Coherent-state path integral versus coarse-grained effective stochastic equation of motion: From reaction diffusion to stochastic sandpiles

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We derive and study two different formalisms used for nonequilibrium processes: the coherent-state path integral, and an effective, coarse-grained stochastic equation of motion. We first study the coherent-state path integral and the corresponding field theory, using the annihilation process $A+A\to A$ as an example. The field theory contains counterintuitive quartic vertices. We show how they can be interpreted in terms of a first-passage problem. Reformulating the coherent-state path integral as a stochastic equation of motion, the noise generically becomes imaginary. This renders it not only difficult to interpret, but leads to convergence problems at finite times. We then show how alternatively an effective coarse-grained stochastic equation of motion with real noise can be constructed. The procedure is similar in spirit to the derivation of the mean-field approximation for the Ising model, and the ensuing construction of its effective field theory. We finally apply our findings to stochastic Manna sandpiles. We show that the coherent-state path integral is inappropriate, or at least inconvenient. As an alternative, we derive and solve its mean-field approximation, which we then use to construct a coarse-grained stochastic equation of motion with real noise.

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I. INTRODUCTION

Stochastic processes are ubiquitous in nature: Think of gold particles suspended in water, which aggregate upon collision [1,2], a beautiful realization of the diffusion aggregation (or annihilation) process $A + A \rightarrow A$. Think of sand grains rolling down a hill, or its cellular automaton representatives, such as the Bak-Tang-Wiesenfeld [3] or the Manna sandpile [4] models. Even simpler, think of a large number of particles diffusing. To understand the physical properties of these systems, several routes are open: One may start with a direct numerical simulation of say the mentioned gold particles. For technical reasons this study would be restricted to a relatively small number of particles. Thus, in a second step, one strives for a more efficient effective description. This could be achieved by dividing the system into boxes of size ℓ , counting the number of particles inside each box, and trying to derive an effective coarse-grained description for the evolution of the number of particles inside each box. The question then arises, how do we do this?

Let us step back and consider an example from equilibrium statistical mechanics: In order to understand the phase transition between the ferromagnetic and the paramagnetic phases in a ferromagnet, or the liquid-gas transition in water, one first reduces these phenomena to the simplest possible model, in both cases the Ising model. The latter can be studied numerically, or through analytic techniques. An analytic treatment may start from the mean-field approximation, and then progress to the construction of a coarse-grained model, also termed effective field theory. What one learns from mean-field theory enters into the effective field theory as the description inside a box, usually with one or a few degrees of freedom (fields). This construction has to be supplied with an additional coupling between boxes, which completes the effective field theory. It can then be analyzed with renormalization-group techniques. The latter are expected to give the correct universal properties, such as the divergence of the specific heat when approaching

the critical point, even though the precise location of the phase transition temperature itself has been lost when constructing the coarse-grained description inside a single box.

Coming back to our discussion of the aggregating gold particles, the key point is the derivation of an effective field theory. There are two general-purpose methods to do so, both with their unique strengths and weaknesses: the coherent-state path integral (CSPI) and the coarse-grained stochastic equation of motion (CGSEM). In these notes, we will study both techniques side by side.

The coherent-state path integral (CSPI) has proven to be a useful tool, both for quantum many-body problems [5,6] and in statistical mechanics [7–10]. Despite its success, e.g., for reaction-diffusion processes, its use led to quite some confusion. Indeed, as we will see below, the CSPI quite naturally introduces an imaginary noise, rendering a physical interpretation difficult. The literature on the subject is vast [5–15], but leaves unanswered key questions the author of these notes asked himself. It is his intention to close this gap.

We start by giving a pedagogical introduction to the CSPI (Sec. II). This is mostly standard, following the work by Doi [7,8], Peliti [9], and the beautiful introduction by Cardy [10].

For concreteness, we then focus on reaction-diffusion processes, such as the gold aggregation process discussed in the beginning, and construct a field-theory action. This kind of processes leads to the appearance of some surprising and seemingly counterintuitive vertices in the field theory, which are a consequence of the conservation of probability. We show how they can be interpreted in terms of a first-passage problem (Sec. III).

The field theory can then be reformulated as a stochastic equation of motion (Sec. IV). It has an imaginary noise, which gives rise to some puzzlement in the literature [11,16–19]. As we will show below, contrary to real noise imaginary noise leads to a narrowing of the probability distribution. As the basis of the CSPI is coherent states, equivalent to Poisson distributions, the presence of an imaginary noise tells us that

over time the probability distribution becomes narrower than a Poissonian distribution. Coding for such narrow distributions with Poissonians is only possible via complex states, i.e., interference. We show, and check numerically, that the stochastic evolution of the coherent states allows us to sample directly the evolution of the discrete probability distribution, starting from the initial Poisson distribution

$$p_{i}(n) = e^{-a_{i}} \frac{(a_{i})^{n}}{n!}$$
 (1)

to the final distribution

$$p_{\mathbf{f}}(n) = \left\langle e^{-a_{\mathbf{f}}} \frac{(a_{\mathbf{f}})^n}{n!} \right\rangle_{\xi}.$$
 (2)

The average goes over end points $a_{\rm f}$ of trajectories of the stochastic equation of motion, for different realizations of the noise $\xi_{\rm f}$. As we will see, this formalism breaks down when $a_{\rm f}$ has diffused "too far" away from the positive real axis. This is an intrinsic problem of the CSPI, and difficult to repair.

As mentioned above, there is a second, alternative approach (Sec. V). To this aim, one replaces the discrete number of particles on a given site by a continuous variable: either as an ad hoc procedure or via coarse graining introducing the particle density in a box of size ℓ . Demanding that the resulting continuous random process has the same drift and variance as the underlying discrete process leads to an effective, coarsegrained stochastic equation of motion (CGSEM) with drift and real noise. As in the CSPI, its amplitude is proportional to the square root of the drift term, with the difference that the latter is real, while the former is imaginary. Contrary to the process with imaginary noise in the CSPI, the real process converges well for all times, and yields an efficient effective description. It is at the basis of most effective stochastic field theories. However, the stochastic equations of motion are rarely derived, even though the procedure given below is quite generally applicable. Most often, the stochastic equations of motion are conjectured on symmetry grounds, more obscuring than enlightening their origin.

We then proceed to a nontrivial example, the stochastic Manna sandpile (Sec. VI). The rules are simple: If the number of grains on a site exceeds one, two grains are redistributed, or toppled, onto randomly chosen neighbors. We study this model numerically, and show that coherent states are not an appropriate basis for a coarse-grained, stochastic description: On one hand, the probability distribution for Manna sandpiles is an exponential, and not a Poissonian, as the coherent state. On the other hand, while a Poisson distribution has one parameter, we observe that Manna sandpiles are characterized by two parameters. Thus, while a description in terms of a CSPI is always possible (and at least for short times exact), it is also plagued by the appearance of complex states, and the corresponding convergence problems. In hindsight, passing into the complex plane may not be surprising, as it can be interpreted as the "trick" of the CSPI to generate a second dynamic variable. We then turn to a more efficient description, and construct an effective stochastic field theory. To this aim, we define a variant of the Manna model, the range-r Manna model: its toppling rules are modified such that grains end up not on neighboring sites but on sites within a distance of r. We show that it converges for $r \to \infty$ to a mean-field model,

which we solve analytically. Using this mean-field model as a coarse-grained description for an elementary box, we derive a stochastic field theory for the Manna model. This field theory is known [20–23]. The advantage of the present scheme is that we do not have to invoke symmetry arguments, and that our scheme fixes all parameters, restricting the classes of models to be considered to a submanifold, which is equivalent to the simplest dissipative dynamics of a driven disordered manifold [24].

After the conclusion (Sec. VII), the reader will find several appendices to which more technical details have been relegated. There we also discuss a canonical change of variables which, when applicable, allows to avoid imaginary noise.

II. THE COHERENT-STATE PATH INTEGRAL

The coherent-state path integral (CSPI) is a formalism which evaluates exactly the evolution of probabilities for a stochastic process. To this aim, the different configurations of the system are represented as in quantum mechanics by n-particle states $|n\rangle$. This allows us to write probabilities p(n) as states, i.e., superpositions $|\psi\rangle := \sum_{n=0}^{\infty} p(n)|n\rangle$. The evolution operator is then encoded into a Hamiltonian, acting on these states. Finally, a path integral is introduced. Its eigenstates are coherent states, i.e., eigenfunctions of the annihilation operator to be defined below.

Having constructed an exact representation of the stochastic process as a coherent-state path integral, the latter can be studied with different methods: either using perturbation theory, possibly coupled with renormalization-group methods (Sec. III), or by rewriting it as a stochastic equation of motion for the states $|\psi\rangle$ (Sec. IV). We will study these techniques in turn.

A. Quantization rules

Consider a single site which can be occupied by n particles (bosons), $n = 0, 1, \ldots$ Denote this n-particle state by

$$|n\rangle := (\hat{a}^{\dagger})^n |0\rangle, \tag{3}$$

where $|0\rangle$ is the normalized vacuum state $\langle 0|0\rangle = 1$. While \hat{a}^{\dagger} is the creation operator, its conjugate \hat{a} is the annihilation operator, $\hat{a}|0\rangle = 0$. They have canonical commutation rules

$$[\hat{a}, \hat{a}^{\dagger}] = 1. \tag{4}$$

The scalar product between two states is

$$\langle n|m\rangle = \langle 0|\hat{a}^n(\hat{a}^\dagger)^m|0\rangle = n!\,\delta_{nm}.\tag{5}$$

This is proven by commuting all \hat{a} to the right, using Eq. (4). Thus $|n\rangle$ is not normalized to 1, but to $\langle n|n\rangle = n!$. The number operator is $\hat{n} := \hat{a}^{\dagger}\hat{a}$, i.e.,

$$\hat{n}|n\rangle \equiv \hat{a}^{\dagger}\hat{a}|n\rangle = n|n\rangle. \tag{6}$$

We note for convenience that

$$\hat{a}\hat{a}^{\dagger}|n\rangle = (n+1)|n\rangle,\tag{7}$$

$$\hat{a}^2(\hat{a}^\dagger)^2|n\rangle = (n+1)(n+2)|n\rangle,\tag{8}$$

$$(\hat{a}^{\dagger})^2 \hat{a}^2 |n\rangle = n(n-1)|n\rangle. \tag{9}$$

B. Master equation and Hamiltonian formalism

We now want to code a master equation for the occupation probability in this formalism. Suppose the probability for having n particles at time t is $p_t(n)$, with $\sum_{n=0}^{\infty} p_t(n) = 1$. We associate with this probability a state

$$|\psi_t\rangle := \sum_{n=0}^{\infty} p_t(n)|n\rangle \equiv \sum_{n=0}^{\infty} p_t(n)(\hat{a}^{\dagger})^n|0\rangle.$$
 (10)

Consider the master equation for the probability $p_t(n)$,

$$\partial_t p_t(n) = \frac{\nu}{2} [(n+1)n \ p_t(n+1) - n(n-1)p_t(n)] + \mu [(n+1)p_t(n+1) - np_t(n)] + \kappa [(n-1)p_t(n-1) - np_t(n)].$$
(11)

In the first process two particles meet and annihilate with rate ν : $A + A \xrightarrow{\nu} A$. In the second process, a particle decays with rate μ : $A \xrightarrow{\mu} \emptyset$. In the third process a particle "gives birth" to two particles with rate κ : $A \xrightarrow{\kappa} A + A$. Note that probability is conserved, $\sum_{n=0}^{\infty} \partial_t p_t(n) = 0$.

We now want to derive the "Hamiltonian" associated with this master equation, in the form

$$\partial_t |\psi_t\rangle = \mathcal{H}|\psi_t\rangle. \tag{12}$$

To this aim we multiply both sides of Eq. (11) with $(\hat{a}^{\dagger})^n |0\rangle$, and then sum over n. The factors of n are expressed using the number operator $\hat{n} = \hat{a}^{\dagger}\hat{a}$,

$$\partial_{t} \sum_{n=0}^{\infty} p_{t}(n) (\hat{a}^{\dagger})^{n} |0\rangle
= \frac{\nu}{2} \sum_{n=0}^{\infty} [p_{t}(n+1) \hat{a}^{\dagger} \hat{a}^{2} \hat{a}^{\dagger} - p_{t}(n) (\hat{a}^{\dagger})^{2} \hat{a}^{2}] (\hat{a}^{\dagger})^{n} |0\rangle
+ \mu \sum_{n=0}^{\infty} [p_{t}(n+1) \hat{a} \hat{a}^{\dagger} - p_{t}(n) \hat{a}^{\dagger} \hat{a}] (\hat{a}^{\dagger})^{n} |0\rangle
+ \kappa \sum_{n=0}^{\infty} [p_{t}(n-1) (\hat{a}^{\dagger} \hat{a} - 1) - p_{t}(n) \hat{a}^{\dagger} \hat{a}] (\hat{a}^{\dagger})^{n} |0\rangle.$$
(13)

Note that we have taken advantage of relations (7) to (9) to simplify the expression. Next we use definition (10) to rewrite this expression in terms of $|\psi_t\rangle$. As an example consider the first term on the right-hand side, $\sum_{n=0}^{\infty} \hat{a}^{\dagger} \hat{a}^2 p_t (n+1) (\hat{a}^{\dagger})^{n+1} |0\rangle \equiv \sum_{n=0}^{\infty} \hat{a}^{\dagger} \hat{a}^2 p_t (n) (\hat{a}^{\dagger})^n |0\rangle = \hat{a}^{\dagger} \hat{a}^2 |\psi_t\rangle$. We extended the sum to n=0, which is possible since the first term on the left-hand side does not contribute, due to the preceding operators \hat{a}^2 .

We thus arrive at

$$\partial_t |\psi_t\rangle = \frac{\nu}{2} [\hat{a}^{\dagger} \hat{a}^2 - (\hat{a}^{\dagger})^2 \hat{a}^2] |\psi_t\rangle + \mu [\hat{a} - \hat{a}^{\dagger} \hat{a}] |\psi_t\rangle + \kappa [(\hat{a}^{\dagger})^2 \hat{a} - \hat{a}^{\dagger} \hat{a}] |\psi_t\rangle.$$
(14)

Using Eq. (12), this identifies the Hamiltonian

$$\mathcal{H} = \frac{\nu}{2} [\hat{a}^{\dagger} \hat{a}^2 - (\hat{a}^{\dagger})^2 \hat{a}^2] + \mu [\hat{a} - \hat{a}^{\dagger} \hat{a}] + \kappa [(\hat{a}^{\dagger})^2 \hat{a} - \hat{a}^{\dagger} \hat{a}]. \tag{15}$$

This Hamiltonian is normal-ordered; i.e., all \hat{a}^{\dagger} stand left of all \hat{a} . It has all the terms expected from quantum mechanics,

except that for each expected term there is a second term which does not change the particle number, and which ensures the conservation of probability. Indeed, conservation of probability can be written as

$$0 = \partial_t \sum_{n=0}^{\infty} p_t(n) \equiv \partial_t \langle 0 | e^{\hat{a}} | \psi_t \rangle$$

$$= \langle 0 | e^{\hat{a}} \mathcal{H}(\hat{a}^{\dagger}, \hat{a}) | \psi_t \rangle$$

$$= \langle 0 | \mathcal{H}(\hat{a}^{\dagger} + 1, \hat{a}) e^{\hat{a}} | \psi_t \rangle. \tag{16}$$

For the first line we used that the 1/n! in the definition of the exponential function cancels the normalization (5). For the second line we used that

$$e^{\lambda \hat{a}} f(\hat{a}^{\dagger}) = f(\hat{a}^{\dagger} + \lambda)e^{\lambda \hat{a}},$$
 (17)

$$e^{\lambda \hat{a}^{\dagger}} f(\hat{a}) = f(\hat{a} - \lambda) e^{\lambda \hat{a}^{\dagger}}.$$
 (18)

Noting that an \hat{a}^{\dagger} inside \mathcal{H} , when acting to the left on $\langle 0|$, gives no contribution, we arrive at the constraint of conservation of probability for the normal-ordered Hamiltonian \mathcal{H} :

$$\mathcal{H}(\hat{a}^{\dagger}, \hat{a})|_{\hat{a}^{\dagger} \to 1} = 0. \tag{19}$$

Equation (19) is a necessary condition to ensure that (16) holds; it is also sufficient since using (17) the state $e^{\hat{a}}|\psi_t\rangle \equiv \sum_n p_t(n)(\hat{a}^{\dagger}+1)^n|0\rangle$ can be chosen arbitrarily. Looking back at Eq. (15), we see that the second term inside each square bracket is such that at $\hat{a}^{\dagger}=1$ the sum of the two terms vanishes; thus as stated it ensures the conservation of probability.

C. Combinatorics

Let us remark that the combinatorics used in the above processes is the basic combinatorics of choosing k out of n particles, $\binom{n}{k}$, relevant, e.g., for the meeting probability of two particles. While this choice is canonic, situations may arise where the combinatorics is different. If the stochastic process was to contain factors nonpolynomial in n, then the Hamiltonian (15) would not be as simple, and might, e.g., become nonanalytic in \hat{a} and \hat{a}^+ ; much of the technology developed here would no longer work. This holds especially true for the stochastic equation of motion to be introduced below, which relies on the fact that, via a suitable decoupling, the Hamiltonian can be rendered linear in \hat{a}^{\dagger} .

D. Observables

Now consider an observable $\mathcal{O}(n)$, which depends only on the occupation number n. Using the same tricks as in Eq. (16), its expectation value can be written as

$$\langle \mathcal{O} \rangle_{\psi_t} := \sum_{n=0}^{\infty} \mathcal{O}(n) p_t(n) = \langle 0 | e^{\hat{a}} \mathcal{O}(\hat{a}^{\dagger} \hat{a}) | \psi_t \rangle$$

$$\equiv \langle 0 | \mathcal{O}(\hat{a}^{\dagger} \hat{a} + \hat{a}) e^{\hat{a}} | \psi_t \rangle = \langle 0 | \mathcal{O}_{N}(\hat{a}^{\dagger} + 1, \hat{a}) e^{\hat{a}} | \psi_t \rangle$$

$$= \langle 0 | \mathcal{O}_{N}(1, \hat{a}) e^{\hat{a}} | \psi_t \rangle. \tag{20}$$

From the first to the second line, we used Eq. (17). In the second line we have introduced the normal-ordered version of the operator \mathcal{O} , obtained by commuting all \hat{a} to the right and

all \hat{a}^{\dagger} to the left. It is generically a function of \hat{a} and \hat{a}^{\dagger} , not $\hat{n} = \hat{a}^{\dagger}\hat{a}$. The last line uses that \hat{a}^{\dagger} acting to the left vanishes.

