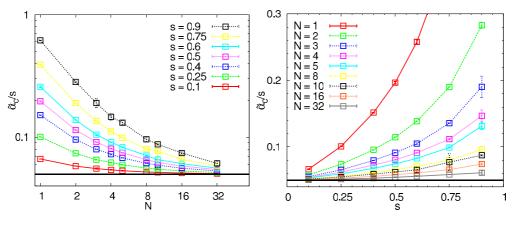


FIG. 4. (Color online) Left: Scaled critical coupling $\tilde{\alpha}_c/s$ versus the number of spins for several bath exponents $s = 0.1, \ldots, 0.9$ (from bottom to top). The solid line indicates the limiting value of $\frac{\Delta}{2\omega_c}$ for $N \to \infty$. Right: $\tilde{\alpha}_c/s$ versus the bath exponent *s*. The parameters are $\Delta = 0.1$ and $\omega_c = 1$.

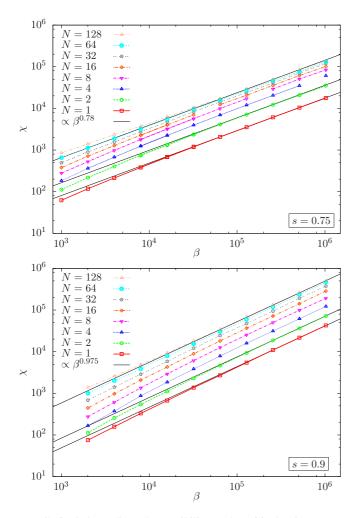


FIG. 6. (Color online) Susceptibility at the critical point versus the inverse temperature β for s = 0.75 (top) and s = 0.9 (bottom) for different *N*. The straight line represents the asymptotic behavior $\chi \propto \beta^{\gamma/\nu}$ with $\gamma/\nu = 0.78$ for s = 0.75 and $\gamma/\nu = 0.975$ for s = 0.9.

to be exact. Since the latter has a phase transition also at nonvanishing temperature (19) we performed extensive QMC simulations for fixed temperatures (fixed β) and $\Delta = 0.1$, $\omega_c = 1$. Figure 7 shows the fourth-order cumulant $Q = \langle m^2 \rangle^2 / \langle m^4 \rangle$ versus the control parameter (i.e., $\tilde{\alpha}$ in our case) for multiple numbers of spins N for two exemplary cases. One

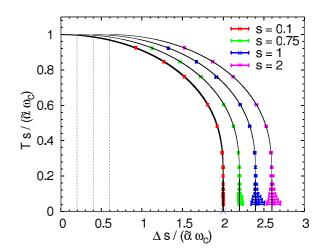


FIG. 8. (Color online) Temperature versus tunneling amplitude Δ (both in unit of $\omega_c \tilde{\alpha}/s$) for different bath exponents *s*. The thick black curve shows the analytic mean-field result (18) and the data points are obtained by QMC in the limit of $N \rightarrow \infty$ (see text). The results for s = 0.75, 1, 2 have been shifted by 0.2, 0.4, and 0.6, respectively (indicated by the dashed lines) for better visibility. The blue cross is the zero-temperature limit (32). The fixed parameters are $\Delta = 0.1$ and $\omega_c = 1$.

can clearly see that the curves intersect in a single point, which determines the phase transition very similarly to the previous zero-temperature analysis, where *N* was held fixed.

The resulting temperature versus tunneling plot (both in units of the scaled coupling $\tilde{J} = \frac{\omega_c \tilde{\alpha}}{s}$) is shown in Fig. 8. All points coincide to a single master curve [cf. Eq. (19)]. One can see an excellent agreement with the expected transition line. All curves tend to the zero-temperature limit of (18)

$$\Delta_c = \frac{2\tilde{\alpha}\omega_c}{s},\tag{32}$$

which is marked as a blue cross in Fig. 8. It should be noted that this transition point exists for arbitrary *s* due to the $N \to \infty$ limit in contrast to the finite *N* case, where a transition only exists for $s \leq 1$. At low temperatures and bath exponents $s \geq 1$, an increase of the size of the statistical error (for constant computational effort) is clearly visible. This is an indication of the onset of the crossover regime between mean-field universality at finite *T* and $N \to \infty$ and

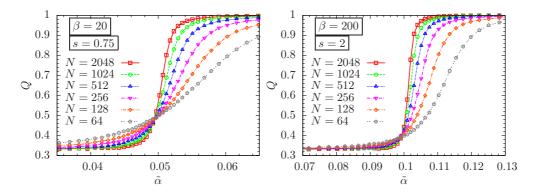


FIG. 7. (Color online) Fourth-order cumulant Q versus the rescaled critical coupling $\tilde{\alpha}$ for several numbers of spin N for the parameters $\beta = 20$ and s = 0.75 (left) and $\beta = 200$ and s = 2 (right).

the nonclassical critical behavior governed by the quantum critical point at T = 0. For the same reason one observes the asymptotic behavior at the quantum critical point for large N only for large β , i.e., low temperatures T.

IV. SPIN-SPIN CORRELATION

In this section, we investigate the effective interaction between the two-level systems mediated by the common bath by calculating the spin-spin correlation functions of the MSBM via QMC. It is defined as

$$\langle \sigma_1^z \sigma_2^z \rangle = \frac{1}{Z} \operatorname{Tr} \left[\sigma_1^z \sigma_2^z \exp(-\beta H) \right],$$
 (33)

which is easily accessible by the quantum Monte Carlo algorithm, since the operator $\sigma_1^z \sigma_2^z$ is diagonal in the used representation.

In Fig. 9 the order parameter $\langle \sigma^z \rangle = \langle \sigma_1^z \rangle = \langle \sigma_2^z \rangle$, the correlation function (33), and the connected correlation function $C_{1,2} = \langle \sigma_1^z \sigma_2^z \rangle - \langle \sigma_1^z \rangle \langle \sigma_2^z \rangle$ are plotted. The vertical lines display the phase transition points α_c with respect to the different bath exponents *s*. As already shown in Sec. III, α_c is shifted towards larger coupling strengths as *s* is increased. At zero temperature ($\beta = \infty$) the order parameter $\langle \sigma^z \rangle$ is zero in the delocalized phase ($\alpha < \alpha_c$) and increases with an algebraic singularity $\propto (\alpha - \alpha_c)^x$ in the localized phase ($\alpha > \alpha_c$). For finite β this sharp transition is smoothed as is visible in Fig. 9(a).

The correlation function $\langle \sigma_1^z \sigma_2^z \rangle$ in Fig. 9(b) is nonzero even in the delocalized phase $\alpha < \alpha_c$ due to the effective ferromagnetic interaction mediated by the common bath. We will quantify this behavior for small coupling strengths at the end of this section. For $\alpha > \alpha_c$ the nonzero order parameter [Fig. 9(a)] superposes with the correlation function. This leads to a kink at α_c that is clearly pronounced if *s* is small, and smeared out if *s* becomes larger.