E. Coherent states

Coherent states play a key role in the path-integral formalism to be developed below. We define them here, and study some of its properties. Coherent states are constructed such that

$$|\phi\rangle := e^{\phi \hat{a}^{\dagger}} |0\rangle \Rightarrow \hat{a} |\phi\rangle = \phi |\phi\rangle.$$
 (21)

Let us start with ϕ real and positive. Then by definition a coherent state has a Poisson probability distribution for n-fold occupation,

$$p(n) = e^{-\phi} \frac{\phi^n}{n!}. (22)$$

Note that the definition (21) does not contain the factor of $e^{-\phi}$; thus it is not normalized. This is for convenience reasons; one may think of it as a histogram.

States $|\phi\rangle$ with a complex ϕ are possible too. Since ϕ is continuous, but the number n an integer, coherent states form an overcomplete basis, even for $\phi \geqslant 0$. However, not all probability distributions can be written as a superposition of coherent states with positive weights, i.e., as

$$|\psi\rangle = \int_0^\infty d\phi \, \rho(\phi) e^{-\phi} \, |\phi\rangle \tag{23}$$

with $\rho(\phi) \ge 0$. There are several ways out of this dilemma: One can use negative (or complex) weights $\rho(\phi)$, states with complex ϕ , or a combination of both. The formalism to be developed below will exploit this freedom.

By definition, the adjoint state is

$$\langle \phi^* | = \langle 0 | e^{\phi^* \hat{a}}. \tag{24}$$

Equation (17) implies that the scalar product is

$$\langle \phi^* | \phi \rangle = e^{\phi^* \phi}. \tag{25}$$

Let us give an interpretation of the adjoint state: Apply $\langle \phi^* |$ to the state $|\psi_t\rangle$ defined in Eq. (10),

$$\rho_t(\phi^*) := \langle \phi^* | \psi_t \rangle = \sum_{n=0}^{\infty} p_t(n) (\phi^*)^n.$$
 (26)

This is nothing but the generating function of the probabilities $p_t(n), n = 0, 1, 2, \dots$

Now consider the expectation value of a normal-ordered observable $\mathcal{O}(\hat{n}) = \mathcal{O}_N(\hat{a}^{\dagger}, \hat{a})$ in a coherent state $|\phi\rangle$,

$$\langle \mathcal{O} \rangle_{\phi} := \frac{\langle 0 | e^{\hat{a}} \mathcal{O}_{N}(\hat{a}^{\dagger}, \hat{a}) e^{\phi \hat{a}^{\dagger}} | 0 \rangle}{\langle 0 | e^{\hat{a}} \mathbb{1} e^{\phi \hat{a}^{\dagger}} | 0 \rangle}$$

$$= \langle 0 | e^{\phi \hat{a}^{\dagger}} \mathcal{O}_{N}((\hat{a}^{\dagger} + 1), (\hat{a} + \phi)) e^{\hat{a}} | 0 \rangle$$

$$= \mathcal{O}_{N}(1, \phi). \tag{27}$$

We used Eqs. (17) and (18), as well as the vanishing of \hat{a} acting on the vacuum to the right, and \hat{a}^{\dagger} to the left. To write the last line \mathcal{O}_N needs to be normal-ordered. Also note that a factor of e^{ϕ} has canceled between the numerator and denominator of the first line; it is necessary, since our coherent states (21) are not normalized to unity.

In coherent states, the number of particles is not fixed. We show in Appendix B 1 that

$$e^{\lambda \hat{n}} \equiv e^{\lambda \hat{a}^{\dagger} \hat{a}} = : e^{(e^{\lambda} - 1)\hat{a}^{\dagger} \hat{a}} : . \tag{28}$$

The right-hand side is called *normal-ordered* and denoted by ":" around the operators in question; it is defined by its Taylor expansion in \hat{a}^{\dagger} and \hat{a} , and then arranging all \hat{a}^{\dagger} to the left and all \hat{a} to the right, as if they were numbers. An example is : $(\hat{a}^{\dagger}\hat{a})^2$: defined to be $(\hat{a}^{\dagger})^2\hat{a}^2$.

Using this relation, or directly the intermediate result Eq. (B3) at $\phi^* = 1$, and the definition of an observable given in Eq. (27) yields

$$\langle e^{\lambda \hat{n}} \rangle_{\phi} = e^{(e^{\lambda} - 1)\phi}. \tag{29}$$

The generating function of connected moments is the logarithm of this function,

$$\langle e^{\lambda \hat{n}} \rangle_{\phi}^{c} = (e^{\lambda} - 1)\phi. \tag{30}$$

This means that the pth connected moment of the number operator n is

$$\langle \hat{n}^p \rangle_{\phi}^{c} = \phi. \tag{31}$$

Let us give some explicit examples:

$$\langle \hat{n} \rangle_{\phi} = \phi, \tag{32}$$

$$\langle \hat{n}^2 \rangle_{\phi} = \phi(1+\phi), \tag{33}$$

$$\langle \hat{n}^3 \rangle_{\phi} = \phi (1 + 3\phi + \phi^2), \tag{34}$$

:

$$\phi = \langle n \rangle_{\phi},\tag{35}$$

$$\phi^2 = \langle \hat{n}(\hat{n} - 1) \rangle_{\phi},\tag{36}$$

$$\phi^3 = \langle \hat{n}(\hat{n} - 1)(\hat{n} - 2)\rangle_{\phi},\tag{37}$$

:

The last set of relations can also be derived directly; see Appendix B 2.

F. Many sites

We now generalize to L sites, denoted $i=1,\ldots,L$, with creation and annihilation operators \hat{a}_i^{\dagger} and \hat{a}_i for site i. The canonical commutation relations are a generalization of Eq. (4):

$$[\hat{a}_i, \hat{a}_i^{\dagger}] = \delta_{ij}. \tag{38}$$

A state is then encoded as

$$|\psi\rangle = \sum_{n_1,\dots,n_L=0}^{\infty} p_t(n_1,\dots,n_L)(\hat{a}_1^{\dagger})^{n_1}\dots(\hat{a}_L^{\dagger})^{n_L}|0\rangle.$$
 (39)

We can also construct a coherent state out of single-particle coherent states.

$$|\psi\rangle := \bigotimes_{i=1}^{L} |\phi_i\rangle. \tag{40}$$

G. Coarse graining

When constructing effective field theories, one often coarse grains, replacing the state variables of several sites by a common effective variable. For coherent states, this is particularly straightforward: Suppose we have two sites with coherent states $|\phi_1\rangle$ and $|\phi_2\rangle$, and we want to know what the probability to have n-fold occupation of the combined two sites is. We evaluate

$$p_{\text{comb}}(n) = \sum_{n_1=0}^{n} \left[e^{-\phi_1} \frac{(\phi_1)^{n_1}}{n_1!} \right] \times \left[e^{-\phi_2} \frac{(\phi_2)^{n-n_1}}{(n-n_1)!} \right]$$
$$= e^{-(\phi_1 + \phi_2)} \frac{(\phi_1 + \phi_2)^n}{n!}. \tag{41}$$

Thus, combining two coherent states leads to a coherent state with the added weights,

$$|\phi_1\rangle \oplus |\phi_2\rangle \longrightarrow |\phi_1 + \phi_2\rangle.$$
 (42)

Finally, if we are interested in the probability for n-particle occupation of our system of size L given in Eq. (40), we get a state

$$|\Phi\rangle = \left|\sum_{i=1}^{L} \phi_i\right\rangle. \tag{43}$$

H. Diffusion

Consider now the hopping of a particle from site i to site j, with diffusion constant (rate) D. The corresponding Hamiltonian is (as expected)

$$\mathcal{H} = D(\hat{a}_i^{\dagger} - \hat{a}_i^{\dagger})\hat{a}_i. \tag{44}$$

Having hopping both from i to j and from j to i with the same rate D leads to

$$\mathcal{H} = -D(\hat{a}_i^{\dagger} - \hat{a}_i^{\dagger})(\hat{a}_j - \hat{a}_i). \tag{45}$$

Note that by this definition the rate to leave a site in dimension d is $2d \times D$, and not D. In the continuum limit, and summing over all nearest-neighbor sites, this becomes the Hamiltonian of diffusion

$$\mathcal{H}_{\text{diffusion}} = -D \int_{x} \nabla \hat{a}_{x}^{\dagger} \nabla \hat{a}_{x}. \tag{46}$$

To avoid overly cumbersome notations, we have set to 1 the lattice cutoff a, which multiplies the lattice diffusion constant D by a factor of a^{2-d} .

I. Resolution of unity

The path-integral representation we wish to establish is based on the coherent states defined in Eq. (21). The key relation which we are going to prove is the resolution of unity:

$$1 = \frac{i}{2\pi} \int d\phi \, d\phi^* e^{-\phi\phi^*} |\phi\rangle\langle\phi^*|. \tag{47}$$

The complex-conjugate pair is $\phi = \phi_x + i\phi_y$, $\phi^* = \phi_x - i\phi_y$; the integration measure is $d\phi d\phi^* = \frac{2}{i}d\phi_x d\phi_y$. Inserting these definitions, the right-hand side of Eq. (47) can be rewritten

as

$$\int \frac{d\phi_x d\phi_y}{\pi} e^{-\phi\phi^*} e^{\phi\hat{a}^{\dagger}} |0\rangle \langle 0| e^{\phi^*\hat{a}}$$

$$= \sum_{n=0}^{\infty} \sum_{m=0}^{\infty} \int_0^{2\pi} \frac{d\theta}{\pi} \int_0^{\infty} dr \, r e^{-r^2} r^{n+m} e^{i\theta(n-m)}$$

$$\times \frac{(\hat{a}^{\dagger})^n}{n!} |0\rangle \langle 0| \frac{\hat{a}^m}{m!}, \tag{48}$$

where in the last line we set $\phi := re^{i\theta}$. The angular integral is vanishing for $n \neq m$, resulting in

$$\sum_{n=0}^{\infty} \int_{0}^{\infty} d(r^{2}) e^{-r^{2}} r^{2n} \frac{(\hat{a}^{\dagger})^{n}}{n!} |0\rangle\langle 0| \frac{\hat{a}^{n}}{n!} = \sum_{n=0}^{\infty} \frac{1}{n!} (\hat{a}^{\dagger})^{n} |0\rangle\langle 0| \hat{a}^{n}.$$
(49)

Applying this expression to the state $|m\rangle = (\hat{a}^{\dagger})^m |0\rangle$, only the term n = m in the sum contributes, and reproduces this state. This completes the proof.

Let us mention another commonly employed trick, namely of analytic continuation. This is most prominently employed in conformal field theory; see, e.g., Ref. [25], to which we refer the reader for the subtleties. The essence is that ϕ and ϕ^* do not have to be complex conjugates, but that one may think of them as two independent variables, which together span $\mathbb{C} \equiv \mathbb{R}^2$. A conceptually convenient choice is ϕ real and ϕ^* imaginary.

J. Evolution operator in the coherent-state formalism, and action

We are now in a position to construct the time evolution in the coherent-state formalism. To this aim write the evolution operator $e^{\delta t\mathcal{H}} \simeq 1 + \delta t\mathcal{H}$ for a small time, and evaluate it in the coherent basis, by applying the resolution of unity (47) to both sides of $e^{\delta t\mathcal{H}}$. To avoid problems with not normal-ordered terms appearing in $(\mathcal{H})^2$, and higher, we choose δt infinitesimally small:

$$e^{\delta t \mathcal{H}(\hat{a}^{\dagger}, \hat{a})} = \frac{i}{2\pi} \int d\phi_{t+\delta t} \, d\phi_{t+\delta t}^* \, \frac{i}{2\pi} \int d\phi_t \, d\phi_t^*$$

$$\times e^{-\phi_{t+\delta t}\phi_{t+\delta t}^* - \phi_t \phi_t^*} |\phi_{t+\delta t}\rangle \langle \phi_t^*|$$

$$\times \langle \phi_{t+\delta t}^* | e^{\delta t \mathcal{H}(\hat{a}^{\dagger}, \hat{a})} |\phi_t\rangle.$$
(50)

We need to evaluate the matrix element in question:

$$\begin{split} \langle \phi_{t+\delta t}^* | e^{\delta t \mathcal{H}(\hat{a}^{\dagger}, \hat{a})} | \phi_t \rangle \\ &\simeq \langle 0 | e^{\phi_{t+\delta t}^* \hat{a}} [1 + \delta t \mathcal{H}(\hat{a}^{\dagger}, \hat{a})] e^{\phi_t \hat{a}^{\dagger}} | 0 \rangle \\ &= e^{\phi_{t+\delta t}^* \phi_t} \langle 0 | e^{\phi_t \hat{a}^{\dagger}} [1 + \delta t \mathcal{H}(\hat{a}^{\dagger} + \phi_{t+\delta t}^*, \hat{a} + \phi_t)] e^{\phi_{t+\delta t}^* \hat{a}} | 0 \rangle. \end{split} \tag{51}$$

We have used Eqs. (17) and (18) to commute the exponential operators. All operators \hat{a} are now acting on the vacuum to

¹Consider the scalar product $\langle n|1|m\rangle \sim \int d\phi \, d\phi^* \, \phi^n (\phi^*)^m e^{-\phi\phi^*} = \int d\phi \, d\phi^* \, \phi^n (-\partial_\phi)^m e^{-\phi\phi^*} = m! \int d\phi \, d\phi^* \, \phi^{n-m} e^{-\phi\phi^*} \, \Theta(n>m)$. For ϕ real and ϕ^* purely imaginary the integral $\int d\phi^* e^{-\phi\phi^*}$ yields $\delta(\phi)$, and the former expression vanishes except for n=m.

the right, thus do not give a contribution. The same holds true for \hat{a}^{\dagger} acting to the left. Further using the normalization of the vacuum state $\langle 0|0\rangle=1$, we finally arrive at

$$\langle \phi_{t+\delta t}^* | e^{\delta t \mathcal{H}(\hat{a}^{\dagger}, \hat{a})} | \phi_t \rangle = e^{\phi_{t+\delta t}^* \phi_t} e^{\delta t \mathcal{H}(\phi_{t+\delta t}^*, \phi_t)} + O(\delta t^2).$$
 (52)

Together with Eq. (50), we identify all terms for a time step from t to $t + \delta t$ as $e^{-S_{t,t+\delta_t}\delta t}$, with

$$-S_{t,t+\delta_t}\delta t = [\phi_{t+\delta_t}^* - \phi_t^*]\phi_t + \mathcal{H}(\phi_{t+\delta_t}^*, \phi_t)\delta t$$

= $\phi_{t+\delta_t}^* [\phi_t - \phi_{t+\delta_t}] + \mathcal{H}(\phi_{t+\delta_t}^*, \phi_t)\delta t$. (53)

The expression $S_{t,t+\delta_t}$ is termed the *action* for the time step from t to $t + \delta t$. The two possible forms were obtained by grouping with either of the factors of $e^{-\phi_{t+\delta_t}\phi_{t+\delta_t}^*-\phi_t\phi_t^*}$ appearing in Eq. (50). Suppose for the following that we evolve from small to large times: Then the second line will be relevant; the unused factor of $e^{-\phi_t\phi_t^*}|_{t=t_i}$ will appear together with the initial state ϕ_i as

$$e^{-\phi_t \phi_t^*} \langle 0 | e^{\phi_t^* \hat{a}} e^{\phi_i \hat{a}^\dagger} | 0 \rangle = e^{-(\phi_t - \phi_i) \phi_t^*}, \tag{54}$$

where we used Eq. (17); when integrated over ϕ_t^* , this identifies ϕ_t as the initial state ϕ_i . Note that when integrating from larger to smaller times, we would use the first line of Eq. (53), evolving from the final state $\langle \phi_f |$ to smaller times; the factor $e^{-\phi_{t+\delta t}^*\phi_{t+\delta t}}|_{t+\delta t=t_f}$ would then fix $\langle \phi_f | = \langle \phi_{t+\delta t} |$. The formalism can thus be used both forward and backward in time, exchanging the role of ϕ and ϕ^* .

We now consider the forward version, evolving from an initial state $|\phi_i\rangle$ at $t=t_i$. In the continuous limit, the action from time $t=t_i$ to time $t=t_f$ becomes

$$\mathcal{S}[\phi^*,\phi] := \int_{t_i}^{t_f} dt \, \phi_t^* \partial_t \phi_t - \mathcal{H}[\phi_t^*,\phi_t]. \tag{55}$$

(Note that we replaced $\phi_{t+\delta t}^* \to \phi_t^*$ in the Hamiltonian $\mathcal{H}[\phi_{t+\delta t}^*, \phi_t]$, which is valid in the small- δt limit.)

The path integral can then be written as

$$|\psi_{\rm f}\rangle = \mathbf{T}e^{\int_{t_{\rm i}}^{t_{\rm f}} dt \mathcal{H}[\hat{a}_{t}^{\dagger}, \hat{a}_{t}]} |\phi_{\rm i}\rangle$$

$$= \int \mathcal{D}[\phi] \, \mathcal{D}[\phi^{*}] e^{-\mathcal{S}[\phi^{*}, \phi]} \Big|_{\phi_{\rm f} = \phi_{\rm i}} |\phi_{t_{\rm f}}\rangle. \tag{56}$$

Note that (for one time slice) the state $|\phi_{t_f}\rangle$ corresponds to the state $|\phi_{t+\delta t}\rangle$ in Eq. (50), thus is part of the path integral (i.e., integrated over). On the other hand, the state $\langle \phi_t^*|$ in Eq. (50) corresponds to $\langle \phi_{t_i}^*|$. When applied to $|\phi_i\rangle$, and integrated over, it yields the boundary condition $\phi_{t_i} = \phi_i$.

The time index t at the operators \hat{a} and \hat{a}^{\dagger} is introduced for bookkeeping purposes, to define the time-ordering operator \mathbf{T} as $\mathbf{T}e^{\int_{t_i}^{t_f}dt\mathcal{H}[\hat{a}_i^{\dagger},\hat{a}_t]} := \prod_{t=t_i}^{t_f,\delta t=\tau}e^{\tau\mathcal{H}[\hat{a}_t^{\dagger},\hat{a}_t]}$, putting smaller times to the right. (This is the same ordering as in the definition of the path integral.)

The final state $|\psi_f\rangle$ is not a coherent state, but the superposition of coherent states $|\phi_f\rangle$. To formalize this better, suppose that the initial state is also a superposition of coherent states, each with weight $\rho(\phi_i)$, and normalized such that $\int_{\phi_i} \rho(\phi_i) := \frac{i}{2\pi} \int d\phi_i d\phi_i^* \rho(\phi_i) = 1$,

$$|\psi_{\rm i}\rangle = \int_{\phi_{\rm i}} \rho(\phi_{\rm i}) \, e^{-\phi_{\rm i}} |\phi_{\rm i}\rangle.$$
 (57)

Restricting support of $\rho(\phi_i)$ to $\phi_i > 0$ is included as a special case, with intuitive physical interpretation. The states $e^{-\phi_i}|\phi_i\rangle$ are normalized, so that together with the normalization of the weight the state $|\psi_i\rangle$ is normalized. Define

$$\mathcal{A}(\phi_{\mathrm{f}}|\phi_{i}) := \int \mathcal{D}[\phi] \, \mathcal{D}[\phi^{*}] \, e^{-\mathcal{S}[\phi^{*},\phi]} \Big|_{\phi_{\mathrm{f}}=\phi_{\mathrm{i}}}^{\phi_{\mathrm{f}}=\phi_{\mathrm{f}}}. \tag{58}$$

Then

$$|\psi_{\rm f}\rangle = \int_{\phi_{\rm f}} |\phi_{\rm f}\rangle \int_{\phi_{\rm i}} \mathcal{A}(\phi_{\rm f}|\phi_{\rm i}) \,\rho(\phi_{\rm i}) \,e^{-\phi_{\rm i}}. \tag{59}$$

By construction, $|\psi_f\rangle$ is normalized; thus $\mathcal{A}(\phi_f|\phi_i)$ defines the transition amplitude.

K. The shift $\phi_t^* \rightarrow \phi_t^* + 1$

In Eq. (20) we had considered expectation values of an observable \mathcal{O} . Suppose we want to measure it at time t_f , evolved from $|\phi_i\rangle$ at time t_i until time t_f ,

$$\langle \mathcal{O}_{t_c} \rangle = e^{-\phi_i} \langle 0 | e^{\hat{a}} \mathcal{O}_{N}(\hat{a}^{\dagger}, \hat{a}) \mathbf{T} e^{\int_{t_i}^{t_f} dt \mathcal{H}[\hat{a}_t^{\dagger}, \hat{a}_t]} | \phi_i \rangle. \tag{60}$$

The factor of $e^{-\phi_i}$ ensures that the initial state is normalized.

We remark now that $\langle 0|e^a=\langle 1|$. Going to the path integral, this can be written as

$$\begin{split} \left\langle \mathcal{O}_{t_{\mathrm{f}}} \right\rangle &= e^{-\phi_{\mathrm{i}}} \langle 1 | \mathcal{O}_{\mathrm{N}}(\hat{a}^{\dagger}, \hat{a}) \mathbf{T} e^{\int_{t_{\mathrm{i}}}^{t_{\mathrm{f}}} dt \mathcal{H}[\hat{a}_{t}^{\dagger}, \hat{a}_{t}]} | \phi_{\mathrm{i}} \rangle \\ &= \int_{\phi_{\mathrm{f}}} \int \mathcal{D}[\phi] \, \mathcal{D}[\phi^{*}] \, \mathcal{O}_{\mathrm{N}}(1, \phi_{\mathrm{f}}) \, e^{\phi_{\mathrm{f}} - \phi_{\mathrm{i}}} \, e^{-S[\phi^{*}, \phi]} \Big|_{\phi_{t_{\mathrm{f}}} = \phi_{\mathrm{f}}}^{\phi_{t_{\mathrm{f}}} = \phi_{\mathrm{f}}}. \end{split}$$

$$(61)$$

The final scalar product yields $\langle 1|\mathcal{O}_N(\hat{a}^\dagger,\hat{a})|\phi_f\rangle = \mathcal{O}_N(1,\phi_f)e^{\phi_f}$. Note that it does not fix $\phi_{t_f}^*=1$, as it would in absence of the operator $\mathcal{O}_N(\hat{a}^\dagger,\hat{a})$.

We now shift all variables $\phi^* \to \phi^* + 1$, to obtain

$$\langle \mathcal{O}_{t_{\rm f}} \rangle = \int \mathcal{D}[\phi] \, \mathcal{D}[\phi^*] \, \mathcal{O}_{\rm N}(1,\phi_{\rm f}) \, e^{-S'[\phi^*,\phi]} \Big|_{\phi_{\rm f} = \phi_{\rm i}}^{\phi_{\rm f_f} = \phi_{\rm f}}, \tag{62}$$

$$S'[\phi^*,\phi] := \int_t^{t_t} dt \, \phi_t^* \partial_t \phi_t - \mathcal{H}'[\phi_t^*,\phi_t], \tag{63}$$

$$\mathcal{H}'[\phi_t^*, \phi_t] := \mathcal{H}[\phi_t^* + 1, \phi_t].$$
 (64)

Note that under this shift

$$\int_{t_{i}}^{t_{f}} dt \, \phi_{t}^{*} \partial_{t} \phi_{t} \longrightarrow \int_{t_{i}}^{t_{f}} dt \, (\phi_{t}^{*} + 1) \partial_{t} \phi_{t}$$

$$= \phi_{t_{f}} - \phi_{t_{i}} + \int_{t_{i}}^{t_{f}} dt \, \phi_{t}^{*} \partial_{t} \phi_{t}. \tag{65}$$

Apart from the obvious change in the argument of \mathcal{H} this accounts for the cancellation of the factor of $e^{\phi_{\rm f}-\phi_{\rm i}}$ in Eq. (61). Cardy in his excellent lecture notes [10] calls this shift the Doi shift. Its main advantage is that the field ϕ^* has expectation zero (at least in the final state). This is particularly useful when interpreting the CSPI graphically, as we will see in the next section. In addition, the formulas are simpler, and more intuitive. Finally, it has advantages when evaluating the CSPI via a stochastic equation of motion; see Sec. IV. To distinguish between shifted and unshifted action, we put a prime on the

shifted one. In the shifted variables, Eqs. (58) and (59) take the form

$$\mathcal{A}'(\phi_{\mathrm{f}}|\phi_{i}) := \int \mathcal{D}[\phi] \, \mathcal{D}[\phi^{*}] \, e^{-S'[\phi^{*},\phi]} \Big|_{\phi_{\mathrm{f}}=\phi_{\mathrm{f}}}^{\phi_{\mathrm{f}}=\phi_{\mathrm{f}}}, \tag{66}$$

$$|\psi_{\rm f}\rangle = \int_{\phi_{\rm f}} e^{-\phi_{\rm f}} |\phi_{\rm f}\rangle \int_{\phi_{\rm i}} \mathcal{A}'(\phi_{\rm f}|\phi_{\rm i}) \,\rho(\phi_{\rm i}). \tag{67}$$

The initial state is still given by Eq. (57). If one starts from a coherent state $|\phi_i\rangle$, then Eq. (67) simplifies to

$$|\psi_{\rm f}\rangle = \int_{\phi_{\rm f}} e^{-\phi_{\rm f}} |\phi_{\rm f}\rangle \mathcal{A}'(\phi_{\rm f}|\phi_{\rm i}).$$
 (68)

III. GRAPHICAL INTERPRETATION OF THE COHERENT-STATE PATH INTEGRAL, FIRST-PASSAGE PROBABILITIES, AND RENORMALIZATION

Field theories of the type introduced above are often evaluated in perturbation theory, and interpreted graphically. To this aim, the part of the action linear in both ϕ and ϕ^* is solved explicitly, yielding a single-particle propagator or response function. One then starts with n particles, draws their trajectories, and studies how they interact via the terms nonlinear in ϕ and ϕ^* . In this process, particles may be destroyed and created. In the following, we show how to derive this picture from the coherent-state path integral, based on the shifted formulation in Eqs. (62)–(68).

A. The initial condition

Start with a general initial state $|\phi_{i,x}\rangle := e^{\int_x \phi_{i,x} \hat{a}_x^{\dagger}} |0\rangle$. Let us create p particles, at positions x_1 to x_p . In the operator picture, this is encoded by

$$\hat{a}_{x_1}^{\dagger} \dots \hat{a}_{x_p}^{\dagger} |0\rangle = \frac{\delta}{\delta \phi_{\mathbf{i}, x_1}} \dots \frac{\delta}{\delta \phi_{\mathbf{i}, x_p}} |\phi_{\mathbf{i}, x}\rangle \Big|_{\phi_{\mathbf{i}, x} = 0}.$$
 (69)

Applying the path-integral formalism developed in Secs. II J and II K, we obtain a field theory with action S', depending on the two fields ϕ and ϕ^* . In Eq. (54) we had derived the factor for the first time slice, on which we still have to shift $\phi^* \to \phi^* + 1$; this has further to be multiplied by the factor of $e^{-\int_x \phi_{i,x}}$ from the normalization; writing both factors explicitly, this results in

$$e^{-\int_{x}(\phi_{x,t_{i}}-\phi_{i,x})(\phi_{x,t_{i}}^{*}+1)} \times e^{\int_{x}-\phi_{i,x}}$$

Note the cancellation for the terms proportional to $\phi_{i,x}$; using Eq. (69), an initial condition with p particles at positions specified above is thus transferred to the path integral as

$$\hat{a}_{x_1}^{\dagger} \dots \hat{a}_{x_n}^{\dagger} |0\rangle \longrightarrow \phi_{x_1,t_1}^* \dots \phi_{x_n,t_n}^* \tag{70}$$

and

$$\phi_{i,x} = 0. \tag{71}$$

Graphically, we draw a particle emanating from position x at time t as a dot at that position in space-time, from which an arrow starts,

$$^{\uparrow}_{\bullet} x, t$$
(72)

B. The propagator

Consider diffusion as given by the Hamiltonian in Eq. (46). According to Eq. (55) the action is

$$S_0'[\phi,\phi^*] = \int_{x,t} \phi_{x,t}^* \partial_t \phi_{x,t} + D \nabla \phi_{x,t}^* \nabla \phi_{x,t}. \tag{73}$$

This yields the propagator, alias Green's or response function (in Fourier space),

Transforming back to real space, this is

$$\oint = G_{x'-x}^{t'-t} := \langle \phi_{x',t'} \phi_{x,t}^* \rangle = \Theta(t'-t) \frac{e^{-\frac{(x-x')^2}{4D(t'-t)}}}{\sqrt{4\pi D(t'-t)}}.$$
(75)

It is solution of the partial differential equation

$$(\partial_{t'} - D\nabla_{x'}^2)G_{x'-x}^{t'-t} = \delta(x - x')\delta(t - t'). \tag{76}$$

Probability is conserved, i.e., $\int_{r'} G_{r'-r}^{t'-t} = 1$.

C. The interactions

To be specific, consider the annihilation process with rate ν ,

$$A + A \xrightarrow{\nu} A. \tag{77}$$

The Hamiltonian of this process was derived in Eq. (15),

$$\mathcal{H}_{\nu}[\hat{a}^{\dagger}, a] = \frac{\nu}{2} [\hat{a}^{\dagger} \hat{a}^2 - (\hat{a}^{\dagger})^2 \hat{a}^2]. \tag{78}$$

The corresponding term in the shifted action is

$$S'_{\nu}[\phi^*, \phi] = -\sum_{x} \int_{t} \mathcal{H}_{\nu}[\phi^*_{x,t} + 1, \phi_{x,t}]$$

$$= \frac{\nu}{2} \sum_{x} \int_{t} (\phi^*_{x,t} + 1) \ \phi^*_{x,t} \phi^2_{x,t}$$

$$= \frac{\nu}{2} \sum_{x} \int_{t} +$$
(79)

Note that both terms have the same sign. Passing to the continuum yields

$$S_{\nu}'[\phi^*, \phi] = \frac{\nu \delta^d}{2} \int_{x, t} (\phi_{x, t}^* + 1) \phi_{x, t}^* \phi_{x, t}^2.$$
 (80)

Since we wish the total number of particles to be $\sum_{x} \phi_{x,t} \rightarrow \int_{x} \phi_{x,t}$, in the discrete version $\phi_{x,t}$ is the number of particles on site x, whereas in the continuum version it is the density of particles, resulting in the additional factor of δ^d in the action. Alternatively, we could keep $\phi_{x,t}$ the number of particles in a box of size δ . To avoid these problems, which are not essential for our discussion, we set $\delta \rightarrow 1$, except when specified otherwise.

D. Perturbation theory

Suppose two particles start at time t = 0 at positions x_1 and x_2 . We want to know the probability $p_1(t_f)$ to find only one of

them at time t_f . Our formalism gives the following perturbative expansion:

$$p_1(t_{\rm f}) = \nu \bigwedge_{}^{} - \nu^2 \bigvee_{}^{} + \nu^3 \bigvee_{}^{} - \dots$$
 (81)

The lower two points are fixed at time t = 0 and at positions x_1 and x_2 . The intermediate times and positions (symbolized by black dots) are integrated over. Let us call t_f the final time.

Naively, one would expect the probability to be given by a path integral, propagating one particle from x_1 at time t = 0, and the other particle from x_2 at the same time to a common position x at time t, and then propagation of a single particle to time t_f . Integrating over x, one thus naively expects that

$$p_{1}(t_{f}) \stackrel{?}{=} \nu \int_{0}^{t_{f}} dt \int_{x,y} G_{x-y}^{t_{f}-t} G_{y-x_{1}}^{t} G_{y-x_{2}}^{t}$$

$$= \nu \int_{0}^{t_{f}} dt \int_{y} G_{y-x_{1}}^{t} G_{y-x_{2}}^{t}. \tag{82}$$

This is but the first diagram in Eq. (81). The question is, where do the remaining terms come from?

E. Interpretation of the "strange" quartic vertex in terms of a first-passage problem

Let us try to construct the probability of annihilation without making reference to the formalism derived above. This is very enlightening, since it will not only re-derive the action (79), but also shed light on necessary ultraviolet cutoffs of the field theory, and their physical interpretation. This procedure is equivalent to renormalization.

Consider two particles propagating. We draw their positions $x_i(t)$ for discrete times $t = n\tau, n \in \mathbb{N}$, represented graphically as



Starting at t = 0 at positions $x_1(0) = x_1$ and $x_2(0) = x_2$, we want to know the probability $p_1(\tau)$ that the two particles meet, or more precisely are within a distance $\delta/2$ at time $t_1 = \tau$:

$$p_{1}(\tau) = \frac{1}{1 + 2}$$

$$= \int_{x,x'} G_{x-x_{1}}^{\tau} G_{x'-x_{2}}^{\tau} \Theta(|x - x'| < \delta/2). \tag{84}$$

Note that since $G(\tau, x - x_1)$ is not a probability, but a probability density, we have to say how close they have to come so that we consider them to "meet"; the probability that the two particles are exactly at the same position is actually zero. With this prescription the above expression is a probability density, as is G(t,x). To simplify our treatment, we will approximate

this by (d is the dimension)

$$p_1(\tau) = \frac{1}{1 + \frac{x}{2}} \approx \delta^d \int_x G_{x-x_1}^{\tau} G_{x-x_2}^{\tau}.$$
 (85)

Let us now calculate the probability that the two particles meet in the second time step: Particle 1 propagates from x_1 at t=0 to x' at time τ , and then to x at time $t=2\tau$. The second particle has intermediate position x''. Thus (with the same approximation as above) we obtain

$$\frac{1}{x} = \delta^d \int_{x,x',x''} G_{x-x'}^{\tau} G_{x'-x_1}^{\tau} G_{x-x''}^{\tau} G_{x''-x_2}^{\tau}.$$
(86)

Now we use that the Green's functions obey the composition property

$$\int_{x'} G_{x-x'}^{\tau} G_{x'-x_1}^{\tau} = G_{x-x_1}^{2\tau}.$$
 (87)

This allows to rewrite this contribution as

$$\frac{x}{1} = \frac{x}{1} \qquad (88)$$

However, this is not the complete result: The particles could already have met in the first time step, at time $t = \tau$. Subtracting this contribution, we have

$$p_1(2\tau) = \frac{x}{x''} - \frac{x}{x''}$$
 (89)

Note that one cannot simply subtract the probability to have met at $t = \tau$,

$$p_1(2\tau) \neq \underbrace{\frac{x}{x''}}_{x''} - \underbrace{\frac{x}{x'''}}_{x''}. \tag{90}$$

Let us now calculate the probability to meet for the first time at $t = 3\tau$,

We subtracted the configurations where the particles met at times $t = \tau$ and $t = 2\tau$; however this subtracts twice the configuration where the particles met at time $t = \tau$ and at time $t = 2\tau$, which have to be added at the end.

Note that once the two particles have met, a single one will continue propagating (not drawn here). We can therefore read

off the action which yields the above perturbation expansion,

$$S'[\phi^*, \phi] = \int_{t_i}^{t_f} dt \, dx \, \phi_{x,t}^* \partial_t \phi_{x,t} + D \nabla \phi_{x,t}^* \nabla \phi_{x,t} + \frac{\delta^d}{2} \sum_{t=n\tau} \int_x (\phi_{x,t}^* + 1) \phi_{x,t}^* \phi_{x,t}^2.$$
(92)

Converting the sum over integer n into an integral, $\sum_{t=n\tau} \rightarrow \frac{1}{\tau} \int_{t_i}^{t_f} dt$, and comparing to Eqs. (73) and (79), we identify the rate ν as

$$v = \frac{1}{\tau}.\tag{93}$$

It is now clear what the quartic vertex is doing: It converts the problem of a meeting of the two particles to the problem of a meeting for the first time, a first passage problem.

One can try to resum explicitly the perturbation series. As we set up the framework, it is well defined for small ν and finite τ . Under these circumstances, resummation is rather tedious, and the author of the present notes has decided to eliminate the corresponding calculations in order to keep the material readable.

We can, however, deduce the result in the limit of $\tau \to 0$ and $\nu \to \infty$: One first realizes that the distance between the two particles is again a random walk with a diffusion constant 2D instead of D. It can thus be described by an action

$$S_{\text{rel}}[\phi, \phi^*] = \int_{x,t} \phi_{x,t}^* \partial_t \phi_{x,t} + 2D\nabla \phi_{x,t}^* \nabla \phi_{x,t}. \tag{94}$$

Second, the field $\phi(x,t)$ is only defined for $x \ge 0$, and zero for x = 0: when the two particles meet, a single particle will propagate from that point on, and their relative position will be zero. This is known as Dirichlet boundary conditions, and can be solved with the method of images [26]. In dimension d = 1, this leads to

$$G_{x',x}^{t'-t} := \langle \phi_{x',t'} \phi_{x,t}^* \rangle = \Theta(t'-t) \frac{e^{-\frac{(x-x')^2}{8D(t'-t)}} - e^{-\frac{(x+x')^2}{8D(t'-t)}}}{\sqrt{8\pi D(t'-t)}}.$$
(95)

Here x is the difference in position at the start, and x' the distance in position at the end. Note the difference from Eq. (75). Integrating over x' from zero to infinity, we obtain the probability that the two particles did not meet up to time t, knowing that they started at distance x at time 0,

$$p_{\text{survive}}(x,t) = \text{erf}\left(\frac{x}{\sqrt{8Dt}}\right).$$
 (96)

The probability $p_1(t)$ given in Eq. (81) then is

$$p_1(t_f) = 1 - p_{\text{survive}}(x, t_f).$$
 (97)

This can be generalized to higher dimensions.