In Fig. 9(c), the connected correlation function $C_{1,2}$ is shown. This property displays the fluctuation around the mean value. One can see that this function grows until the phase transition point is reached and decreases rapidly to zero after passing it. The larger the bath exponent is, the higher the maximum of the curve is until the phase transition to the localized phase suppresses the fluctuations. For s > 1, where no phase transition takes place, $C_{1,2}$ is monotonically increasing and saturates at $C_{1,2} = 1$ for large α . In this case, the spins are strongly correlated, but not able to perform a localization. In the path-integral representation, the two world lines arrange themselves in a synchronized way, where their kinks occur at nearly equal times, separating identically orientated world line segments from each other.

Figure 10 displays $C_{1,2}$ for a larger range of α , where the behavior for an Ohmic bath spectrum is also covered. Plotted are curves for different inverse temperatures. One observes that even the largest system size with $\beta = 4.096\,000$ still shows strong finite β effects in contrast to systems in the sub-Ohmic regime, where, e.g., for s = 1/2 the curves for $\beta = 32\,000$ and $\beta = 64\,000$ collapse already quite well. The origin of this slow convergence to the infinite β limit can be understood by the particular critical behavior at s = 1. At this point, the

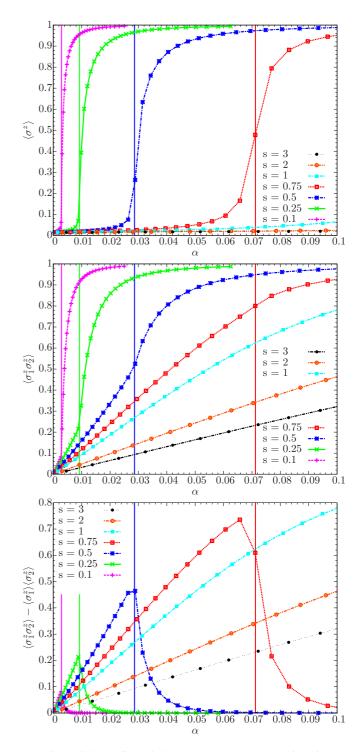


FIG. 9. (Color online) Order parameter $\langle \sigma^z \rangle$ (top), spin-spin correlation function $\langle \sigma_1^z \sigma_2^z \rangle$ (middle), and connected correlation function $C_{1,2} = \langle \sigma_1^z \sigma_2^z \rangle - \langle \sigma_1^z \rangle \langle \sigma_2^z \rangle$ versus the coupling strength α for various bath exponents *s*. The vertical lines correspond to the phase transition points. The inverse temperature is $\beta = 32\,000$.

kernel $K_{\beta}(\tau)$ falls off in imaginary time with an inversesquare law. This kind of long-range interaction is known to produce a Kosterlitz-Thouless transition, where the spin-spin correlation function decays logarithmically with the distance

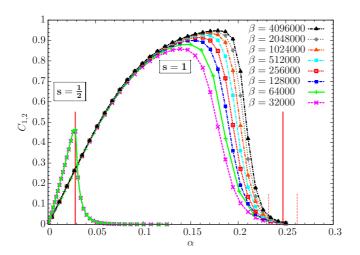


FIG. 10. (Color online) Connected correlation function $C_{1,2}$ versus the coupling strength for the Ohmic case s = 1, where data from $\beta = 32\,000$ to $\beta = 4\,096\,000$ are shown and for $s = \frac{1}{2}$ with $\beta = 32\,000$ and $\beta = 64\,000$. The vertical solid line marks the phase transition point and the dashed lines its error.

in imaginary time [29,30]. Therefore, finite size effects survive over many decades causing much larger deviations in the numerically determined properties. The determination of the phase transition point is performed by an extrapolation of the fourth-order cumulant $Q = \langle m^2 \rangle^2 / \langle m^4 \rangle$ to infinite β and is for s = 1 accompanied by a much larger error bar than for the sub-Ohmic regime (s < 1).

The observed behavior of the connected correlation function is reminiscent of the entanglement entropy

$$\mathcal{E} = -\operatorname{Tr}[\rho_s \log_2 \rho_s],\tag{34}$$

with the reduced density matrix $\rho_s = \text{Tr}_B \rho$, where the bath degrees of freedom were traced out. The entanglement entropy was recently studied in Ref. [12], where Fig. 6 shows also a cusp for the sub-Ohmic case s = 1/2 whereas Fig. 5 therein shows a smooth maximum for the Ohmic spectrum (note that vanishing direct spin-spin coupling K = 0 corresponds to the case considered by us).

We will now discuss some particular regimes of the model. In Ref. [12] it was shown that a polaron transformation $U = \exp[-\frac{1}{2}(\sigma_1^z + \sigma_2^z) \sum_k \frac{\lambda_k}{\omega_k} (b_k^{\dagger} - b_k)]$ applied to the twoimpurity spin-boson model will renormalize a bare (antiferromagnetic) Ising coupling term $\frac{K}{4}\sigma_1^z\sigma_2^z$ to $(\frac{K}{4} - \frac{\alpha\omega_c}{s})\sigma_1^z\sigma_2^z$. Since we consider no bare direct coupling between the spins (K = 0), it is convenient to plot the data against the rescaled coupling strength $J = \alpha\omega_c/s$ to eliminate the global $\propto s^{-1}$ scaling of the bath-induced spin-spin interaction, if one is comparing the behavior for different *s*. In Fig. 11 the data of Fig. 9(b) are plotted versus *J*. By means of this plot, one can see that the curves for high s > 1 tend to a master curve. This master curve corresponds to the analytic result of two nondissipative and Ising-coupled spins driven by a transverse field

$$H = \frac{\Delta}{2} \left(\sigma_1^x + \sigma_2^x \right) - J \sigma_1^z \sigma_2^z.$$
(35)

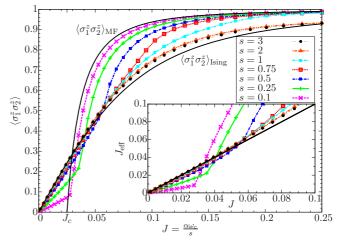


FIG. 11. (Color online) Correlation function of two spins in a common bath versus the rescaled coupling parameter $J = \alpha \omega_c / s$ for various bath exponents *s*. The solid lines are the analytic results of the pure Ising coupling or the mean-field interaction, respectively (see text). The inset shows the effective direct interaction J_{eff} versus *J* for small coupling strengths ($\alpha = 0...0.1$). The solid line is the identity $J_{\text{eff}} = J$.

This toy model is exactly solvable and one could calculate the spin-spin correlation to be

$$\left\langle \sigma_1^z \sigma_2^z \right\rangle_{\text{Ising}} = \frac{J}{\sqrt{J^2 + \Delta^2}} \tag{36}$$

in the zero-temperature limit.

For $s \rightarrow 0$ the data are compatible with mean-field behavior as can be seen by solving the MSBM in the mean-field approximation [cf. Eq. (17)] for finite N. Within this approximation, the model shows a phase transition at $J_c = \Delta/(2N)$ as discussed in Sec. II C, and the correlation function, which is in this case simply the order parameter squared, reads

$$\left\langle \sigma_1^z \sigma_2^z \right\rangle_{\rm MF} = \left\langle \sigma^z \right\rangle^2 = \begin{cases} 0 & \text{if } J \leqslant J_c = \frac{\Delta}{2N}, \\ 1 - \frac{\Delta^2}{4N^2 J^2} & \text{if } J > J_c. \end{cases}$$
(37)

The data for $\langle \sigma_1^z \sigma_2^z \rangle$ for the lowest bath exponent s = 0.1 are quite close to the mean-field results (37) in the localized phase. In the delocalized phase $\langle \sigma_1^z \sigma_2^z \rangle$ does not vanish, in contrast to mean-field behavior, and also deviates from the behavior of the two coupled spins without bath (36). To shed light on this weak-coupling regime, we recast Eq. (36) and define an effective direct coupling

$$J_{\rm eff} = \frac{\Delta \left\langle \sigma_1^z \sigma_2^z \right\rangle}{\sqrt{1 - \left\langle \sigma_1^z \sigma_2^z \right\rangle^2}},\tag{38}$$

which is a measure for the effective ferromagnetic coupling mediated by the common bath.

The inset of Fig. 11 shows the effective direct coupling (38) versus the rescaled coupling $J = \alpha \omega_c/s$. Again, one sees the approach of the curves for high *s* to the identity $(J_{\text{eff}} = J)$, whereas for small *s* the effective direct coupling develops significantly slower than *J*, but still linear. The data imply that the ratio J_{eff}/J decreases to zero for $s \rightarrow 0$, which is compatible with the mean-field approximation, since the

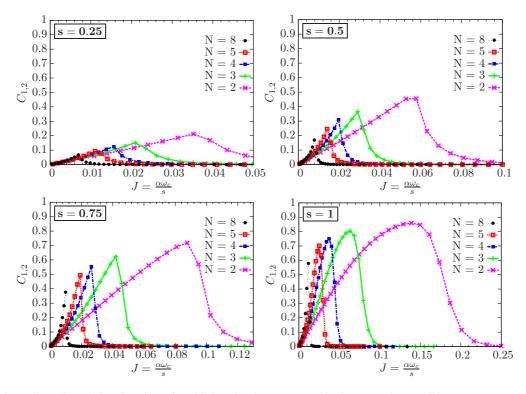


FIG. 12. (Color online) Correlation function of multiple spins in a common bath versus the coupling parameter $J = \alpha \omega_c / s$ for bath exponents s = 0.25, 0.5, 0.75 and s = 1 at an inverse temperature of $\beta = 32\,000$.

results (37) predict $J_{\text{eff}} = 0$ in the delocalized phase. For s > 1 the ratio J_{eff}/J approaches 1; the effective interaction mediated by the common bath is just $J_{\text{eff}} = J = \alpha \omega_c/s$. Note that even for $s \to \infty$, i.e., short-ranged interactions in the imaginary time direction, the two-level systems are still ferromagnetically correlated due to the effective bath interaction at equal imaginary times. The $s \to \infty$ limit of the MSBM for N = 2 is thus simply the pair of two-level systems without bath but with a ferromagnetic coupling $J = \alpha \omega_c/s$.

Finally, we had a look at the correlations for more than two spins. Figure 12 shows the connected correlation function for N = 2 to N = 8 for different bath exponents. One can see that the maximum becomes lower and more sharp as the number of spins is increased. For larger N the maximum takes place at the mean-field prediction of $J_c = \Delta/(2N)$ (37). The initial slope for $J \rightarrow 0$ is quite similar for all curves, but raises super-linearly for N > 2 instead of linearly for N = 2, when the coupling is increased. That means that the spins build up their cooperative polarization successively.

V. SPATIAL SEPARATION

In this section we consider spatially separated two-level systems in a common bath. As originally proposed in Ref. [6] the two-level systems are now coupled to a phase-shifted polarization of the bosonic modes. The phase shift arises from the time t = R/v that is needed to propagate the information from one spin to another spin at a distance R with the propagation velocity v. Such a situation arises for instance if the bosonic modes are represented by standing waves in a box with a linear size L and periodic boundary conditions (Fig. 13). In this case, the bath is fully described by a complete set of harmonic modes

 $\mathbf{k} = \mathbf{n} \cdot 2\pi/L$ with $\mathbf{n} = (n_1, \dots, n_d)$, and $n_i = 1, 2, 3, \dots$ for $i = 1, \dots, d$, and *d* the dimension of the box. After performing $L \to \infty$, this approach is basically a continuous version of phonon-induced interaction between localized electrons [32]. The same distance-dependent interaction has also been used for two harmonic oscillators in a common bath instead of two spins [33].

A. The dissipative spin array

The generalization of interaction term of the MSBM (4) to a system where the N spins are located at the positions $\mathbf{r_m}$ is given by [13,31]

$$H_I = \sum_{m=1}^{N} \frac{1}{2} \sum_{\mathbf{k}} \lambda_{\mathbf{k}} (a_{\mathbf{k}}^{\dagger} e^{i\mathbf{k}\cdot\mathbf{r}_{\mathbf{m}}} + a_{\mathbf{k}} e^{-i\mathbf{k}\cdot\mathbf{r}_{\mathbf{m}}}) \sigma_m^z.$$
(39)

Without loss of generality, the coefficients λ_k are assumed to be real numbers, because a complex phase carries no information. In contrast to the previous model without any relation to spatial variables, the indices have now the meaning of actual wave numbers rather than a simple counter for the frequencies. In principle, this would involve a dispersion

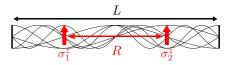


FIG. 13. (Color online) Sketch of two spins at a distance R embedded in a one-dimensional common bosonic bath. The bath is fully described by a complete set of harmonic modes, that are restricted to a box with linear size L.

relation $\omega \to \omega(\mathbf{k})$, which would also affect the couplings $\lambda_k \to \lambda_{\omega(\mathbf{k})}$. Equation (39) describes hence a bath of modes propagating in vacuum, where $\omega \propto |\mathbf{k}|$.

For a pair of spins at a distance $R = |\mathbf{R}| = |\mathbf{r} - \mathbf{r}'|$, one can show (cf. Appendix A) that the dependence on **R** can be completely integrated into the spectral function

$$J(\omega, \mathbf{R}) = \pi \sum_{\mathbf{k}>0} \lambda_{\mathbf{k}}^2 \cos\left(\mathbf{k} \cdot \mathbf{R}\right) \delta(\omega - \omega_{\mathbf{k}})$$
(40)

such that the form of the kernel (7) remains unchanged. The *N* spin problem is just a superposition of all of these pair kernels for every distinct spin pair. The sum in (40) can be replaced by an integral in the limit $L \to \infty$ and the couplings are assumed to follow a power-law function $\lambda_{k_i} \propto k_i^{(2+s-d)/2}$ (i = 1, ..., d).