IV. STOCHASTIC EQUATION OF MOTION FOR THE COHERENT-STATE PATH INTEGRAL

A. General formulation

We established in Eq. (66) that the transition amplitude between the coherent states $|\phi_i\rangle$ and $|\phi_f\rangle$ is given by

$$\mathcal{A}'(\phi_{\rm f}|\phi_{\rm i}) = \int \mathcal{D}[\phi] \, \mathcal{D}[\phi^*] \, e^{-S'[\phi^*,\phi]} \Big|_{\phi_{\rm f}=\phi_{\rm i}}^{\phi_{\rm f}=\phi_{\rm f}}. \tag{98}$$

Note that we use the shifted action, thus shifted fields ϕ^* , since then both ϕ and ϕ^* have zero expectation values, rendering all following considerations simpler.

Suppose now that the shifted Hamiltonian and thus the shifted action have only linear and quadratic terms in ϕ_t^* ; a term independent of ϕ^* is absent due to the conservation of probability, Eq. (19),

$$\mathcal{H}'[\phi_t^*, \phi_t] = \phi_t^* \mathcal{L}[\phi_t] + \frac{1}{2} (\phi_t^*)^2 \mathcal{B}[\phi_t].$$
 (99)

First consider $\mathcal{B}[\phi_t] = 0$, i.e., only a term linear in ϕ_t^* . Then the saddle point obtained by variation with respect to ϕ^* gives the exact solution to the path integral, encoded in the equation of motion, of ϕ_t

$$\partial_t \phi_t = \mathcal{L}[\phi_t], \quad \phi_{t_i} = \phi_i,$$
 (100)

$$\mathcal{A}'(\phi_{\rm f}|\phi_{\rm i}) = \delta(\phi_{\rm f} - \phi_{t_{\rm f}}). \tag{101}$$

Quite amazingly, an explicit solution for a nonlinear path integral has been given.

This simple solution is no longer possible if $\mathcal{B}[\phi_t] \neq 0$. To nevertheless use an equation of motion, we introduce a Gaussian random variable, i.e., white noise, ξ_t , to write $e^{\frac{1}{2}\mathcal{B}[\phi_t](\phi_t^*)^2}$ as an expectation value over the noise,

$$e^{\frac{1}{2}\int_{t}\mathcal{B}[\phi_{t}](\phi_{t}^{*})^{2}} = \left\langle e^{\int_{t}\phi_{t}^{*}\sqrt{\mathcal{B}[\phi_{t}]}\xi_{t}}\right\rangle_{\varepsilon}$$
(102)

$$\langle \xi_t \xi_{t'} \rangle_{\varepsilon} = \delta(t - t'). \tag{103}$$

Note that if $\mathcal{B}[\phi_t]$ is negative, then the noise is imaginary. The sign of the root is irrelevant, since ξ_t is statistically invariant under $\xi_t \to -\xi_t$. With the noise, the equation of motion (100) changes to

$$\partial_t \phi_t = \mathcal{L}[\phi_t] + \sqrt{\mathcal{B}[\phi_t]} \xi_t, \tag{104}$$

$$\phi_{t} = \phi_{i}. \tag{105}$$

The interpretation is as follows: The transition amplitude $\mathcal{A}'(\phi_f|\phi_i)$ can be sampled by simulating the Langevin equation (104), with initial condition (105) and noise (103),

$$\mathcal{A}'(\phi_{\rm f}|\phi_{\rm i}) = \langle \delta(\phi_{\rm f} - \phi_{t_{\rm f}}) \rangle_{\xi}. \tag{106}$$

According to Eq. (62) an observable \mathcal{O} has then expectation at time t_f :

$$\langle \mathcal{O}_{t_{\rm f}} \rangle = \langle \mathcal{O}_{\rm N}(1, \phi_{t_{\rm f}}) \rangle_{\varepsilon}.$$
 (107)

This is an intuitive result, with some caveats: First, we recall the replacement of $\hat{a}^{\dagger} \rightarrow 1$. Second, $\mathcal{O}_{N}(\hat{a}^{\dagger},\hat{a})$ is the normal-ordered version of the operator, e.g., is $\hat{n}^{2} = (\hat{a}^{\dagger}\hat{a})^{2} = (\hat{a}^{\dagger})^{2}\hat{a}^{2} + \hat{a}^{\dagger}\hat{a}$, so that $\langle \hat{n}_{t_{f}}^{2} \rangle = \langle \phi_{t}^{2} + \phi_{t} \rangle_{\xi}$; see Eq. (33).

In Appendix B 3 we give a formal proof of this relation, based uniquely on the CSPI. The formalism produces two terms, a linear term proportional to $\frac{\delta}{\delta\hat{a}_x}\mathcal{O}(1,\hat{a})$, and a quadratic term proportional to $\frac{\delta}{\delta\hat{a}_x}\frac{\delta}{\delta\hat{a}_y}\mathcal{O}(1,\hat{a})$. These two terms can then be interpreted as drift and diffusion terms in the Itô formalism. This gives an independent derivation of the process (104) and the relation (107).

If $\sqrt{\mathcal{B}[\phi_t]}$ is real, one may think of equation (104) as describing what is "going on" in the system. This is not

the case if $\mathcal{B}[\phi_t]$ is negative, thus $\sqrt{\mathcal{B}[\phi_t]}$ purely imaginary: Then generically, states sampled by the path integral are "nonphysical" in the sense that they do not correspond to a probability density, even though the transition amplitude is given by Eq. (106). We will come back to this question in Sec. IV G: There we will show that in the case of imaginary noise, the formalism works for short times, but breaks down for longer times.

In Sec. V we will propose a different physically motivated treatment, leading to a coarse-grained effective stochastic equation of motion with a real noise.

B. Example: Diffusion

Let us start with simple diffusion, with Hamiltonian

$$\mathcal{H}'[\hat{a}^{\dagger}, \hat{a}] := \mathcal{H}[\hat{a}^{\dagger} + 1, \hat{a}] = -D \int_{x} \nabla \hat{a}_{x}^{\dagger} \nabla \hat{a}_{x}. \tag{108}$$

Note that the shift has no effect since only $\nabla \hat{a}_x^{\dagger}$ appears. This implies the action, already given in Eq. (73),

$$S'[\phi^*, \phi] = \int_{x,t} \phi_{x,t}^* \partial_t \phi_{x,t} + D \nabla \phi_{x,t}^* \nabla \phi_{x,t}. \tag{109}$$

Variation with respect to ϕ^* leads to the equation of motion

$$\partial_t \phi_{x,t} = D \nabla^2 \phi_{x,t}. \tag{110}$$

This is a very simple and actually quite remarkable equation: While diffusion is a noisy process, leading to fluctuations of the number of particles on a given site, Eq. (110) is a an exact, noiseless equation. It tells us how the distribution of the number of particles on a given site evolves with time.

As a test, let us check that it keeps the total-particle-number distribution fixed. Equation (43) implies that at, say $t = t_i$, the total-particle-number distribution is given by the coherent state

$$|\Phi\rangle = \left| \int_{r} \phi_{x,t_{i}} \right\rangle. \tag{111}$$

Since for periodic boundary conditions $\int_x \partial_t \phi_{x,t} = D \int_x \nabla^2 \phi_{x,t} = 0$, the state $|\Phi\rangle$ does not change over time. In particular, this implies particle-number conservation.

C. Example: Reaction diffusion

Consider the reaction-diffusion process $A + A \xrightarrow{\nu} A$ with (shifted) action defined by Eqs. (73) and (79):

$$S'[\phi^*, \phi] = \int_{x,t} \left\{ \phi_{x,t}^* \partial_t \phi_{x,t} + D \nabla \phi_{x,t}^* \nabla \phi_{x,t} + \frac{\nu}{2} \left[\phi_{x,t}^* \phi_{x,t}^2 + (\phi_{x,t}^*)^2 \phi_{x,t}^2 \right] \right\}.$$
(112)

The corresponding equations of motion and noise are

$$\partial_t \phi_{x,t} = -\frac{\nu}{2} \phi_{x,t}^2 + D \nabla^2 \phi_{x,t} + i \sqrt{\nu} \phi_{x,t} \xi_{x,t},$$
 (113)

$$\langle \xi_{x,t} \xi_{x',t'} \rangle = \delta(t - t') \delta(x - x'). \tag{114}$$

This noise is imaginary. It has puzzled many researchers whether this is unavoidable [11,16,17,19], or could even be beneficial [18]. We will come back to this question later.

D. Dual formulation: Equation of motion for ϕ_t^*

Note that one can define the dual process of Eq. (104) by exchanging in the dynamical action the roles of ϕ and ϕ^* : Suppose the Hamiltonian can be written in the form

$$\mathcal{H}[\phi_t^*, \phi_t] = \mathcal{L}^*[\phi_t^*]\phi_t + \frac{1}{2}\mathcal{B}^*[\phi_t^*](\phi_t)^2. \tag{115}$$

The path integral for the generating function at time $t_{\rm f}$ then becomes

$$\rho_{\mathbf{f}}(\phi^*) := \langle e^{\phi^* \phi_{t_{\mathbf{f}}}} \rangle
= \int \mathcal{D}[\phi] \mathcal{D}[\phi^*] e^{\phi^* \phi_{t_{\mathbf{f}}} - \int_{t} \phi_{t}^* \partial_{t} \phi_{t} - \mathcal{H}[\phi_{t}^*, \phi_{t}]} \Big|_{\phi_{t_{\mathbf{i}}} = \phi_{\mathbf{i}}}^{\phi^* = \phi_{t_{\mathbf{f}}}^*} .$$
(116)

Note that this equation is written in terms of the unshifted Hamiltonian. Contrary to Eq. (61) it is normalized, since the leftmost state is not $\langle 0|e^a=\langle 1|$. Integrating $\int_t \phi_t^* \partial_t \phi_t$ by part, and noting that the boundary term changes in the exponential $\phi^*\phi_{t_f} \to \phi_t^*\phi_i$, yields

$$\rho_{\rm f}(\phi^*) = \int \mathcal{D}[\phi] \, \mathcal{D}[\phi^*] e^{\phi_{\rm f}^* \phi_{\rm i} + \int_t \phi_t \, \partial_t \phi_t^* + \mathcal{H}[\phi_t^*, \phi_t]} \Big|_{\phi_{\rm f} = \phi_{\rm i}}^{\phi^* = \phi_{\rm f}^*}. \tag{117}$$

This path integral is sampled by the stochastic process

$$-\partial_t \phi_t^* = \mathcal{L}^*[\phi_t^*] + \sqrt{\mathcal{B}^*[\phi_t^*]} \xi_t, \tag{118}$$

$$\phi_{t_c}^* = \phi^*. \tag{119}$$

It evolves the (dual) state ϕ_t^* from t_f to t_i , backward in time, as is suggested by the sign of Eq. (118).

Consider now $\mathcal{B}^* \equiv 0$, such that the evolution becomes deterministic, $-\partial_t \phi_t^* = \mathcal{L}^*[\phi_t^*]$. Denote by $\Psi_{t,t_{\rm f}}: \phi^* = \phi_{\rm f}^* \to \phi_t^*$ this time evolution, i.e.,

$$\phi_t^* = \Psi_{t,t_f}(\phi^*). \tag{120}$$

Note that $\Psi_{t,t_f}(0) = 0$ and $\Psi_{t,t_f}(1) = 1$.

As a concrete example, consider the branching process, including a possible annihilation $A \rightarrow 0$:

$$A \xrightarrow{\lambda_n} nA. \tag{121}$$

Then

$$\mathcal{H}[\hat{a}^{\dagger}, \hat{a}] = \sum_{n} \lambda_{n} [(\hat{a}^{\dagger})^{n} \hat{a} - \hat{a}^{\dagger} \hat{a}]$$

$$\equiv f(\hat{a}^{\dagger}) \hat{a} - f(1) \hat{a}^{\dagger} \hat{a}, \qquad (122)$$

where we defined $f(x) := \sum_{n} \lambda_{n} x^{n}$. The equation of motion (118) then becomes (there is no Doi shift)

$$-\partial_t \phi_t^* = f(\phi_t^*) - f(1)\phi_t^*, \quad \phi_{t_t}^* = \phi^*. \tag{123}$$

To be explicit, choose $\lambda_2=1$, and all other $\lambda_i=0$. We have to solve (backward in time) the equation $\partial_t \phi_t^* = \phi_t^* - (\phi_t^*)^2$. It has solution $\phi_t^* = 1/[1 - e^{t_t - t}(1 - 1/\phi^*)]$. The function Ψ then reads

$$\Psi_{t_f,t}(x) = \frac{x}{x + e^{t_f - t}(1 - x)}. (124)$$

Using Eq. (117), this yields the generating function, evaluated at $t = t_i$,

$$\rho_{f}(\phi^{*}) = \rho_{i}(\Psi_{t_{f},t_{i}}(\phi^{*})). \tag{125}$$

This is a classical result; see, e.g., [9,27]. Suppose one starts with a single particle at time t=0; then $\rho_i(\phi^*)=\phi^*$, and the above becomes

$$\rho_{\rm f}(\phi^*) = \frac{\phi^*}{\phi^* + e^{t_{\rm f} - t_{\rm i}} (1 - \phi^*)}.$$
 (126)

The probability to have n particles at time t_f is given by the nth series coefficient, namely

$$p_n(t_f) = e^{t_i - t_f} (1 - e^{t_i - t_f})^{n-1}, \quad n \geqslant 1.$$
 (127)

This is a rather simple expression.

We could also try to solve the problem by varying with respect to ϕ^* , inducing the stochastic equation of motion

$$\partial_t \phi_t = \phi_t + i\sqrt{2\phi_t}\xi_t. \tag{128}$$

This equation talks about the evolution of the state $|\phi_t\rangle$, which will become complex. We will discuss in the next section how this can be interpreted. Compared to the latter approach, the solution (126) is much more elegant and explicit.

E. Testing the coherent-state path integral

Consider the annihilation equation (77),

$$A + A \xrightarrow{\nu} A,$$
 (129)

with stochastic equation of motion (113). For simplicity, we concentrate on a single site,²

$$\partial_t \phi_t = -\frac{v}{2} \phi_t^2 + i \sqrt{v} \phi_t \xi_t, \tag{130}$$

$$\langle \xi_t \xi_{t'} \rangle = \delta(t - t'). \tag{131}$$

Let us use as initial distribution coherent state $|\phi_i\rangle$, i.e., a Poisson distribution with parameter ϕ_i ,

$$p_{i}(n) = e^{-\phi_{i}} \frac{\phi_{i}^{n}}{n!}.$$
 (132)

There are two ways to study this process.

F. Direct simulation of the reaction process

Let us start by directly simulating the reaction process (129): First, use the probability distribution (132) to obtain an integer n (the occupation number at $t = t_i$), and then evolve Eq. (129) for a time $T = t_f - t_i$. The latter is best done by remarking that if at a given time t there are n particles, the probability that they have not decayed up to time $t + \delta t$ (with arbitrary δt) is

$$p_n^{\text{survive}}(\delta t) = \exp\left(-\frac{n(n-1)}{2}\nu\delta t\right).$$
 (133)

Thus one can draw a random number $r_n \in [0,1]$, sampling the decay probability; solving $r_n = p_n^{\text{survive}}(\delta t)$ for δt then yields

$$\delta t_n := -\frac{2}{n(n-1)} \ln r_n, \quad t_n := \sum_{i \ge n} \delta t_i, \quad t_1 := \infty.$$
 (134)

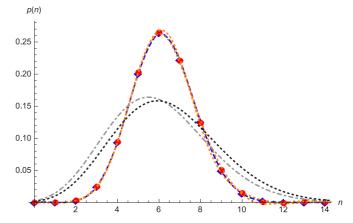


FIG. 1. Results for the process (129), using $a_i = 15$, $t_f - t_i = 0.2$, and $\nu = 1$. Blue diamonds: Direct numerical simulation with 5×10^5 samples (partially hidden behind red dots). The blue dashed line is a guide for the eye. The statistical error bars are smaller than this linewidth. Red dots: Integration of the stochastic equation of motion, using Re[$\mathcal{P}_f^{cs}(n)$] from Eq. (136), for 5×10^4 samples, $\delta t = 10^{-3}$. The orange dotted lines are Re[$\mathcal{P}_f^{cs}(n)$] \pm |Im[$\mathcal{P}_f^{cs}(n)$]|, defined for all real n, which is an estimate of the error. Both simulations ran about 100 s. Within these errors, the agreement is excellent. The dot-dashed gray line is a Poisson distribution with $a_f = 6$, which would be the result of Eq. (130) in the absence of noise. Taking into account the drift term $va_t/2$ induced by the noise in Eq. (129) (see Appendix C) leads to a = 6.43, (black, dotted). The real distribution is centered around this value, but is much narrower than a Poisson distribution.

The t_n (with n decreasing) are a series of times when one of the n particles decays. Thus the number n_f of particles at t_f is given by

$$n_{\rm f}: t_{n_{\rm f}} \leqslant t_{\rm f} - t_{\rm i} < t_{n_{\rm f}-1}.$$
 (135)

Repeating this procedure, one obtains a histogram for the final number of particles, and an associated normalized probability distribution $p_f^{DS}(n)$, where DS stands for "direct simulation." The result, for $\phi_i = 15$, $t_f - t_i = 0.2$, and $\nu = 1$, is presented in Fig. 1 (blue diamonds, partially hidden behind the red circles).

Alternatively, one can numerically integrate the master equation (11) (with $\mu = \kappa = 0$) using $p_i(n)$ given by Eq. (132) for $n \le n_{\text{max}} = 3\phi$, and setting $p(n) \to 0$ for $n > n_{\text{max}}$. This is the fastest and most precise solution.

G. Integration of the stochastic equation of motion (130)

Integrating the stochastic equation of motion (130) for different noise realizations ξ_t , and with initial condition $\phi_{t_i} = \phi_i$, one obtains a (complex) result for $\phi_f = \phi_{t_f}$. One then measures the final distribution, as an average over all noise realizations,

$$p_{\rm f}^{\rm SEM}(n) := \left\langle e^{-\phi_{\rm f}} \frac{\phi_{\rm f}^n}{n!} \right\rangle_{\xi}. \tag{136}$$

The result is again shown in Fig. 1 (red circles). One sees several things: First, the agreement between the direct numerical simulation of the decay process (blue diamonds, and blue dashed line as guide for the eye) and the stochastic equation of motion (red dots, and orange dashed lines, with

²A complementary study was performed in [18] for the process $A + A \leftrightarrows 0$.