Absorbing all constants into the bath coupling strength α leads to the form

$$J(\omega, R) = 2\pi \alpha \omega^s \omega_c^{1-s} f^d(\bar{R}), \qquad (41)$$

where we have introduced the scaled distance $\bar{R} = R\omega_c/v$. The dependence on the spatial separation is thus absorbed into a single function $f^d(\bar{R})$. For the different dimensions d = 1,2,3 of the bath, this function is

$$f^{d}(\bar{R}) = \begin{cases} \cos(\bar{R}) & d = 1, \\ J_{0}(\bar{R}) & d = 2, \\ \sin(\bar{R})/\bar{R} & d = 3, \end{cases}$$
(42)

where $J_0(\bar{R})$ is the Bessel function of the first kind.

Let us focus on the separation-dependent part and consider the zero-tunneling limit $\Delta = 0$. The approach of Sec. II B can easily be generalized and the resulting partition function corresponds to the one of a long-range Ising model

$$Z = \sum_{\{\sigma_1,...,\sigma_N\}} \exp\left[\frac{\beta}{2} \sum_{m=1}^N \sum_{m'=1}^N \sigma_m \sigma_{m'} \frac{\alpha \omega_c}{s} X^d \times \left(s, \frac{\omega_c}{v} |\mathbf{r}_m - \mathbf{r}_{m'}|\right)\right],$$
(43)

where the Ising interaction is defined as

$$X^{d}(s,\bar{R}) = \int_{0}^{1} sx^{s-1} f^{d}(x\bar{R}) dx.$$
 (44)

For every dimension d = 1,2,3, these functions reproduce the MSBM case from Sec. II B, if the scaled distance $\bar{R} \to 0$. In Fig. 14, $X^d(s, \bar{R})$ is plotted for different *s* in one, two, and three dimensions. For small *s* they decay monotonically, whereas for higher *s* oscillations become dominant. The asymptotic behavior of the former case can be determined by recasting (44) to $X^d(s, \bar{R}) = \bar{R}^{-s} I^d(s, \bar{R})$ and noting that the remaining integral $I^d(s, \bar{R}) = \int_0^{\bar{R}} sx^{s-1} f^d(x) dx$ converges to a finite (nonzero) value as $\bar{R} \to \infty$. It turns out that

$$X^d(s,\bar{R}) \simeq \bar{R}^{-s} \tag{45}$$

for 0 < s < (d+1)/2.

Above this value s > (d + 1)/2 this analysis breaks down and one should rather focus on the envelop function of $X^d(s, \bar{R})$. One way is to use the form (e.g., in d = 1) $X^1(s, \bar{R}) =$

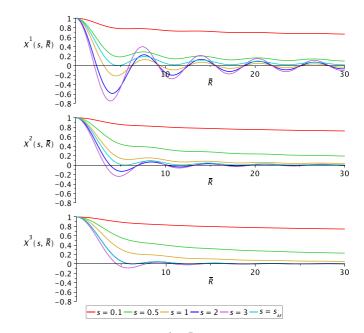


FIG. 14. (Color online) $X^d(s, \overline{R})$ in one, two, and three dimensions for various *s*. For d = 3, the curve for s = 2 coincides with one for $s = s_{AF}$.

 $\bar{R}^{-1} \int_{0}^{\bar{R}} s(x/\bar{R})^{s-1} \cos(x) dx$ and analyze the asymptotic behavior of the maxima at $\bar{R}_n = 2\pi n + \pi/2$ for large *n*. One obtains the *s*-independent power-law decay

$$X^{d}(s,\bar{R}) \simeq \bar{R}^{-(d+1)/2},$$
 (46)

for the envelop function, if s > (d + 1)/2.

In addition to the asymptotic behavior another upper bound for the bath exponent is important that also depends on the dimension of the bath. We define s_{AF}^d such that for all $s > s_{AF}^d$ the coupling interaction $X^d(s, \bar{R})$ is not restricted to positive values for all distances any more. Intervals of the distance \bar{R} occur, for which antiferromagnetic interactions occur. The onsets of antiferromagnetism are in this sense $s_{AF}^1 = 0.6923(9)$, $s_{AF}^2 = 1.3545(2)$, and $s_{AF}^3 = 2.001(2)$. For bath exponents smaller than these values the spin-spin interactions are always ferromagnetic and decay asymptotically as discussed above.

The lower bound for feasible values of the bath exponents follows from the following consideration: It is known that a *D*-dimensional Ising model with algebraic decaying long-range interaction does only have a nondiverging energy per spin, if the decay is faster than $1/R^D$ (see Ref. [25] and references therein). That means for example that in the present case the one-dimensional infinite spin chain in a *d*-dimensional bosonic bath can only exist if $s \ge 1$.

These considerations are summarized in Fig. 15. For $1 \leq s \leq s_{AF}^d$, the system has strict ferromagnetic interactions and is equivalent to the classical Ising spin chain with algebraically decaying long-ranged interactions [25]. This regime exists in the range of bath exponents 1 < s < 2 in the case of a three-dimensional bath and vanishes for d = 1. For $s \geq s_{AF}$ the behavior depends on the details of the lattice spacings $|\mathbf{r}_m - \mathbf{r}_{m+1}|$ between the spins. Two possible cases are for instance when the spins are either arranged near the maxima of the underlying oscillation of $X^d(s, \bar{R})$ or, on the other hand,

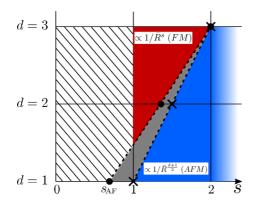


FIG. 15. (Color online) Sketch of the phases of the dissipative spin array in the zero-tunneling limit for a one-, two-, and threedimensional bath. White lined: A decay slower than $\propto R^{-1}$ is thermodynamically not possible in an infinite chain. Red: Ferromagnetic interaction (FM) with $\propto R^{-s}$ decay. Blue: Antiferromagnetic interaction (AFM) with *s*-independent $\propto R^{-\frac{d+1}{2}}$ decay. Gray: Intermediate regime between FM and AFM.

when ferromagnetic and antiferromagnetic contributions are mostly canceling each other resulting in alternating signs of the interactions.

B. A pair of spatially separated spins

In this section we study the MSBM for N = 2 with spatially separated spins in a distance \overline{R} . For a pair of spins, the action (6) separates into four parts $S = S_1 + S_2 + 2S_R$, namely the self-energy for each spin

$$S_m = \int_0^\beta \int_0^\tau \sigma_m^z(\tau) \sigma_m^z(\tau') K_\beta(\tau - \tau', 0) d\tau' d\tau \qquad (47)$$

for m = 1, 2, and twice the interaction between the two spins

$$S_R = \int_0^\beta \int_0^\tau \sigma_1^z(\tau) \sigma_2^z(\tau') K_\beta(\tau - \tau', \bar{R}) d\tau' d\tau.$$
(48)

The Monte Carlo cluster algorithm (see Appendix B) can be adapted to the present case by taking all of these actions into account. The key step is to evaluate the integral kernel $K_{\beta}(\tau, R)$ (7) with the spectral function (41), which is modified by the spatial function $f^d(\bar{R})$. The kernel must be nonnegative in order to apply the algorithm, which is designed for ferromagnetic interaction only. The condition for that can be obtained by calculating the values of *s*, for which $K_{\beta}(\tau, R)$ begins to exhibit roots with respect to τ . At these onset values of *s*, the roots are at $\tau = 0$ and $\tau = \beta$. In the limit $\beta \to \infty$, the values of kernels at these points are

$$K_{\beta}(0,R) = K_{\beta}(\beta,R) = \int_{0}^{\omega_{c}} 2\alpha \omega^{s-1} \omega_{c}^{1-s} f^{d} \left(\frac{\omega R}{v}\right) \frac{\beta}{2} d\omega$$
$$= \frac{2\omega_{c}^{2}\alpha}{s+1} X^{d}(s+1,\bar{R}) \frac{\beta}{2}.$$
(49)

The function $X^d(s, \bar{R})$ exhibits roots at $s \ge s_{AF}^d$ (Sec. V A). This means that the kernel is nonnegative only for bath exponents $s < s_{AF}^d - 1$. If we restrict ourselves to the case of a three-dimensional bath, we can safely explore the complete (sub-)Ohmic regime ($0 < s \le 1$). The integral kernel is plotted

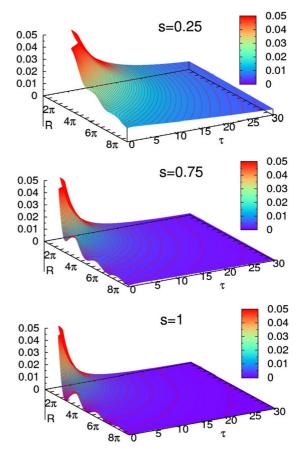


FIG. 16. (Color online) Integral kernel $K_{\beta}(\tau, \bar{R})$ at d = 3 for low imaginary time $0 \le \tau \le 30$ and scaled distances $0 \le \bar{R} \le 8\pi$ for different bath exponents [s = 0.25 (top), s = 0.75 (middle), and s = 1 (bottom)].

in Fig. 16 for several parameters for d = 3. At the s = 1, one can see the roots at $\bar{R} = 2\pi, 4\pi, 6\pi, \dots$

1. Phase transition point

As already known from [12,13], the transition point from delocalization to localization of two spins in a common bath is lowered in comparison to the single spin case. If one introduces a finite spatial separation, the transition point increases up to the single spin case, if $\bar{R} = \infty$. We determined the phase diagram as in Sec. III B via finite β scaling of the fourth-order cumulant Q. Figure 17 shows the phase diagram for various distances. One can see that the spins are influencing each other over relatively large distances. Even at $\bar{R} = 1000$, the transition point is still clearly distinguishable from the single spin case. The localization is strongly enhanced by the presence of a second impurity in the bath. Note that the model neglects retardation effects, which are more important the further the spins are separated. At this point retardation effects mean all corrections that are caused by the finiteness of the propagation speed of the modes and change the effective interaction between spins. Therefore, the bath-mediated spin-spin influence at large distances may be less prominent in a real experimental setup.

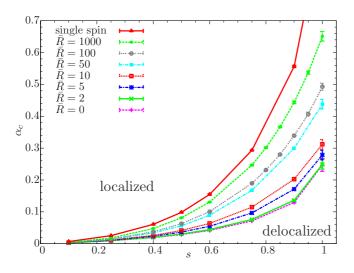


FIG. 17. (Color online) Phase diagram of the MSBM for N = 2 with spatial separation for various distances \bar{R} . The case $\bar{R} = 0$ corresponds to the MSBM with N = 2 from Sec. II and the case $\bar{R} = \infty$ represents the common single spin-boson model.

2. Order parameter and spin-spin correlation

We will now focus on the observables. First we will address extrapolation to zero temperature. In Fig. 18 the order parameter $\langle \sigma^z \rangle$, the fourth-order cumulant Q, and the correlation function $\langle \sigma_1^z \sigma_2^z \rangle$ versus the distance \bar{R} are shown for several inverse temperatures and for two different points in the parameter space, namely $s = 0.75, \alpha = 0.1$ and s = $1, \alpha = 0.3$. There are two things to realize: The fourth-order cumulants intersect for s = 0.75 and merge with each other smoothly for s = 1, which is a characteristic that the former case is an ordinary continuous phase transition whereas the latter one is expected to be a phase transition of a Kosterlitz-Thouless type. The second things is that the correlation function is nearly not affected by the finite size.

In Fig. 19 we plot the order parameter and the correlation function versus the distance for several coupling strengths and for the two selected parameters s = 0.75 and s = 1. Basically the observables behave typically when driven through the phase transition by increasing the control parameter \bar{R} . As a special feature, the curves show an underlying wiggling that is

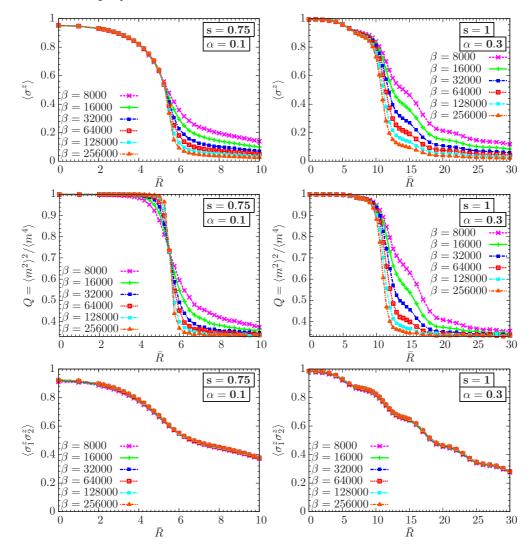


FIG. 18. (Color online) Order parameter $\langle \sigma^z \rangle$ (top), fourth-order cumulant Q (middle), and the correlation function $\langle \sigma_1^z \sigma_2^z \rangle$ (bottom) versus the distance \bar{R} for different inverse temperatures. In the left column the parameters s = 0.75 and $\alpha = 0.1$ are used whereas in the right column s = 1 and $\alpha = 0.3$ are assumed.

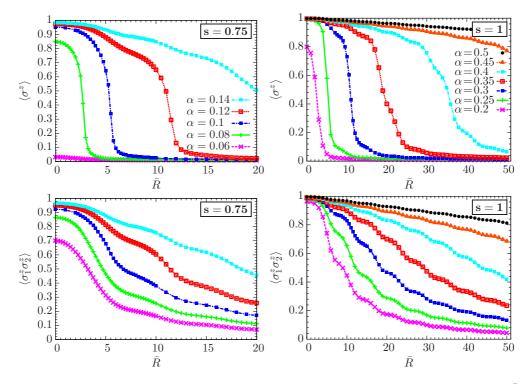


FIG. 19. (Color online) Order parameter $\langle \sigma^z \rangle$ (top) and the correlation function $\langle \sigma_1^z \sigma_2^z \rangle$ (bottom) versus the distance \bar{R} for the parameter s = 0.75 (left) and the Ohmic spectrum s = 1 (right). The inverse temperature for all data is $\beta = 256\,000$. The coupling strengths used in the lower row correspond to values of the upper row.

caused by the oscillations of the spatially dependent interaction (cf. Sec. V A). Also worth mentioning is the observation that the phase transition is not really recognizable by the correlation function curves despite the significant jump of the order parameter. This fact is however also present without spatial separation as one can see Fig. 9(b) for the same s.