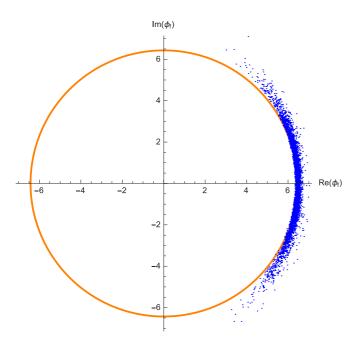


FIG. 2. 5000 samples for the result of the integration of Eq. (130), with $\nu = 1, t_{\rm f} - t_{\rm i} = 0.2, \phi_{\rm i} = 15$. These samples lie approximately on a circle of size $\phi_{\rm f} = 6.43$. For larger times, the samples extend further around the circle.

error estimate) is quite good. This confirms that Eq. (136) is indeed applicable, even though the final states ϕ_f are complex.

Second, the final distribution is much narrower than a Poisson distribution: both the Poissonian obtained for $\phi = 6$ [gray dashed line, result of integrating Eq. (130), dropping the noise term], plotted in Fig. 1, or the one including a drift term $\nu \phi_t/2$ in Eq. (130) (black dashed line; see Appendix C for discussion). Having a distribution narrower than a Poissonian is possible only with imaginary noise, which leads to a diffusion of the phase of ϕ_t ; see Fig. 2. In contrast, real noise leads to a widening of the distribution. We study this in more detail below.

Third, using the stochastic equation of motion has its limits: Indeed, already for $t_f - t_i = 0.5$, the stochastic equation of motion gets appreciable error bars, even with a large number s of samples, and for $t_f - t_i = 1$ convergence is no longer assured. We tried an improved algorithm as follows: Instead of starting s "particles" at $\phi(t_i) = \phi_i$, and evolving them until time t_f , whenever one of these particles gets too large a phase (which promises to give a large value in $e^{-\phi_f}$), we "split" the particle in two, each of which carries half of the weight (the original weight is w = 1/s) of its "father." If the phase is still too large, we split it again, propagating two particles with half the weight each. This procedure is repeated recursively. We have not been able to find parameters to improve the precision at constant execution time. We suspect that when splitting points, it becomes more probable that "bad regions" are reached, and while the weight of the corresponding points is reduced, the probability that they appear is increased. This is illustrated in Fig. 3. We also note that for the toy model studied in Appendix C 2 (pure phase) this problem is present. This indicates that the convergence problem is severe, and no algorithm to overcome it has been found yet. We refer

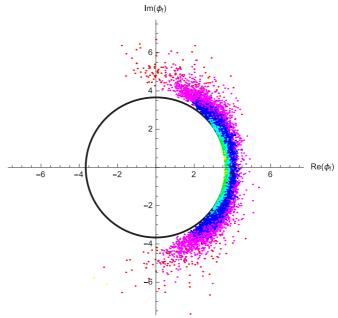


FIG. 3. Result of the integration of Eq. (130), with $\nu = 1$, $t_f - t_i = 0.5$, $\phi_i = 15$. The black circle has radius $\phi_f = 3.6614$, obtained by integrating the drift term $\partial_t \phi_t = -\phi_t^2 + \phi_t + t/2$. The color codes the number of splittings, from yellow over green, cyan, blue, magenta to red. (Thus a red point has 2^{-5} times the weight of a yellow point.)

the reader to [18,19,28] for a more detailed discussion of the problems and partially successful attempts at their solution.

H. Integrating a stochastic equation of motion with multiplicative noise: Real vs imaginary noise

Let us integrate the following stochastic equations of motion:

$$\partial_t a_t = \xi_t a_t - \frac{1}{2} a_t, \tag{137}$$

$$\partial_t a_t = i\xi_t' a_t + \frac{1}{2}a_t. \tag{138}$$

They are constructed such that $\langle |a_t| \rangle$ does not change with time; see Appendix C. In Fig. 4 one sees that, as expected, a real noise leads to a broadening of the distribution (green dots), whereas an imaginary noise leads to a narrowing of the latter (blue diamonds). For imaginary noise, the phase portrait looks similar to Fig. 2. As one can easily observe in a numerical simulation, this leads to problems for larger times, since then a_f can become imaginary, and both a_f^n and the factor of e^{-a_f} will lead to strong interference between the different samples, and to large fluctuations of the such obtained averages.

Analytic solutions for both processes are given in Appendices C 1 (real noise) and C 2 (imaginary noise).

I. Integrating a stochastic equation of motion with multiplicative noise: "Canceling" real and imaginary noises

It is instructive to consider an equation of motion with two noises,

$$\partial_t a_t = \xi_t a_t + i \xi_t' a_t. \tag{139}$$

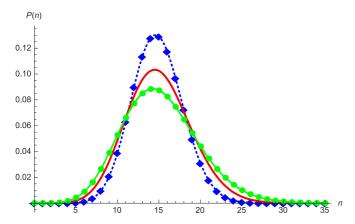


FIG. 4. Result for $p_{\rm f}^{\rm CGSEM}(n)$ in Eq. (136), after integrating Eqs. (137) (real noise, green circles) and (138) (imaginary noise, blue diamonds) for $a_{\rm i}=15$, $t_{\rm f}-t_{\rm i}=0.025$. One sees that real noise leads to a broadening of the distribution, while imaginary noise leads to a narrowing.

Writing the effective action and averaging over the noises ξ_t and ξ_t' yields an exact cancelation of the terms generated in the dynamic action: $\frac{1}{2} \int_t [a_t^* a_t]^2$ generated from the average over ξ_t cancels with $-\frac{1}{2} \int_t [a_t^* a_t]^2$ generated from the average over ξ_t' . Nevertheless, our equation of motion does not vanish. This is possible only since the coherent states are overcomplete; i.e., we do not need the states $|a\rangle$ with complex a to code all possible states. This might allow us to define a projection algorithm, which eliminates the states with complex a.

Let us check this cancellation. To do so write

$$a_t = a_0 e^{\phi_t}; \tag{140}$$

this implies

$$\partial_t \phi_t = \xi_t + i \xi_t'. \tag{141}$$

For details see Appendix C. Note that there is no drift term (as compared to the purely real or purely imaginary cases). The probability to find $\phi = \phi_x + i\phi_y$ at time t is given by the diffusion propagator,

$$P_t(\phi,\phi^*)d\phi_x d\phi_y = \frac{e^{-\frac{\phi\phi^*}{2t}}}{2\pi t}d\phi_x d\phi_y.$$
 (142)

This implies

$$P_{t}(a, a^{*})da_{x}da_{y} = \left| \frac{\partial(\phi_{x}, \phi_{y})}{\partial(a_{x}, a_{y})} \right| P_{t}(\phi, \phi^{*})da_{x}da_{y}$$

$$\simeq \frac{1}{2\pi a a^{*}t} \exp\left(-\frac{\ln\left(\frac{a}{a_{0}}\right)\ln\left(\frac{a^{*}}{a_{0}}\right)}{2t}\right) da_{x}da_{y}.$$
(143)

The approximation is due to the fact that the larger ϕ_y in Eq. (142) are not summed over; for pedagogical reasons we content ourselves with this approximation. Let us now take Eq. (143), and try to integrate over a. Writing $a = \frac{1}{2} \left(\frac{1}{2} \right) \left(\frac{1}{2} \right)$

$$(x+iy)a_{i}, a^{*} = (x-iy)a_{i}, \text{ we get}$$

$$\int P_{t}(a,a^{*})e^{(a_{x}+ia_{y})\hat{a}^{\dagger}}da_{x}da_{y}$$

$$\simeq \int \frac{dx\,dy}{2\pi t}\,e^{(x+iy)a_{i}\hat{a}^{\dagger}}e^{-\frac{(x-1)^{2}+y^{2}}{2t}}$$

$$= e^{a_{i}\hat{a}^{\dagger}}.$$
(144)

This result says that real and imaginary noise have canceled; thus the superposition of all states gives back the original coherent state. This is of course what one expects, knowing that the two terms in the effective action cancel.

V. ALTERNATIVE STOCHASTIC MODELING: AN EQUATION OF MOTION WITH REAL NOISE

A. Stochastic noise as a consequence of the discreteness of the states

In Sec. IV F, we have simulated directly the random process $A + A \xrightarrow{\nu} A$. Each simulation run gave one possible realization of the process, in the form of an integer-valued monotonically decreasing function n(t). Averaging over these runs, one samples the final distribution $P_f(n)$, or, equivalently, moments of n_f . Let us now start with a fixed number of particles instead of a coherent state, $n(t_i) = n_i$. We then want to ask the question: Is there a continuous random process $\hat{n}(t)$ which has the same statistics as n(t)?

Let us consider a little more general problem: Let n_t be the number of particles at time t. With rate r_+ the number of particles increases by one, and with rate r_- it decreases by one. This implies that after one time step, as long as $r_\pm \delta t$ are small,

$$\langle n_{t+\delta t} - n_t \rangle = (r_+ - r_-)\delta t, \tag{145}$$

$$\langle (n_{t+\delta t} - n_t)^2 \rangle = (r_+ + r_-)\delta t. \tag{146}$$

The following continuous random process \hat{n}_t has the same first two moments as n_t ,

$$d\hat{n}_t = (r_+ - r_-)dt + \sqrt{r_+ + r_-} \,\xi_t dt,\tag{147}$$

$$\langle \xi_t \xi_{t'} \rangle = \delta(t - t'). \tag{148}$$

This procedure can be modified to include higher cumulants of $n_{t+\delta t} - n_t$, leading to more complicated noise correlations. Results along these lines were obtained in Ref. [29] by considering cumulants generated in the effective field theory.

B. Example: The reaction-annihilation process

In the case of the reaction-annihilation process, the rate $r_+ = 0$, and $r_- = \frac{\nu}{2} \hat{n}_t (\hat{n}_t - 1)$; the latter, in principle, is only defined on integer \hat{n}_t , but we will use it for all \hat{n}_t . Thus the best we can do to replace the discrete stochastic process with a continuous one is to write

$$\frac{d\hat{n}_t}{dt} = -\frac{\nu}{2}\hat{n}_t(\hat{n}_t - 1) + \sqrt{\frac{\nu}{2}\hat{n}_t(\hat{n}_t - 1)}\,\xi_t. \tag{149}$$

Using $n_i = 15$ and $\nu = 1$, we have shown two typical trajectories in Fig. 5, one for the process n_t (red, with jumps), and one for the process \hat{n}_t (blue-gray, rough). While by construction

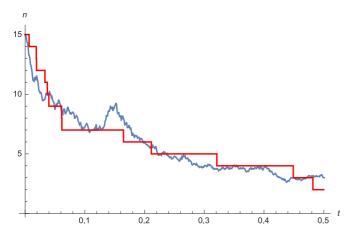


FIG. 5. One trajectory each for process n_t , i.e., a direct numerical simulation of $A + A \rightarrow A$ (red, with jumps) and \hat{n}_t , Eq. (149) (bluegray, continuous, rough). The rate is $\nu = 1$. We have chosen two trajectories which look "similar." Note that \hat{n}_t is not monotonically decreasing.

both processes have (almost) the same first two moments, clearly \hat{n}_t looks different: It is continuous, which n_t is not, and it can increase in time, which n_t cannot. One can also compare the distribution for $t_f - t_i = 0.5$; see Fig. 6. While the distribution of n_f is discrete (blue diamonds), that for \hat{n}_f is continuous (cyan). Rounding n_f to the nearest integer gives a different distribution (red). We have also drawn (black lines) the size of the boxes which would produce p(n) from $p(\hat{n})$. Clearly, there are differences. On the other hand, it is also evident that these differences will diminish when increasing n_i .

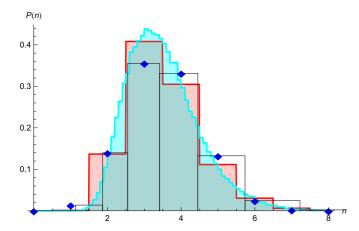


FIG. 6. Result of a numerical simulation, starting with $n_i = 15$ particles, and evolving for $t_f - t_i = 0.025$. Blue diamonds: Direct numerical simulation of the process $A + A \rightarrow A$ with rate $\nu = 1$. Cyan: Distribution of the continuous random walk (149). Red: The latter distribution, when rounding n_f to the nearest integer. Black boxes: The size of the boxes in the n direction to obtain the result of the direct numerical simulation of the process $A + A \rightarrow A$. Both processes have first moment 3.511 ± 0.001 , and second connected moment 1 ± 0.05 ; the third connected moments already differ quite substantially, 0.75 versus 0.2.

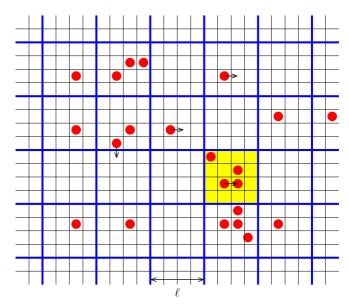


FIG. 7. A coarse-grained lattice with box size $\ell = 4$. The yellow box contains n = 4 particles.

C. Diffusion

We now derive the effective stochastic equation of motion for diffusion, i.e., hopping of grains from site i to site $i \pm 1$ with rate D. We cannot directly write an equation of the form (147) for the particle number \hat{n}_i on site i, since it does not respect the conservation of the number of particles. The latter is realized by introducing the current $J_{i+\frac{1}{2},i}$: A positive current $J_{i+\frac{1}{2},i} = 1$ represents a particle hopping from site i to i+1. A negative current $J_{i+\frac{1}{2},i} = -1$ corresponds to a particle hopping from site i+1 to site i. Each hopping has a rate D; the rate for a given particle to leave a site is the coordination number 2d times D. We thus arrive at the rate equations (with the hat again denoting the variables of the continuous process)

$$d\hat{n}_{i,t} = \left(\hat{J}_{i-\frac{1}{2},t} - \hat{J}_{i+\frac{1}{2},t}\right) dt, \tag{150}$$

$$\hat{J}_{i+\frac{1}{2},t} = D(\hat{n}_{i,t} - \hat{n}_{i+1,t}) + \sqrt{D(\hat{n}_{i,t} + \hat{n}_{i+1,t})} \, \eta_{i+\frac{1}{2},t}.$$
(151)

The white noise has correlations

$$\langle \eta_{i+\frac{1}{2},t} \rangle = 0, \quad \langle \eta_{i+\frac{1}{2},t} \eta_{j+\frac{1}{2},t'} \rangle = \delta_{i,j} \delta(t-t').$$
 (152)

To perform the continuum limit, let us introduce the density of particles inside a box $B_{\ell}(x)$ centered at x and of linear size ℓ , as well as the (d-dimensional) current,

$$\rho(x,t) := \sum_{i \in B_{\ell}(x)} \frac{\hat{n}_{i,t}}{\ell^d}, \quad J(x,t) := \sum_{i \in B_{\ell}(x)} \frac{\hat{J}_{i+1/2,t}}{\ell^d}, \quad (153)$$

which (in first approximation) is independent of the size of the box (see Fig. 7). (We dropped the hat for convenience of notation.) In terms of ρ , the stochastic equations become³

³These equations are standard, undisputed, and appear frequently in the literature; see, e.g., [30]. They are a special case of Eq. (17) of [31], itself equivalent to Eq. (4) of [17].

(generalized to d dimensions)

$$\partial_t \rho(x,t) = -\nabla \vec{J}(x,t), \tag{154}$$

$$\vec{J}(x,t) = -D\nabla \rho(x,t) + \sqrt{2D\rho(x,t)}\,\vec{\eta}(x,t),\tag{155}$$

$$\langle \eta^i(x,t)\eta^j(x',t')\rangle = \delta^{ij}\delta^d(x-x')\delta(t-t'). \tag{156}$$

Note that there is no ℓ -dependent factor either for the current or the noise term. Combining the first two equations yields

$$\partial_t \rho(x,t) = D\nabla^2 \rho(x,t) + \nabla[\sqrt{2D\rho(x,t)}\vec{\eta}(x,t)]. \quad (157)$$

Let us step back and analyze the above findings; for simplicity of notation we again set d=1. First of all, the diffusion process is constructed such that particles do not interact. A given particle will be on a chosen site with probability 1/L, where L is the system size. If $N=\bar{n}L$ is the total number of particles, and \bar{n} the mean particle number per site, then the probability to find n particles on a given site is

$$p(n) = {N \choose n} \left(\frac{1}{L}\right)^n \left(1 - \frac{1}{L}\right)^{N-n} \simeq e^{-\bar{n}} \frac{(\bar{n})^n}{n!}.$$
 (158)

The last relation is valid in the limit of L large. We recuperate our old friend, the normalized coherent state $e^{-\phi}|\phi\rangle$, with $\phi = \bar{n}$. Note that in the CSPI, the analog of Eq. (157) is given by Eq. (110), namely

$$\partial_t \phi(x,t) = D \nabla^2 \phi(x,t). \tag{159}$$

Since the CSPI works with coherent states, it does not need the noise of Eq. (157). What diffuses is the "weight" $\phi(x,t)$ of the coherent state, which is the mean particle number per site, termed \bar{n} above. Thus in the CSPI, both mean and variance of the number of grains on a site tends to \bar{n} . We checked with a numerical simulation that Eq. (157) indeed leads to a distribution of grains per site with mean and variance \bar{n} .

D. An effective stochastic field theory of the reaction process

Let us now construct an effective field theory of the reaction-diffusion process. Consider a lattice of size L^d , with particles on it, which can hop from one site to a neighboring one with rate D. To simplify our considerations, let us suppose that we take the limit of $v \to \infty$: if a particle jumps on an occupied site, only one of them survives. To construct an effective field theory, we introduce boxes of size ℓ . Each of these boxes contains n(x,t) particles at time t.

With rate D, a particle hops. Thus the probability with which a particle in a given box will hop is $D n_{x,t} \delta t$; that it will land on an occupied site and thus annihilate is

$$\delta_{\rm an} n(x,t) \simeq D \, n(x,t) \delta t \times \frac{n(x,t)-1}{\ell^d}.$$
 (160)

The second factor is an approximation which neglects the correlations inside the box. If the particle had hopped out of its box, then the second factor should involve the density in the neighboring box; writing the density in the same box is another approximation. Last but not least, if the box is sufficiently large, then one can replace $n(x,t) - 1 \rightarrow n(x,t)$.

To perform the continuum limit, we use the density $\rho(x,t)$ defined in Eq. (153). The equation of motion of this density

then becomes

$$\partial_t \rho(x,t) = -D\rho(x,t)^2 + \sqrt{D}\rho(x,t)\xi(x,t) + D\nabla^2 \rho(x,t) + \nabla[\sqrt{2D}\rho(x,t)\vec{\eta}(x,t)], \quad (161)$$

$$\langle \xi(x,t)\xi(x',t')\rangle = \delta(t-t')\delta^d(x-x'),\tag{162}$$

$$\langle \eta^i(x,t)\eta^j(x',t')\rangle = \delta^{ij}\delta(t-t')\delta^d(x-x'), \tag{163}$$

$$\langle \xi(x,t) \rangle = \langle \eta^i(x,t) \rangle = \langle \xi(x,t) \eta^i(x',t') \rangle = 0. \tag{164}$$

Note that all factors of δt and ℓ have disappeared, absorbed into a nontrivial dimension of $\xi(x,t)$ and $\vec{\eta}(x,t)$. Note that the diffusive noise is usually dropped. The reason is that the coarse-grained density $\rho(x,t)$ varies smoothly, thus

$$\nabla[\sqrt{2D\,\rho(x,t)}\vec{\eta}(x,t)] \simeq \sqrt{2D\,\rho(x,t)}\nabla\vec{\eta}(x,t). \tag{165}$$

Integrating the letter over a box of size ℓ will yield $\eta(x,t)$ on the boundary, making it less relevant by a factor of $1/\ell$. It is customarily dropped as subdominant. Thus the effective stochastic description of the annihilation-diffusion process is

$$\partial_t \rho(x,t) = -D\rho(x,t)^2 + \sqrt{D}\rho(x,t)\xi(x,t) + D\nabla^2 \rho(x,t). \tag{166}$$

E. Comparison between the coherent-state path integral and the effective coarse-grained stochastic equation of motion

Let us rewrite the equations of motion for the effective field theory for the annihilation-diffusion process, first in coherent states, and second in the coarse-grained stochastic-equationof-motion formalism, choosing similar conventions. First, the stochastic equation of motion for the coherent-state formalism reads

$$\partial_t \phi_{x,t} = -\frac{\nu}{2} \phi_{x,t}^2 + i \sqrt{\nu} \phi_{x,t} \, \xi_{x,t} + D \nabla^2 \phi_{x,t}, \tag{167}$$

$$\langle \xi_{x,t} \rangle = 0, \quad \langle \xi_{x,t} \xi_{x',t'} \rangle = \delta(t - t') \delta^d(x - x'). \tag{168}$$

Second, the coarse-grained stochastic equation of motion reads (we dropped the noise from diffusion)

$$\partial_t \rho_{x,t} = -\frac{\nu}{2} \rho_{x,t}^2 + \sqrt{\frac{\nu}{2}} \rho_{x,t} \xi_{x,t} + D \nabla^2 \rho_{x,t}, \qquad (169)$$

$$\langle \xi_{x,t} \rangle = 0, \quad \langle \xi_{x,t} \xi_{x',t'} \rangle = \delta(t - t') \delta^d(x - x').$$
 (170)

Let us summarize our findings, based on what we have done so far:

- (1) Both equations look rather similar.
- (2) In the formulation with a coherent state $\phi_{x,t}$, the noise $\xi_{x,t}$ is imaginary. This indicates that the distribution becomes *narrower* than a coherent state.
- (3) In the formulation with the effective density $\rho_{x,t}$, which starts with a sharp distribution, the latter *widens* by the presence of the real noise $\xi_{x,t}$.
- (4) The noises are both proportional to $\sqrt{\nu}$, the rate, times the state variable; they differ by a factor of $i\sqrt{2}$.
- (5) In the coherent-state formulation, perturbation theory can be interpreted in terms of particle trajectories and first-meeting probabilities; see Sec. III. This interpretation is not possible for the coarse-grained stochastic equation of motion.

(6) The coarse-grained stochastic equation of motion in principle has an additional noise term; see the last term of Eq. (161). This term becomes relevant if we coarse-grain with very small boxes, but is irrelevant for large boxes: Since it is a total derivative, its integral over a volume element is a boundary (surface) term, down by a factor of $1/\ell$, with ℓ the box size.

F. Other approaches

As we saw above, the appearance of an imaginary, or more generally complex, noise and its physical interpretation are puzzling. This is reflected in the literature; see, e.g., [11,16–19]. One basic reason is that in the CSPI the variables are coherent states, i.e., (rather broad) distributions instead of sharp δ distributions. The second reason is that they are microscopic, not coarse-grained, variables. We then discussed a physically motivated approach based on coarse-grained variables. In the field-theoretic context, these questions were considered in [29], and more specifically for branching annihilation in [32] (see also [11], Sec. 9.2).

An alternative approach is to search a change of variables for the coherent states. A beautiful proposal was made in Ref. [17]. This approach can be interpreted as rewriting creation and annihilation operators of the CSPI on each site as

$$\hat{a}^{\dagger} = e^{\hat{\rho}^{\dagger}}, \quad \hat{a} = e^{-\hat{\rho}^{\dagger}}\hat{\rho}.$$
 (171)

The operator $\hat{\rho}$ is the particle-number operator (and not a particle annihilation operator as \hat{a}). A derivation of the basic equations and its consequences are discussed in Appendix F. As can be seen from the transformation (171), if the process does not respect particle-number conservation, we obtain additional factors of $e^{\pm \hat{\rho}^{\dagger}}$; these terms are nonlinear in $\hat{\rho}^{\dagger}$, and cannot simply be decoupled with the help of a Gaussian noise. The situation may be different if particle numbers are conserved. We come back to this question in Sec. VIF.

VI. A PHENOMENOLOGICAL DERIVATION OF THE STOCHASTIC FIELD THEORY FOR THE MANNA MODEL

In this section, we apply our considerations to a nontrivial example, the stochastic Manna model. We will see that our formalism permits a systematic derivation of the effective stochastic equations of motion. While the result is known in the literature [20–23], it was there derived by symmetry principles, which are not always convincing. Furthermore, they leave undetermined all coefficients. While many of them can be eliminated by rescaling, our derivation will "land" on a particular line of parameter space, characterized by the absence of additional memory terms; see Sec. VID.

A. Basic definitions

The Manna sandpile was introduced in 1991 by Manna [4] as a stochastic version of the Bak-Tang-Wiesenfeld (BTW) sandpile [3]. It is defined as follows.

Manna model. Randomly throw grains on a lattice. If the height at one point is greater or equal to two, then with rate 1 move two grains from this site to randomly chosen neighboring sites. Both grains may end up on the same site.

We start by analyzing the phase diagram. We denote by a_i the fraction of sites with i grains. It satisfies the sum rule⁴

$$\sum_{i} a_i = 1. \tag{172}$$

In these variables, the number of grains n per site can be written as

$$n := \sum_{i} a_i i. \tag{173}$$

The empty sites are

$$e := a_0.$$
 (174)

The fraction of active sites is

$$a := \sum_{i \ge 2} a_i. \tag{175}$$

We also define the (weighted) activity as

$$\rho := \sum_{i \geqslant 2} a_i (i - 1). \tag{176}$$

Note that ρ satisfies the sum rule

$$n - \rho + e = 1. \tag{177}$$

In order to take full advantage of this definition, one may change the toppling rules of the Manna model to those of the weighted Manna model.

Weighted Manna model. If a site contains $i \ge 2$ grains, randomly move these grains to neighboring sites with rate (i-1).

In Fig. 8 (thick lines), we show a numerical simulation of the Manna model in a 2-dimensional system of size $L \times L$, with L = 150. There is a phase transition at $n = n_c = 0.702$. Close to n_c , the fraction of doubly occupied sites a_2 grows linearly with $n - n_c$, and higher occupancy is small. Indeed, we checked numerically that for $n > n_c$ the probability p_i to find i grains on a site decays exponentially with i, i.e., $p_i \sim \exp(-\alpha_n i)$, where α_n depends on n; see Fig. 9. This is to be contrasted with the initial condition, where we randomly distribute $n \times L^2$ grains on the lattice of size $L \times L$, and which yields a Poisson distribution, the coherent state $|n\rangle$, for the number of grains on each site; see inset of Fig. 9 (left). This result suggests that coherent states may not be the best representation for this system. It further implies that close to the transition, $\rho \approx a$, and we expect that the weighted Manna model and the original Manna model have the same critical behavior. We come back to this question below.

B. MF solution

In order to make analytical progress, we now study the *topple-away* or mean-field solution of the stochastic Manna sandpile, which we can solve analytically.

Mean-field Manna model. If a site contains two or more grains, move these grains to any randomly chosen other site of the system.

⁴Note that a_i has nothing to do with the operator \hat{a}_i used earlier.

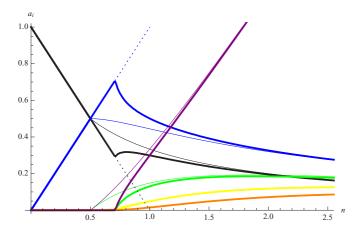


FIG. 8. Thick lines: The order parameters of the Manna model, as a function of n, the average number of grains per site, obtained from a numerical simulation of the stochastic Manna model on a grid of size 150×150 with periodic boundary conditions. We randomly update a site for 10^7 iterations, and then update the histogram 500 times every 10^5 iterations. Plotted are the fractions of sites that are unoccupied (black), singly occupied (blue), double occupied (green), triple occupied (yellow), and quadruple occupied (orange). The activity $\rho = \sum_{i>1} a_i (i-1)$ is plotted in purple. No data were calculated for n < 0.5, where $a_0 = e = 1 - n$, $a_1 = n$, and $a_{i>2} = 0$ (inactive phase). Note that before the transition, $a_0 = 1 - n$ and $a_1 = n$. The transition is at $n = n_c = 0.702$. Thin lines: The MF phase diagram, as given by Eqs. (181) and following for $n \leq \frac{1}{2}$, and by Eqs. (184) and following for $n \geq \frac{1}{2}$. We checked the latter with a direct numerical simulation.

The rate equations are, setting for convenience $a_{-1} := 0$,

$$\partial_t a_i = -a_i \Theta(i \ge 2) + a_{i+2} + 2 \left[\sum_{j \ge 2} a_j \right] (a_{i-1} - a_i).$$
(178)

They can be rewritten as

$$\partial_t a_i = -a_i \Theta(i \geqslant 2) + a_{i+2} + 2(1 - a_0 - a_1)(a_{i-1} - a_i).$$
(179)

We are interested in the steady state $\partial_t a_i = 0$. One can solve these equations by introducing a generating function. An alternative solution consists in realizing that for $i \ge 2$, Eq. (179) admits a steady-state solution of the form

$$a_i = a_2 \kappa^{i-2}, \quad i > 2.$$
 (180)

This reduces the number of independent equations $\partial_t a_i = 0$ in Eq. (179) from infinity to three. Furthermore, there are the equations $\sum_{i=0}^{\infty} a_i = 1$ and $\sum_{i=0}^{\infty} i \ a_i = n$. Thus there are 5 equations for the 4 variables a_0, a_1, a_2 , and κ . The reason we apparently have one redundant equation is due to the fact that we already used the normalization condition (172) to go from Eq. (178) to Eq. (179).

These equations have two solutions: For 0 < n < 1, there is always the solution for the inactive or absorbing state,

$$a_0 = 1 - n, (181)$$

$$a_1 = n, (182)$$

$$a_{i\geqslant 2}=0. (183)$$

For n > 1/2, there is a second nontrivial solution:

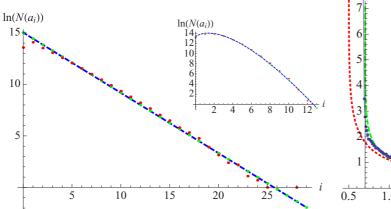
$$a_0 = \frac{1}{1+2n},\tag{184}$$

$$a_{i>0} = \frac{4n\left(\frac{2n-1}{2n+1}\right)^i}{4n^2 - 1}.$$
 (185)

(Note that a_2/a_1 has the same geometric progression as a_{i+1}/a_i for i > 2, which we did not suppose in our ansatz.) Thus the probability to find i > 0 grains on a site is given by the exponential distribution

$$p(i) = \frac{4n}{4n^2 - 1} \exp(-i\alpha_n), \quad \alpha_n = \log\left(\frac{2n + 1}{2n - 1}\right).$$
 (186)

Using these two solutions, we get the mean-field (MF) phase diagram plotted in Fig. 8 (thin lines). This has to be compared



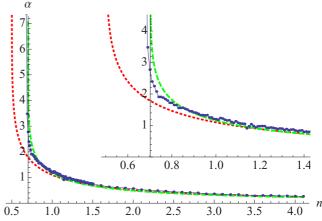


FIG. 9. Left: (Unnormalized) histogram after many topplings for n=2; the probability that a site has i grains decays as $e^{-0.585i}$, for all $i \ge 1$. Inset: The initial distribution, a Poissonian. Right: The exponential decay coefficient α as a function of n. The dots are from a numerical simulation. The dashed red line is the MF result (214). The green dashed line is a fit corresponding to $\alpha \approx \frac{2}{3} \ln[(n+n_c)/(n-n_c)]$. Inset: Blowup of main plot.

with the simulation of the Manna model on the same figure (thick lines). One sees that for $n \ge 2$, the MF solution and simulation are becoming almost indistinguishable. We have also checked with simulations that the Manna model has a similar exponentially decaying distribution of grains per site, with a decay constant α plotted on the right of Fig. 9.

A similar MF analysis can be performed for the weighted Manna model and the Abelian sandpile model (ASM); this is discussed in Appendix D.

There is a series of models which interpolates between the Manna model and its MF version: the range-r Manna model, where grains are not deposited on a random neighbor, but on any site within a distance r. In Appendix E it is discussed how this model converges for large r to the MF Manna model.

C. The complete effective equations of motion for the Manna model

In this section, we will give the effective equations of motion for the Manna model. Let us start from the mean-field equations for $\rho(t)$ and n(t). For simplicity of expressions, we use the weighted Manna model. The physics close to the transition should not depend on it. Let us start from the hierarchy of MF equations (D1), similar to Eq. (179) for the Manna model, and which can be rewritten as

$$\partial_t a_i = (1 - i)a_i \Theta(i \ge 2) + (i + 1)a_{i+2} + 2\rho(a_{i-1} - a_i).$$
(187)

For convenience, let us write explicitly the rate equation for the fraction of empty sites $e \equiv a_0$,

$$\partial_t e = a_2 - 2\rho e. \tag{188}$$

The first term, the gain $r_+ = a_2$, comes from the sites with two grains, toppling away, and leaving an empty site. The second term, the loss term, is the rate at which one of the toppling grains lands on an empty site, $r_- = 2\rho e$.

We now follow the formalism developed in Sec. VA, Eqs. (145)–(148). This yields

$$\partial_t e = \rho (1 - 2e) + \sqrt{a_2 + 2\rho e} \,\bar{\xi}_t,$$
 (189)

where $\langle \bar{\xi}_t \bar{\xi}_{t'} \rangle = \delta(t - t')/l^d$, and l is the size of the box which we consider. Now remark that close to the transition, $a_2 \approx \rho$. Inserting this into the above equation, we arrive at

$$\partial_t e \approx \rho (1 - 2e) + \sqrt{\rho} \sqrt{1 + 2e} \,\bar{\xi}_t.$$
 (190)

Due to Eq. (177), the combination $n - \rho + e = 1$, and since n is conserved this implies $\partial_t e \equiv \partial_t \rho$. It is customary to write Eq. (190) for $\partial_t \rho$, instead of $\partial_t e$. Next we approximate $\sqrt{1 + 2e}$ by the value of e at the transition, i.e., $e \rightarrow e_c^{\text{MF}} = \frac{1}{2}$; see the mean-field phase diagram in Fig. 8. We thus arrive at

$$\partial_t \rho \approx (2n-1)\rho - 2\rho^2 + \sqrt{2\rho}\,\bar{\xi}_t.$$
 (191)

Note that this equation gives back $n_c^{\rm MF}=\frac{1}{2}$, and as a consequence of the conservation law $n-\rho+e=1$ also $e_{\rm c}^{\rm MF}=\frac{1}{2}$, used above in the simplification of the noise term.

Finally, let us suppose we have not a single box of size ℓ , but a lattice of boxes, labeled by a d-dimensional label x. Each toppling event moves two grains from a site to the neighboring

sites, equivalent to a current

$$J(x,t) = -D\nabla \rho(x,t) + \sqrt{2D\rho(x,t)}\xi(x,t)$$
 (192)

with diffusion constant $D = 2 \times \frac{1}{2d} = \frac{1}{d}$. The first factor of 2 is due to the fact that two grains topple. The factor of $\frac{1}{2d}$ is due to the fact that each grain can topple in any of the 2d directions; thus the rate D per direction is $\frac{1}{2d}$, resulting in D = 1/d. As discussed above, we will drop the noise term as subdominant.

This current changes both the activity $\rho(x,t)$, as the number of grains n(x,t), resulting in $\partial_t \rho(x,t) = \partial_t n(x,t) = -\nabla J(x,t)$. It does not couple to the density of empty sites. Using the sum rule (177) $n - \rho + e = 1$, implies the consistency relation $\partial_t \rho(x,t) \equiv \partial_t n(x,t) + \partial_t e(x,t)$ for the current; this confirms that both $\rho(x,t)$ and n(x,t) must couple to the same current.

Thus, we finally arrive at the following set of equations:

$$\partial_{t}\rho(x,t) = \frac{1}{d}\nabla^{2}\rho(x,t) + [2n(x,t) - 1]\rho(x,t) - 2\rho(x,t)^{2} + \sqrt{2\rho(x,t)}\,\xi(x,t),$$
(193)
$$\partial_{t}n(x,t) = \frac{1}{d}\nabla^{2}\rho(x,t),$$
(194)

$$\langle \xi(x,t)\xi(x',t')\rangle = \delta^d(x-x')\delta(t-t'). \tag{195}$$

This is known as the conserved directed percolation class. Instead of writing coupled equations for $\rho(x,t)$ and n(x,t), we can also write coupled equations for e(x,t) and $\rho(x,t)$:

$$\partial_t e(x,t) = [1 - 2e(x,t)]\rho(x,t) + \sqrt{2\rho(x,t)}\,\xi(x,t),$$
 (196)

$$\partial_t \rho(x,t) = \frac{1}{d} \nabla^2 \rho(x,t) + \partial_t e(x,t). \tag{197}$$

The above equations for ρ and n were obtained in the literature [20–23] by means of symmetry principles, but never properly derived. Evoking symmetry principles also leaves all coefficients undefined, and does not ensure that Eq. (196) is valid on a single site, i.e., is free of spatial derivatives. This locality will prove essential in the next section.