VI. CONCLUSIONS

With the help of analytic considerations and large-scale QMC simulations we have shown that the quantum phase transition in the (unbiased) multi-spin-boson model, describing N independent two-level systems coupled to a common bosonic bath, falls in the universality class of the single spin-boson model, i.e., is independent of the number of two-level systems for any finite N. In the limit $N \rightarrow \infty$ the universality class of the transition changes to mean-field behavior and the system has a phase transition also at finite temperatures. Consequently for large values of N one observes a crossover from mean-field to finite N MSBM critical behavior for decreasing temperature.

The leading *N* dependence of the critical bath coupling is shown to be $\alpha_c \simeq \frac{1}{N} \frac{s\Delta}{2\omega_c} + \frac{a(s,\Delta)}{N^2}$; i.e., asymptotically the critical bath coupling is proportional to 1/N. This confirms that the common bath mediates a ferromagnetic interaction between the two-level systems, reduces fluctuations and thus also the critical bath coupling. We showed that the effective interaction between the two-level systems can be quantified by the spin-spin correlation function, which interpolates smoothly between mean-field behavior for $s \to 0$ (including a transition from uncorrelated to correlated two-level systems) and a simple ferromagnetically coupled two-spin system without bath for $s \gg 1$. For $s \gg 1$ the short-ranged interactions in imaginary time still mediate an instantaneous ferromagnetic interaction of strength $J = \alpha \omega_c / s$. For s < 1 the behavior of the connected spin-spin correlation function is reminiscent of the entanglement entropy of the two-level systems: approximately linearly increasing with the bath coupling α in the delocalized phase $\alpha < \alpha_c$ and then rapidly decreasing in the localized phase ($\alpha > \alpha_c$).

For spatially separated two-level systems the critical bath coupling monotonically increases with the distance *R* from the value for the MSBM without spatial separation for R = 0 to the value for the single spin-boson model for infinite separation $R \rightarrow \infty$. No indication is found for an alteration of the universality class of the transition by the spatial separation. The order parameter as well as the spin-spin correlation function decreases systematically with the distance *R* but shows superimposed oscillations.

The results reported in this paper are for the MSBM without a direct ferromagnetic interaction between the twolevel systems, and one obvious question arises immediately: Will the universality class of the transition change if a direct coupling between the two-level systems is introduced? NRG results for a pair of spins (N = 2) reported in [12] are confined to the transition triggered by the ferromagnetic coupling for fixed bath coupling and predict exponents (e.g., $\beta = 0.09$ for s = 0.9 and $\beta = 0.2$ for s = 0.75), which cannot be compared directly with the bath coupling triggered critical behavior of the MSBM (for which we found $\beta = 0.037$ for s = 0.9 and $\beta = 0.23$ for s = 0.75). An alternative view was obtained recently [14] with a renormalization group calculation for a pair of spins each with its own bath, which might behave differently than the case we are considering here. Note that the quantum critical point of a chain of ferromagnetically coupled spins, each with its own bath, was studied in the limit $N \to \infty$ in Ref. [15]. Again the question is whether the reported universality class changes for a common bath. Certainly it will be worthwhile to study these questions with the method that we used in this paper.

ACKNOWLEDGMENT

This work was financially supported by the German Research Foundation DFG within GRK 1276.

APPENDIX A: DERIVATION OF THE ACTION

We are using the Caldeira-Leggett formulation [2] for quantum dissipative systems, where the 1/2-spins are represented by spinless particles at positions Q_m and momentum P_m in a symmetric double-well potential $V(Q_m)$. The full Hamiltonian of N spatially separated spins in a common bosonic bath can written as

 $H = H_S + H_B + H_I + H_c$

(A1)

$$H_{S} = \sum_{m=1}^{N} \left(\frac{P_{m}^{2}}{2M} + V(Q_{m}) \right),$$
(A2)

$$H_B = \sum_{\mathbf{k}=-\infty}^{\infty} \frac{p_{\mathbf{k}}^2}{2m_{\mathbf{k}}} + \sum_{\mathbf{k}=-\infty}^{\infty} \frac{m_{\mathbf{k}}\omega_{\mathbf{k}}^2}{2} x_{\mathbf{k}}^2, \qquad (A3)$$

$$H_I = \sum_{\mathbf{k}=-\infty}^{\infty} c_{\mathbf{k}} x_{\mathbf{k}} \sum_{m=1}^{N} Q_m \exp(i\mathbf{k} \cdot \mathbf{r}_m), \qquad (A4)$$

$$H_{C} = \sum_{\mathbf{k}=-\infty}^{\infty} \frac{|c_{\mathbf{k}}|^{2}}{2m_{\mathbf{k}}\omega_{\mathbf{k}}^{2}} \sum_{m=1}^{N} \sum_{m'=1}^{N} Q_{m} Q_{m'} \exp[i\mathbf{k} \cdot (\mathbf{r}_{m} - \mathbf{r}_{m'})],$$
(A5)

where H_S is the (inner) system part, H_B denotes the bath degrees of freedom, H_I is the interaction term, and H_C the so-called counterterm, that compensates a global energy shift. The linear coupling between the particles and the bath mode is imposed by the coefficients c_k . The common procedure is to express the partition function as a path integral [1]

$$Z \propto \int \prod_{m} \mathcal{D}[\mathcal{Q}_{m}] \mathcal{D}[P_{m}] \prod_{\mathbf{k}} \mathcal{D}[x_{\mathbf{k}}] \mathcal{D}[p_{\mathbf{k}}] \exp(-S) \quad (A6)$$

and switch to Fourier space $Q_m(\tau) = \sum_n Q_{m,n} \exp(i\omega_n \tau)$ $(P_{m,n}, p_{\mathbf{k},n}, \text{ and } x_{\mathbf{k},n} \text{ similarly})$, where $\omega_n = 2\pi n/\beta$ are the Matsubara frequencies and τ the imaginary time coordinate. The bath degrees of freedom and the particle momenta can now be eliminated by Gaussian integration. The resulting interacting (nonlocal) part of the action is then given by

$$S_{I} = \beta \sum_{m,m'=1}^{N} \left[\sum_{n=-\infty}^{\infty} \alpha(n, \mathbf{r}_{m} - \mathbf{r}_{m'}) Q_{m,n} Q_{m',-n} \right]$$
(A7)

with the kernel

$$\alpha(n,\mathbf{R}) = \sum_{\mathbf{k}} \frac{|c_{\mathbf{k}}|^2 \omega_n^2 \exp(-i\mathbf{k}\cdot\mathbf{R})}{2m_{\mathbf{k}}\omega_{\mathbf{k}}^2 (\omega_n^2 + \omega_{\mathbf{k}}^2)}.$$
 (A8)