D. Excursion: Mapping to disordered elastic manifolds

In [24] it had been proposed to use these equations as a basis for mapping the effective field theory of the Manna model derived above onto driven disordered elastic systems. The identifications are

$$\rho(x,t) = \partial_t u(x,t)$$
 the velocity of the interface, (198)

$$e(x,t) = \mathcal{F}(x,t)$$
 the force acting on the interface. (199)

The second equation (197) is the time derivative of the equation of motion of an interface, subject to a random force $\mathcal{F}(x,t)$,

$$\partial_t u(x,t) = \frac{1}{d} \nabla^2 u(x,t) + \mathcal{F}(x,t). \tag{200}$$

Since $\rho(x,t)$ is positive, u(x,t) is for each x monotonically increasing. Instead of parametrizing $\mathcal{F}(x,t)$ by space x and time t, it can be written as a function of space x and interface position u(x,t). Setting $\mathcal{F}(x,t) \to F(x,u(x,t))$, the

first equation (196) becomes

$$\partial_{t} \mathcal{F}(x,t) \to \partial_{t} F(x,u(x,t))$$

$$= \partial_{u} F(x,u(x,t)) \partial_{t} u(x,t)$$

$$= [1 - 2F(x,u(x,t))] \partial_{t} u(x,t)$$

$$+ \sqrt{2 \partial_{t} u(x,t)} \xi(x,t). \tag{201}$$

For each x, this equation is equivalent to the Ornstein-Uhlenbeck [33] process F(x,u), defined by

$$\partial_u F(x, u) = 1 - 2F(x, u) + \sqrt{2} \,\xi(x, u),$$
 (202)

$$\langle \xi(x, u)\xi(x', u')\rangle = \delta^d(x - x')\delta(u - u'). \tag{203}$$

It is a Gaussian Markovian process with mean $\langle F(x,u) \rangle = 1/2$, and variance in the steady state of

$$\langle [F(x,u) - \frac{1}{2}][F(x',u') - \frac{1}{2}] \rangle = \frac{1}{2} \delta^d(x' - x') e^{-2|u - u'|}.$$
(204)

Writing the equation of motion (200) as

$$\partial_t u(x,t) = \frac{1}{d} \nabla^2 u(x,t) + F(x,u(x,t)), \qquad (205)$$

it can be interpreted as the motion of an interface with position u(x,t), subject to a disorder force F(x,u(x,t)). The latter is δ -correlated in the x direction, and short-ranged correlated in the u direction. In other words, this is a disordered elastic manifold subject to random-field disorder. It can be treated via field theory. The latter relies on functional RG (see [34] for an introduction) for the renormalized version of the force-force correlator (204). Functional RG is nowadays well developed, and predicts not only a plethora of critical exponents [35] but also size, velocity, and duration distributions [36], as well as the shape of avalanches [37].

Also note that Eq. (193) has a quite peculiar symmetry, namely the factor of 2 in front of both $n(x,t)\rho(x,t)$ and $-\rho(x,t)^2$. As a consequence, Eq. (196) does not contain a term $\sim \rho^2(x,t)$, which would spoil the simple mapping presented above. The absence of this term cannot be induced on symmetry arguments only. How this additional term, if present, can be treated is discussed in Ref. [24].

E. Coherent states for the Manna model?

In the last section, we derived an effective field theory for the stochastic Manna model, in the coarse-grained stochastic-equation-of-motion formalism (CGSEM). The reader might wonder why we did not try to use the CSPI formalism. Well, we tried, and here we will share the rather disappointing outcome: One starts from the discrete Hamiltonian, with rate 1, and for simplicity in dimension d=1

$$\mathcal{H}_{\text{Manna}}[\hat{a}^{\dagger}, \hat{a}] = \sum_{i} \left[\frac{1}{4} (\hat{a}_{i+1}^{\dagger} + \hat{a}_{i-1}^{\dagger})^{2} - (\hat{a}_{i}^{\dagger})^{2} \right] \hat{a}_{i}^{2}. \quad (206)$$

The first term, proportional to $(\hat{a}_{i+1}^{\dagger} + \hat{a}_{i-1}^{\dagger})^2 \hat{a}_i^2$, checks whether there are two or more grains on site i, and then moves two grains to randomly chosen neighbors. The last term $-(\hat{a}_i^{\dagger})^2 \hat{a}_i^2$ is responsible for the conservation of probability. One can show that this is equivalent to the stochastic equation

of motion

$$\partial_t a_{t,j} = \nabla^2 a_{t,j}^2 + \frac{1}{\sqrt{2}} (\eta_{t,j-1} a_{t,j-1} + \eta_{t,j+1} a_{t,j+1}) + \sqrt{2} i \xi_{t,j} a_{t,j}$$
(207)

with noise $\langle \eta_{t,j} \eta_{t',j'} \rangle = \langle \xi_{t,j} \xi_{t',j'} \rangle = \delta(t-t') \delta_{j,j'}$ and $\langle \eta_{t,j} \xi_{t',j'} \rangle = 0$. Note that the first (diffusive) term comes from the shift $\hat{a}_i^\dagger \to \hat{a}_i^\dagger + 1$, to be done before decoupling the path integral with noise terms.

This equation is quite ugly: Not only does it have two multiplicative noises; one of them is imaginary, inducing the convergence problems mentioned in Sec. IV G.

As we do not know how to proceed from here, let us try to address this question on a more abstract level: Could coherent states indeed help us? Let us also state our prerogative, namely that one wants to derive an effective field theory for a local field. First of all, one has to realize that what the coherent state does is to rewrite the number of grains on a given site as a superposition of Poisson distributions (also known as coherent states) on this site. This complicates matters; after all, we want an effective variable which gives us some intuition.

Let us take a step back and look at the derivation of the field theory of the Ising model. One realizes that there one needs coarse graining. What we have done in the preceding section was to construct coarse-grained effective variables for a block of $\ell \times \ell$ sites. As for Ising, this necessitates making approximations inside a block. We did this, supposing that we are close to the transition, and that we can use the MF approximation for the block.

But maybe this block can be described by a coherent state? From Fig. 9 we see that the distribution inside a block is, at least approximately, an exponential, represented by (we normalized)

$$|a_0,\beta\rangle = a_0|0\rangle + \frac{(1-a_0)(1-\beta)}{1-\beta\hat{a}^{\dagger}}\hat{a}^{\dagger}|0\rangle.$$
 (208)

This implies $n = \frac{1-a_0}{1-\beta}$ and $\rho = \frac{1-a_0}{1-\beta}\beta$; in the example of Fig. 9, $a_0 \approx 0.195$, $\beta \approx e^{-0.585}$, n = 2, and $\rho \approx 1.115$. This has to be contrasted to a coherent state, with on average \bar{n} particles (also normalized),

$$|\bar{n}\rangle = e^{\bar{n}(\hat{a}^{\dagger} - 1)}|0\rangle. \tag{209}$$

This state has $a_0 = e^{-\bar{n}}$, possesses \bar{n} particles, and activity $\rho = \bar{n} + e^{-\bar{n}} - 1$. There are two problems: First of all, the tail is different; this should not be an issue close to the transition, where triple and higher occupancy are unimportant. The second and more fundamental problem is that while the state (208) has two independent parameters, the coherent state (209) has only one. Still, the coherent-state functional integral will correctly propagate a state. More precisely, it will calculate a transition probability from an initial state to a final state, after decomposition into a coherent-state representation. It will do so by passing through complex intermediate states. The imaginary part of these complex variables provides the second, "missing" variable. Thus thinking of a single (real) coherent state per site is not appropriate.

One could also try to work with coherent states for the number of empty, once, twice, or triple,... occupied sites

inside a box. This would yield a state of the form

$$|a_0, a_1, a_2, \dots\rangle := e^{\ell^2 (a_0 \hat{a}_0^{\dagger} + a_1 \hat{a}_1^{\dagger} + a_2 \hat{a}_2^{\dagger} + \dots)} |0\rangle,$$
 (210)

where the \hat{a}_i^{\dagger} create an i times occupied site. If $\sum_i a_i = 1$, the expectation of the number of sites will still be ℓ^2 , but it will be a fluctuating variable, with variance ℓ . While this is probably acceptable, the author does not see what would be gained in terms of simplicity of derivation, knowing that all the approximations necessary for the MF treatment of Sec. VI B would have to be made as well.

Coherent states as a basis for a stochastic description of the Manna model have also been proposed by Pastor-Satorras [20], based on the field theory of Wijland-Oerding-Hilhorst [38]. The idea there was to introduce two species A and B, where B particles are associated with the activity, and can diffuse with rate D, whereas A particles are stationary. The particle number in the Manna model is associated with the total number of A and B particles. The rate equations proposed to mimic the Manna model are $B \xrightarrow{k_1} A$ and $A + B \xrightarrow{k_2} B + B$. This yields a Hamiltonian

$$\mathcal{H}[\hat{a}^{\dagger}, \hat{a}, \hat{b}^{\dagger}, \hat{b}] = \int_{x} D\hat{b}^{\dagger} \nabla^{2} \hat{b} + k_{1} (\hat{a}^{\dagger} - \hat{b}^{\dagger}) \hat{b}$$
$$+ k_{2} [(\hat{b}^{\dagger})^{2} \hat{b} \hat{a} - \hat{b}^{\dagger} \hat{a}^{\dagger} \hat{b} \hat{a}]. \tag{211}$$

This theory was then analyzed in terms of shifted fields, see Eq. (1.15) of [38], which somehow obscures the analysis. What was not realized is that the proposed stochastic interpretation necessitates an imaginary noise. This can be seen from the fact that the terms quadratic in \hat{a}^{\dagger} and \hat{b}^{\dagger} are of the form

$$\mathcal{H}[\hat{a}^{\dagger}, \hat{a}, \hat{b}^{\dagger}, \hat{b}] = k_2 \int_{x} (\hat{a}^{\dagger}, \hat{b}^{\dagger}) \begin{pmatrix} 0 & -\frac{1}{2} \\ -\frac{1}{2} & 1 \end{pmatrix} \begin{pmatrix} \hat{a}^{\dagger} \\ \hat{b}^{\dagger} \end{pmatrix} \hat{a}\hat{b}$$
+ linear terms in $\hat{a}^{\dagger}, \hat{b}^{\dagger}$. (212)

The matrix has both a positive eigenvalue corresponding to a real noise as well as a negative eigenvalue corresponding to an imaginary noise, as is the case for Eq. (207). From our above considerations this is not surprising. It shows once more that the CSPI formalism does not yield stochastic equations of motion with a purely real noise. Remark that the equations of Sec. VIC are recovered within the same approximations, when applying the CGSEM formalism developed here to the two-species model.

Finally, let us mention another peculiarity of the Manna Hamiltonian in Eq. (206): The rate for a site with $n \ge 2$ grains to topple is proportional to n(n-1); this is dictated by the demand to have a Hamiltonian as simple as possible. The original Manna model has a constant rate, while the weighted Manna model we defined above has a rate proportional to n-1. While this should be no problem at the transition, it imposes a specific rate not present in the original formulation, a rate found in pairing (or meeting) probabilities of n particles in a box. The latter is indeed the framework in which the combinatorics of the coherent-state path integral is appropriate.

F. A canonical transformation for the Manna model?

Some efforts were spent to find exact representations of a stochastic process, without using the CSPI formalism. The hope was that such a reformulation could be interpreted as a stochastic process with real noise. In Sec. VF we discuss, and in Appendix F we rederive, the proposal of Ref. [17]. This approach can be interpreted as rewriting creation and annihilation operators of the CSPI on each site as

$$\hat{a}^{\dagger} = e^{\hat{\rho}^{\dagger}}, \quad \hat{a} = e^{-\hat{\rho}^{\dagger}}\hat{\rho}.$$
 (213)

Applying this transformation to the Manna Hamiltonian (206) vields

$$\mathcal{H}_{\text{Manna}}^{\rho}[\hat{\rho}^{\dagger}, \hat{\rho}] = \frac{1}{4} \int_{r} (e^{\hat{\rho}_{i+1}^{\dagger} - \hat{\rho}_{i}^{\dagger}} + e^{\hat{\rho}_{i-1}^{\dagger} - \hat{\rho}_{i}^{\dagger}})^{2} \hat{\rho}_{i}(\hat{\rho}_{i} - 1). \tag{214}$$

Taking the continuum limit, and dropping higher-order terms in the lattice cutoff, one arrives at

$$\mathcal{H}_{\text{Manna}}^{\rho,\text{cont}}[\hat{\rho}^{\dagger},\hat{\rho}] \simeq \int_{x} \left\{ [\nabla \hat{\rho}^{\dagger}(x)]^{2} + \nabla^{2} \hat{\rho}^{\dagger}(x) \right\} \hat{\rho}(x) [\hat{\rho}(x) - 1].$$
(215)

The action $S_{\text{Manna}}^{\rho} = \int_{x,t} \rho^*(x,t) \partial_t \rho(x,t) - \mathcal{H}[\rho^*,\rho]$ allows for an interpretation as a stochastic equation of motion with a real noise $\eta^i(x,t)$,

$$\partial_t \rho(x,t) \simeq \nabla^2 \{ \rho(x,t) [\rho(x,t) - 1] \}$$

$$+ \sqrt{2} \nabla [\vec{\eta}(x,t) \sqrt{\rho(x,t) [\rho(x,t) - 1]}], \quad (216)$$

$$\langle \eta^i(x,t)\eta^j(x',t')\rangle = \delta^{ij}\delta^d(x-x')\delta(t-t'). \tag{217}$$

This equation is very similar to the linear diffusion equation (157), except that on the right-hand side the particle number ρ has been replaced by $\rho(\rho - 1)$.

The question remains whether passing from Eq. (214) to (215) is justified. In the path integral, $\hat{\rho}$ is the number of particles. As a discrete number, $\hat{\rho}$ strongly fluctuates between nearest neighbors, and one expects $\hat{\rho}^{\dagger}$ to do the same. Thus the approximation from Eq. (214) to (215) is probably not justified. One should first construct coarse-grained variables, which would probably lead to a stochastic equation of motion different from Eq. (216). We leave exploitation of these ideas for future research.

VII. CONCLUSION

In this article, we started with the coherent-state path integral (CSPI) for stochastic systems, which we then reformulated as a stochastic equation of motion. We showed how the evolution of the probability distribution can be followed, despite the appearance of imaginary noise. Limitations of this formalism were discussed, especially its (at least practical) breakdown at finite times. We also showed how some of the appearing vertices can be interpreted as transforming a simple diffusion probability into a first-meeting probability.

We then constructed a complementary formalism, based on an effective coarse-grained stochastic equation of motion (CGSEM) for a continuous variable. Demanding that drift and variance for the underlying discrete system be correctly reproduced by the CGSEM fixes the latter continuous process uniquely.

We should stress again that while both the CSPI and the CGSEM formalism share some common features, they should not be confounded: It is tempting to derive stochastic equations of motion in the CSPI formalism, and then to interpret the coherent state $|\phi\rangle$, i.e., a Poisson distribution with expectation ϕ , by the state ϕ itself, equivalent to a δ distribution at ϕ , as is used in the CGSEM formalism [20]. We remarked on the example of the reaction process $A+A\to A$, that starting from a Poisson distribution, the probability distribution becomes narrower than a Poissonian, which in the CSPI is only possible with complex coherent states, thus imaginary noise in the equation of motion. On the other hand, the probability distribution becomes broader than a δ distribution, necessitating a real noise in the CGSEM. Both noises have, up to a factor of $\sqrt{2}$, the same strength.

We concluded our considerations by analyzing the stochastic Manna model, and gave a straightforward derivation of its effective stochastic field theory. Our procedure, based on coarse graining, fixes all amplitudes, including the noise strength. Compared to earlier derivations, this derivation is simple and transparent; that it fixes all constants is an additional advantage. It also ensures the simplest mapping on disordered elastic manifolds.

We hope that our work helps to clarify the origin and interpretation of stochastic field theories, and that the techniques presented here are more broadly useful.

Note added in proof. Recently we found that rate equations similar to Eqs. (178) and following were proposed in the literature for a closely related process, the conserved threshold transfer process; see [12], page 229, recursively citing [39], [40], and [41]. We have found an application to neither the Manna model nor the solution (185) and following.

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APPENDIX A: FERMIONIC COHERENT-STATE PATH INTEGRAL

The coherent-state path integral can also be applied to fermionic degrees of freedom, i.e., to states which can only be unoccupied, or simply occupied. This can be done via the coherent-state path integral with Grassmann numbers. Here we only write down the basic equations.

We introduce fermion creation and annihilation operators with

$$\{c, c^{\dagger}\} = 1, \quad \{c, c\} = \{c^{\dagger}, c^{\dagger}\} = 0.$$
 (A1)

The states are

$$|1\rangle = c^{\dagger}|0\rangle, \quad c|0\rangle = 0.$$
 (A2)

This implies

$$\langle 1|1\rangle = \langle 0|cc^{\dagger}|0\rangle = \langle 0|\{c,c^{\dagger}\}|0\rangle = \langle 0|0\rangle = 1. \tag{A3}$$

Coherent states

$$|\psi\rangle = e^{\psi c^{\dagger}}|0\rangle = (1 + \psi c^{\dagger})|0\rangle,$$
 (A4)

$$\langle \psi | = \langle 0 | e^{c\psi^*} = \langle 0 | (1 + c\psi^*),$$
 (A5)

$$\langle \psi^* | \psi \rangle = 1 + \psi^* \psi. \tag{A6}$$

Resolution of unity

$$\int d\psi^* d\psi |\psi\rangle\langle\psi^*|e^{-\psi^*\psi} = 1.$$
 (A7)

This relation can be checked by applying it to $|0\rangle$ and to $\psi'|1\rangle$.