Transforming back by means of the identity [34]

$$\sum_{n} \frac{\omega_k^2}{(\omega_n)^2 + \omega_k^2} e^{i\omega_n(\tau - \tau')} = \frac{\beta}{2} \omega_k \frac{\cosh\left[\frac{\beta\omega_k}{2}\left(1 - \frac{2|\tau - \tau'|}{\beta}\right)\right]}{\sinh\left(\frac{\beta}{2}\omega_k\right)}$$
(A9)

leads to the representation of Eqs. (6), (7), (40) with $|\lambda_k|^2 = |c_k|^2/(2m_k\omega_k)$. The (local) system part (A2) becomes in the spin-boson limit a two-state system whose path-integral representation is a Poissonian distribution of an even number of kinks [36–38].

APPENDIX B: THE CONTINUOUS TIME CLUSTER QMC ALGORITHM

In this appendix we describe the essential steps of the QMC procedure we are using. To demonstrate the principle of the algorithm we start with a discretized (Trotter) representation of the action

$$S = -\frac{1}{2} \sum_{i,j=1}^{L'} J_{|i-j|} s_i s_j - K \sum_{i=1}^{L'} s_i s_{i+1}$$
(B1)

with a ferromagnetic long-range and a nearest-neighbor part

$$J_{i} = \frac{\Delta \tau^{2}}{2\beta} K_{\beta}(i\Delta\tau) \quad \text{and} \quad K = -\frac{1}{2} \log \left[\tanh\left(\frac{\Delta}{2} \cdot \Delta\tau\right) \right]$$
(B2)

and perform the continuous time limit afterwards. Within this discretization, the world line is divided into spins with infinitesimal lattice spacing $\Delta \tau = \beta/L' \rightarrow 0$ in contrast to the approximative discretization of Sec. III A which would correspond to $\Delta \tau = 2/\Delta$. Cluster Monte Carlo methods for classical Ising spins models involve a cluster building procedure, that adds equally aligned (up or down) spins s_i and s_j with a bond activation probability (BAP)

$$p_{|i-j|} = 1 - \exp(-2J_{|i-j|} - 2K\delta_{1,|i-j|})$$
(B3)

to a cluster [35]. The two parts of the interaction can be treated separately in the cluster building process since the separation of long-range and nearest-neighbor BAPs

$$p_i^{(lr)} = 1 - \exp(-2J_i),$$
 (B4)

$$p_i^{(nn)} = 1 - \exp(-2K\delta_{1,i})$$
 (B5)

yields the correct global BAP (B3) via

$$p_i = p_i^{(lr)} + p_i^{(nn)} - p_i^{(lr)} p_i^{(nn)}$$
(B6)

$$= 1 - (1 - p_i^{(\text{lr})})(1 - p_i^{(\text{nn})}).$$
 (B7)

We will now focus on the long-range interaction and will describe how the cluster building process can be performed quickly [25]. The probability for activating a bond to the kth

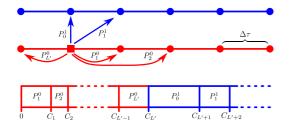


FIG. 20. (Color online) Sketch of the arrangement of the skipping probabilities P_k^n and the array of the cumulative probabilities C_k for two world lines. The neighbors are counted to the right meaning that the left neighbor of the same world line has the label L'. In this example, the reference spin to which the bonds shall be connected is depicted by the square. The skipping probabilities to the spins of the second world line are concatenated with the cumulative probability array. The draw of a random number $C_{L'+1} < g \leq C_{L'+2}$ would correspond to the selection of the nearest-neighbor spin to the right of the *other* world line.

neighbor without making any bond to neighbors in between is

$$P_{k} = p_{k}^{(\mathrm{lr})} \prod_{i=1}^{k-1} \left(1 - p_{i}^{(\mathrm{lr})}\right)$$
(B8)

and will be called skipping probability in the following. Note that only the equal alignment of the reference spin and its *k*th neighbor is required to add the candidate s_k to the cluster, so that the spins s_1, \ldots, s_{k-1} can be skipped without being touched. The cumulative probability

$$C_l = \sum_{k=1}^{l} P_k = 1 - \exp\left(-2\sum_{i=1}^{l} J_i\right)$$
 (B9)

enables one to draw directly (without rejections) the next candidate l for making a bond by a formal inversion

$$l = C^{-1}(g),$$
 (B10)

where g = ran(0, C(L')) is a random number uniformly distributed between 0 and C(L'). This rejection-free procedure can build up for arbitrary ordered probabilities. If we label the different world lines in a way that we can address them by a fixed and ordered sequence, every world line possesses its well-defined neighbor world lines. The extended skipping probability for activating a bond to the *l*th neighbor spin but of the *n*th neighbor world line where all spins of all world lines in between have been skipped can defined by (cf. Fig. 20)

$$P_k^n = \left(\prod_{m=0}^{n-1} \prod_{i=1}^{L'} \left(1 - p_i^{(\mathrm{lr}),m}\right)\right) p_k^{(\mathrm{lr}),n} \prod_{i=1}^{k-1} \left(1 - p_i^{(\mathrm{lr}),n}\right).$$
(B11)

Note that for n = 0 the first product is empty and (B8) is recovered. We can now append these probabilities (B11) to the cumulative probability

$$C_{n \cdot L'+l} = \sum_{m=0}^{n} \sum_{k=1}^{l} P_k^m$$
(B12)

and process as in the single world line case. For later purposes a cumulative probability for groups of spin has to be defined. Consider a certain number r with the labels s_{-r+1}, \ldots, s_0 of adjacent and equally orientated spins. The probability for activating at least one bond between these spins and another one s_i (i > 0) is

$$p_i^{(\mathrm{lr})}(r) = 1 - \prod_{j=0}^{r-1} \left(1 - p_{i+j}^{(\mathrm{lr})} \right) = 1 - \exp\left(-2\sum_{j=0}^{r-1} J_{i+j} \right).$$
(B13)

Based on this probability modified skipping probabilities and cumulative probabilities can be defined in the same way as described above leading to

$$C_{l,r} = 1 - \exp\left(-2\sum_{i=1}^{l}\sum_{j=0}^{r-1}J_{i+j}\right).$$
 (B14)

We will now perform the continuous time limit $\Delta \tau \rightarrow 0$, which means that a sequence of equally aligned spins $s_i, s_{i+1}, \ldots, s_j$ is represented as a continuous segment [s(i), s(j)] of length t = |j - i|. The nearest-neighbor BAP becomes $p_i^{nn} = 1 - (\Delta/2)\Delta\tau$ in the $\Delta\tau \rightarrow 0$ limit and the probability for adding *all* equally aligned spins up to a distance *t* is

$$p^{\text{seg}}(t) = \lim_{\Delta \tau \to 0} \left(1 - \frac{\Delta}{2} \Delta \tau \right)^{t/\Delta \tau} = \exp\left(-\frac{\Delta}{2} t \right). \quad (B15)$$

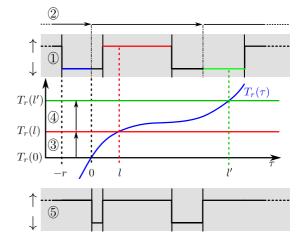


FIG. 21. (Color online) An exemplary part with four kinks of an arbitrary world line is shown in step (1). In step (2) exponentially distributed random numbers according to (B15) are drawn and sequentially inserted into the world line (two dash-dotted lines). These "cuts" together with the already existing kinks divide the world line into segments. Every segment is able to initiate a cluster building process. Without loss of generality the first (\downarrow) segment to the left (blue) is chosen and its end point defines the origin of coordinates. Drawing an exponentially distributed random number in step (3) gives the value $T_r(l) = T_r(0) - \log [ran(0,1)] [cf. (B18)].$ Determining numerically the point l and its corresponding segment (red) leads to the next candidate for adding to the cluster, which will be rejected in the present case due to its incompatible (\uparrow) orientation. Proceeding likewise with l as the starting point leads to another candidate (green) in step (4), which is this time actually added to the cluster. Finally, if this cluster is flipped and the redundant "cuts" are removed the world line would look like in step (5).

The cumulative long-range probability can be expressed by inserting (B2) in (B14) and reads

$$C(l,r) = 1 - \exp\left(-\frac{1}{\beta} \int_{0}^{l} d\tau \int_{-r}^{0} d\tau' K_{\beta}(\tau - \tau')\right).$$
(B16)

This is the cumulative distribution probability for activating a bond between a reference segment at [-r,0] and the segment which is located at a distance *l*. Note that for multiple world lines $K_{\beta}(\tau - \tau')$ has the meaning of a concatenated function of the different K_{β} 's. Without giving the derivation, the cumulative distribution function for activating the segment at the distance *l'* under the condition that the point *l* has been selected before is

$$C_{l}(l',r) = 1 - \exp\left(-\frac{1}{\beta} \int_{l}^{l'} d\tau \int_{-r}^{0} d\tau' K_{\beta}(\tau - \tau')\right).$$
(B17)

For the required inversion we recast (B17) to the form

$$-\log[1 - C(l,r)] = T_r(l') - T_r(l),$$
(B18)

where the function $T_r(l) = [K_{\beta}^{(2)}(l+r) - K_{\beta}^{(2)}(l)]/\beta$ contains the second integral of the kernel $K_{\beta}(t)$.

In practice, we pre-calculate $K_{\beta}^{(2)}(\tau)$ by means of an adaptive spline interpolation for quicker evaluation during the simulation. The essential steps of the algorithm are summarized in Fig. 21.

APPENDIX C: LOW-TEMPERATURE MEAN-FIELD SOLUTION

We start with the discretized form of the partition function (20) with the abbreviation for the interaction (23)

$$Z = \sum_{\{s_{m,i}\}} \exp\left[\frac{1}{2N} \sum_{m,m'}^{N} \sum_{i,j}^{L} J_{i,j} s_{m,i} s_{m',j}\right]$$

Introducing continuous variables m_i by means of delta functions leads to

$$Z = \sum_{\{s_{m,i}\}} \exp\left[\frac{N}{2} \sum_{i,j}^{L} J_{i,j} \left(\sum_{m=1}^{N} \frac{s_{m,i}}{N}\right) \left(\sum_{m'=1}^{N} \frac{s_{m',j}}{N}\right)\right]$$
$$= \sum_{\{s_{m,i}\}} \int \prod_{i=1}^{L} \left[dm_i \delta\left(m_i - \sum_{m=1}^{N} \frac{s_{m,i}}{N}\right)\right]$$
$$\times \exp\left[\frac{N}{2} \sum_{i,j}^{L} J_{i,j} m_i m_j\right].$$

The delta functions can substituted by

$$\delta\left(m_{i}-\sum_{m=1}^{N}\frac{s_{m,i}}{N}\right)=N\delta\left(Nm_{i}-\sum_{m=1}^{N}s_{m,i}\right)$$
$$=\frac{N}{2\pi}\int\exp\left[i\hat{m}_{i}\left(Nm_{i}-\sum_{m=1}^{N}s_{m,i}\right)\right]d\hat{m}_{i},$$

which leads to

$$Z = \int \prod_{i=1}^{L} \left[\frac{d\hat{m}_i dm_i}{2\pi N^{-1}} \right] \sum_{\{s_{m,i}\}} \exp\left(iN \sum_{i=1}^{L} \hat{m}_i m_i\right)$$
$$\times \exp\left(-i \sum_{i,m} \hat{m}_i s_{m,i}\right) \exp\left[\frac{N}{2} \sum_{i,j}^{L} J_{i,j} m_i m_j\right].$$

Now, the trace over the spin variables $s_{m,i}$ can performed directly:

$$Z = \int \prod_{i=1}^{L} \left[\frac{d\hat{m}_i dm_i}{2\pi N^{-1}} \right] \prod_i \left[2\cos\left(\hat{m}_i\right) \right]^N$$
$$\times \exp\left(iN \sum_{i=1}^{L} \hat{m}_i m_i \right) \exp\left[\frac{N}{2} \sum_{i,j}^{L} J_{i,j} m_i m_j \right]$$
$$= \int \prod_{i=1}^{L} \left[\frac{d\hat{m}_i dm_i}{2\pi N^{-1}} \right] \exp\left[N f(\{m_i\}, \{\hat{m}_i\}) \right],$$

where in the last line the function

$$f(\{m_i\},\{\hat{m}_i\}) = \ln[2\cos(\hat{m}_i)] + i\sum_{i=1}^{L} \hat{m}_i m_i + \frac{1}{2}\sum_{i,j}^{L} J_{i,j} m_i m_j$$

was defined. In the limit $N \rightarrow \infty$, a saddle point integration leads finally to the common mean-field relation for the magnetization

$$\frac{\partial f}{\partial m_i} = 0 = \frac{\partial f}{\partial \hat{m}_i} \implies m_i = \tanh\left(\sum_{j=1}^L J_{i,j}m_j\right). \quad (C1)$$

In equilibrium, all $m_i = m$ are equal and the phase transition point can calculated from (C1) by the condition $\sum_j J_{i,j} = 1$. Substituting back $J_{i,j}$ (23), one obtains

$$\sum_{j=1}^{L} J_{i,j} = \int_{0}^{\beta} \frac{1}{\Delta \beta} \tilde{K}_{\beta}(\tau - \tau') d\tau = \frac{2\omega_{c}\tilde{\alpha}}{\Delta s}, \qquad (C2)$$

where the replacement of the sum by the integral is valid for $O(1) \ll \Delta\beta$. Setting Eq. (C2) equal to 1 recovers the critical point.

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