APPENDIX B: PROOFS OF SOME RELATIONS USED IN THE MAIN TEXT

1. Proof of Eq. (28)

Equation (28) reads

$$e^{\lambda \hat{n}} \equiv e^{\lambda \hat{a}^{\dagger} \hat{a}} = :e^{(e^{\lambda} - 1)\hat{a}^{\dagger} \hat{a}} : .$$
 (B1)

The proof of this equation consists of two steps: Applying the left-hand side of Eq. (B1) to a coherent state yields

$$e^{\lambda \hat{a}^{\dagger} \hat{a}} e^{\phi \hat{a}^{\dagger}} |0\rangle = \sum_{n=0}^{\infty} e^{\lambda \hat{a}^{\dagger} \hat{a}} \frac{\phi^{n} (\hat{a}^{\dagger})^{n}}{n!} |0\rangle$$

$$= \sum_{n=0}^{\infty} e^{\lambda n} \frac{\phi^{n} (\hat{a}^{\dagger})^{n}}{n!} |0\rangle$$

$$= e^{e^{\lambda \hat{a}^{\dagger}} |0\rangle}.$$
 (B2)

Thus

$$\langle \phi^* | e^{\lambda \hat{a}^{\dagger} \hat{a}} | \phi \rangle = \langle 0 | e^{\phi^* \hat{a}} e^{e^{\lambda} \phi \hat{a}^{\dagger}} | 0 \rangle = e^{e^{\lambda} \phi^* \phi}. \tag{B3}$$

On the other hand,

$$\begin{aligned} \langle \phi^* | : e^{(e^{\lambda} - 1)\hat{a}^{\dagger}\hat{a}} : | \phi \rangle \\ &= \sum_{n=0}^{\infty} \langle 0 | e^{\phi^* \hat{a}} (e^{\lambda} - 1)^n \frac{(\hat{a}^{\dagger})^n \hat{a}^n}{n!} e^{\phi \hat{a}^{\dagger}} | 0 \rangle \\ &= \sum_{n=0}^{\infty} \langle 0 | e^{\phi^* \hat{a}} (e^{\lambda} - 1)^n \frac{(\phi^* \phi)^n}{n!} e^{\phi \hat{a}^{\dagger}} | 0 \rangle = e^{e^{\lambda} \phi^* \phi}. \end{aligned} \tag{B4}$$

This proves relation (B1).

2. A different derivation of Eq. (35)

The set of relations (35) and following can also be derived from

$$\phi^{p} = \phi^{p} \langle 0 | e^{\hat{a}} | \phi \rangle e^{-\phi} = \langle 0 | e^{\hat{a}} \hat{a}^{p} e^{\phi \hat{a}^{\dagger}} | 0 \rangle e^{-\phi}
= \langle 0 | (\hat{a}^{\dagger} + 1)^{p} e^{\hat{a}} \hat{a}^{p} e^{\phi \hat{a}^{\dagger}} | 0 \rangle e^{-\phi}
= \langle 0 | e^{\hat{a}} (\hat{a}^{\dagger})^{p} \hat{a}^{p} e^{\phi \hat{a}^{\dagger}} | 0 \rangle e^{-\phi}
= \langle 0 | e^{\hat{a}} (\hat{n} - p + 1) \dots (\hat{n} - 1) \hat{n} e^{\phi \hat{a}^{\dagger}} | 0 \rangle e^{-\phi}.$$
(B5)

3. Formal derivation of the evolution of expectation values in the coherent-state path integral

Consider the expectation value (60) at time t_f :

$$\langle \mathcal{O}_{t_{\rm f}} \rangle = e^{-\phi_{\rm i}} \langle 0 | e^{\hat{a}} \mathcal{O}(\hat{a}^{\dagger}, \hat{a}) \mathbf{T} e^{\int_{t_{\rm i}}^{t_{\rm f}} dt \, \mathcal{H}[\hat{a}_{t}^{\dagger}, \hat{a}_{t}]} | \phi_{\rm i} \rangle. \tag{B6}$$

We are interested in its temporal evolution. At a slightly smaller time, the observable \mathcal{O} had expectation

$$\langle \mathcal{O}_{t_f - \delta t} \rangle = e^{-\phi_i} \langle 0 | e^{\hat{a}} e^{\delta t \mathcal{H}[\hat{a}^{\dagger}, \hat{a}]} \mathcal{O}(\hat{a}^{\dagger}, \hat{a}) \mathbf{T} e^{\int_{t_i}^{t_f - \delta t} dt \, \mathcal{H}[\hat{a}_t^{\dagger}, \hat{a}_t]} | \phi_i \rangle. \tag{B7}$$

Note that we have been able to add the factor $e^{\delta t \mathcal{H}[\hat{a}^{\dagger}, \hat{a}]} \triangleq e^{\delta t \mathcal{H}[\hat{a}^{\dagger}_{i_{\rm f}}, \hat{a}_{i_{\rm f}}]}$ since $\mathcal{H}[\hat{a}^{\dagger}, \hat{a}]$, when applied to the left to $\langle 0|e^{\hat{a}}$ vanishes; see Sec. II B. Thus the time derivative of the expectation value of an operator is given by its commutator with $\mathcal{H}[\hat{a}^{\dagger}, a]$,

$$\frac{d}{dt_{\rm f}}\langle \mathcal{O}_{t_{\rm f}}\rangle = \langle [\mathcal{O}(\hat{a}^{\dagger}, \hat{a}), \mathcal{H}[\hat{a}^{\dagger}, \hat{a}]]\rangle, \tag{B8}$$

where the expectation is as in Eq. (B6). To simplify the calculations, we show that if $\mathcal{O}[\hat{a}^{\dagger},\hat{a}]$ is normal-ordered, $\mathcal{O}[\hat{a}^{\dagger},\hat{a}] \equiv \mathcal{O}_{N}[\hat{a}^{\dagger},\hat{a}]$, then $\mathcal{O}[\hat{a}^{\dagger},\hat{a}]$ can be replaced by $\mathcal{O}[1,\hat{a}]$; indeed, Eq. (B8) is proportional to

$$\langle 0|e^{\hat{a}}\{\mathcal{O}(\hat{a}^{\dagger},\hat{a})\mathcal{H}[\hat{a}^{\dagger},\hat{a}] - \mathcal{H}[\hat{a}^{\dagger},\hat{a}]\mathcal{O}(\hat{a}^{\dagger},\hat{a})\}\dots$$

$$= \langle 0|e^{\hat{a}}\{\mathcal{O}(1,\hat{a})\mathcal{H}[\hat{a}^{\dagger},\hat{a}] - \mathcal{H}[\hat{a}^{\dagger},\hat{a}]\mathcal{O}(1,\hat{a})\}\dots$$
(B9)

since $\langle 0|e^{\hat{a}}\mathcal{O}(\hat{a}^{\dagger},\hat{a})=\langle 0|e^{\hat{a}}\mathcal{O}(1,\hat{a}),$ and $\langle 0|e^{\hat{a}}\mathcal{H}[\hat{a}^{\dagger},\hat{a}]=0.$ Thus

$$\frac{d}{dt_{\epsilon}} \langle \mathcal{O}_{t_{\epsilon}} \rangle = \langle [\mathcal{O}(1, \hat{a}), \mathcal{H}[\hat{a}^{\dagger}, \hat{a}]] \rangle. \tag{B10}$$

Next, the commutator can be calculated by remarking that

$$[\mathcal{O}(1,\hat{a}),\mathcal{H}[\hat{a}^{\dagger},\hat{a}]] = \mathbf{W}(\mathcal{O}(1,\hat{a}),\mathcal{H}[\hat{a}^{\dagger},\hat{a}]). \tag{B11}$$

The operator **W** denotes all possible Wick contractions between $\mathcal{O}(1,a)$ and $\mathcal{H}[\hat{a}^{\dagger},a]$. To proceed we suppose that the Hamiltonian has only a linear and quadratic term in \hat{a}^{\dagger} , i.e.,

$$\mathcal{H}[\hat{a}^{\dagger}, \hat{a}] = \int_{x} \hat{a}_{x}^{\dagger} \mathcal{L}_{x}[\hat{a}] + \frac{1}{2} \int_{x, y} \hat{a}_{x}^{\dagger} \hat{a}_{y}^{\dagger} \mathcal{B}_{x, y}[\hat{a}], \qquad (B12)$$

with $\mathcal{B}_{xy}[\hat{a}] = \mathcal{B}_{yx}[\hat{a}]$. Then the commutator will be

$$[\mathcal{O}(1,\hat{a}),\mathcal{H}[\hat{a}^{\dagger},\hat{a}]] = \int_{x} \frac{\delta}{\delta \hat{a}_{x}^{\dagger}} \mathcal{H}[\hat{a}^{\dagger},\hat{a}] \frac{\delta}{\delta \hat{a}_{x}} \mathcal{O}(1,\hat{a})$$

$$+ \frac{1}{2} \int_{x,y} \frac{\delta}{\delta \hat{a}_{x}^{\dagger}} \frac{\delta}{\delta \hat{a}_{y}^{\dagger}} \mathcal{H}[\hat{a}^{\dagger},\hat{a}]$$

$$\times \frac{\delta}{\delta \hat{a}_{x}} \frac{\delta}{\delta \hat{a}_{y}} \mathcal{O}(1,\hat{a}). \tag{B13}$$

As a consequence, the expectation (B10) evaluates to

$$\frac{d}{dt_{f}}\langle \mathcal{O}_{t_{f}} \rangle = \left\langle \int_{x} \left(\mathcal{L}_{x}[\hat{a}] + \int_{y} \mathcal{B}_{x,y}[\hat{a}] \right) \frac{\delta}{\delta \hat{a}_{x}} \mathcal{O}(1,\hat{a}) \right\rangle
+ \frac{1}{2} \left\langle \int_{x,y} \mathcal{B}_{x,y}[\hat{a}] \frac{\delta}{\delta \hat{a}_{x}} \frac{\delta}{\delta \hat{a}_{y}} \mathcal{O}(1,\hat{a}) \right\rangle.$$
(B14)

We can give an interpretation in terms of a stochastic process ϕ_t defined by

$$\partial_t \phi_{x,t} = \mathcal{L}_x[\phi] + \int_{\mathcal{V}} \mathcal{B}_{x,y}[\phi] + \xi_{x,t}, \qquad (B15)$$

$$\langle \xi_{x,t} \xi_{x',t'} \rangle = \delta(t - t') \mathcal{B}_{x,x'}[\phi], \tag{B16}$$

$$\mathcal{O}_{t_{\rm f}} = \mathcal{O}(1, \phi_{t_{\rm f}}). \tag{B17}$$

Indeed, applying the Itô formalism (see, e.g., [42]) to the expectation $\langle \mathcal{O}_{t_f} \rangle_{\xi}$ yields Eq. (B14). Comparing Eq. (B15) to Eq. (104), we note that the drift term also contains a term linear in \mathcal{B} . The reason is that to arrive at Eq. (104) the shifted Hamiltonian had been used. Indeed, this is accounted for by shifting in Eq. (B12) $\hat{a}^{\dagger} \rightarrow \hat{a}^{\dagger} + 1, \mathcal{L}_x \rightarrow \mathcal{L}_x + \int_y \mathcal{B}_{x,y}$, while $\mathcal{B}_{x,y}$ remains unchanged. Thus Eqs. (104) and (B15) are equivalent.

APPENDIX C: STOCHASTIC EQUATIONS OF MOTION WITH PURELY MULTIPLICATIVE NOISE

The terms in the action which cause us trouble are of the form

$$\delta S[a^*, a] = \frac{1}{2} \int_t (a_t^*)^2 a_t^2.$$
 (C1)

Decoupling with a noise ξ_t , with $\langle \xi_t \rangle = 0$, $\langle \xi_t \xi_t' \rangle = \delta(t - t')$, we have

$$e^{-\delta \mathcal{S}[a^*,a]} = \langle e^{i\int_t a_t^* a_t \xi_t} \rangle_{\mathcal{E}}.$$
 (C2)

Thus the equation of motion will have a term of the form

$$\partial_t a_t = i a_t \xi_t + \cdots . \tag{C3}$$

This is a multiplicative noise. In the following, we will solve this equation, dropping all further terms, both for a real noise, and for an imaginary noise. To reserve the notation ϕ_t for the phase, we noted this variable a_t ; it reminds the coherent-state variable \hat{a} .

1. Integrating a stochastic equation of motion with purely multiplicative noise: Real case

To start, consider the easier case of a real noise

$$\partial_t a_t = \xi_t a_t. \tag{C4}$$

We can make the ansatz

$$a_t = a_0 e^{\phi_t}. (C5)$$

This leads to

$$\partial_t \phi_t = \xi_t - \frac{1}{2}.\tag{C6}$$

Note the drift term; using Itô calculus with $dB_t = \xi_t dt$, the latter equation reads

$$d\phi_t = dB_t - \frac{1}{2}dt. \tag{C7}$$

We infer, as claimed,

$$da_{t} = a_{0}de^{\phi_{t}}$$

$$= a_{0}e^{\phi_{t}} \left[d\phi_{t} + \frac{1}{2}d\phi_{t}^{2} + \cdots \right]$$

$$= a_{0}e^{\phi_{t}} \left[dB_{t} - \frac{1}{2}dt + \frac{1}{2}dB_{t}^{2} + \cdots \right]$$

$$= a_{t}dB_{t}.$$
(C8)

Therefore, for arbitrary λ , the generating function is

$$\left\langle \left(\frac{a_t}{a_0}\right)^{\lambda}\right\rangle = \left\langle e^{\lambda\phi_t}\right\rangle = \left\langle e^{\lambda\int_0^t d\tau(\xi_{\tau} - \frac{1}{2})}\right\rangle = e^{\lambda(\lambda - 1)t/2}.$$
 (C9)

The probability is obtained by inverse-Laplace transforming,

$$P(\ln(a/a_0),t) = \int \frac{d\lambda}{2\pi} \langle e^{i\lambda[\ln(a/a_0) - \ln(a_t/a_0)]} \rangle$$

$$= \int \frac{d\lambda}{2\pi} e^{i\lambda \ln(a/a_0) - \lambda(\lambda + i)t/2}$$

$$= \frac{e^{-\frac{[\ln(a/a_0) + t/2]^2}{2t}}}{\sqrt{2\pi t}}.$$
 (C10)

This leads to the probability as a function of a,

$$P(a,t) = \frac{d \ln(a/a_0)}{da} P(\ln(a/a_0),t) = \frac{e^{-\frac{[\ln(a/a_0)+t/2]^2}{2t}}}{\sqrt{2\pi t}a}.$$
 (C11)

Integrating $\int_a a^n P(a,t)$ we find

$$\langle a_t^n \rangle = a_0^n e^{n(n-1)t/2}.$$
 (C12)

This yields, as it should,

$$\partial_t \langle a_t^n \rangle = \frac{n(n-1)}{2} \langle a_t^n \rangle.$$
 (C13)

2. Integrating a stochastic equation of motion with purely multiplicative noise: Imaginary case

For an imaginary noise as in Eq. (C3), we start from

$$\partial_t a_t = i\xi_t a_t. \tag{C14}$$

We make the ansatz

$$a_t = a_0 e^{i\phi_t + t/2}. (C15)$$

This leads to

$$\partial_t \phi_t = \xi_t. \tag{C16}$$

Note the drift term, analogous to the one in Eq. (C7), but with opposite sign. The probability to find ϕ is then given by

$$P(\phi,t) = \frac{e^{-\frac{\phi^2}{2t}}}{\sqrt{2\pi t}}.$$
 (C17)

The generating function for a_t , and its complex conjugate a_t^* , or more precisely for their logarithms, reads

$$Z(\lambda, \lambda^*) = \left\langle a_t^{\lambda^*} (a_t^*)^{\lambda} \right\rangle = \left(a_0 e^{t/2} \right)^{\lambda + \lambda^*} e^{-(\lambda + \lambda^*)^2 t/2}.$$
 (C18)

In Sec. IV G, we had seen that the probability for n-particle occupation is still given by the normalized coherent state. Even though here we are only dealing with a toy version of the real equation of motion, namely Eq. (C14), we can still study the same observable. This will shed light on the convergence issues noted in Sec. IV G.

$$p_n(t) = \left\langle e^{-a_t} \frac{a_t^n}{n!} \right\rangle_{a_t}$$
$$= \sum_{t=0}^{\infty} (-1)^k \left\langle \frac{a_t^{n+k}}{k! \, n!} \right\rangle_{a_t}$$

$$= \sum_{k=0}^{\infty} \frac{(-1)^k}{n!k!} (a_0 e^{t/2})^{n+k} e^{-(n+k)^2 t/2}.$$
 (C19)

For large times t, this converges towards $p_0(t) \to 1 - a_0$, $p_1(t) \to a_0$, $p_{n>2} \to 0$, which is no longer a probability for $a_0 > 1$. Another option is to write the above expectation as an integral over $P(\phi, t)$,

$$p_n(t) = \int_{-\infty}^{\infty} d\phi \, \frac{e^{-\frac{\phi^2}{2t}}}{\sqrt{2\pi t}} \, e^{-a_0 e^{i\phi + t/2}} \, \frac{(a_0 e^{i\phi + t/2})^n}{n!}. \tag{C20}$$

When the time t becomes large, the integral starts to oscillate. As an example, for $a_0 = 2$ it can no longer be done for $t \gtrsim 5$. Using (C17), we can also study the probability distribution for the real part

$$a_{\rm r} = \frac{a+a^*}{2}.\tag{C21}$$

Defining $b := e^{-t/2}a_r/a_0$, this is

$$P_{\rm r}^{b}(b,t) = \frac{1}{\sqrt{2\pi t (1 - b^{2})}}$$

$$\times \sum_{n = -\infty}^{\infty} \left[e^{-\frac{[a\cos(b) + 2\pi n]^{2}}{2t}} + e^{-\frac{[a\cos(b) - 2\pi n]^{2}}{2t}} \right]$$

$$= \frac{\vartheta_{3}\left(\frac{a\cos(b)}{2}, e^{-t/2}\right) + \vartheta_{3}\left(\frac{a\cos(b)}{2}, e^{-t/2}\right)}{\sqrt{2\pi t (1 - b^{2})}}. \quad (C22)$$

For small times, it is dominated by the term n = 0,

$$P_{\rm r}^b(b,t) \approx P_{\rm r}^{b,{\rm app}}(b,t) := \frac{\sqrt{2} e^{-\frac{{\rm acos}(b)^2}{2t}}}{\sqrt{\pi t (1-b^2)}}.$$
 (C23)

Note that

$$P_{\rm r}^a(a_{\rm r},t) = \frac{e^{-t/2}}{a_0} P_{\rm r}^b \left(\frac{a_{\rm r}e^{-t/2}}{a_0},t\right).$$
 (C24)

We checked this result numerically; see Fig. 10.

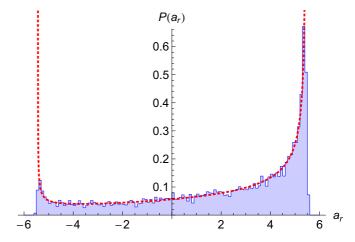


FIG. 10. The probability (C24) (red dashed), compared to a direct simulation of Eq. (C14) with 10^4 samples, $\delta t = 10^{-4}$, t = 2, $a_0 = 2$ (blue); the thin vertical red lines delimit the domain of the analytic solution. Some simulation points lie outside, due to a finite δt .

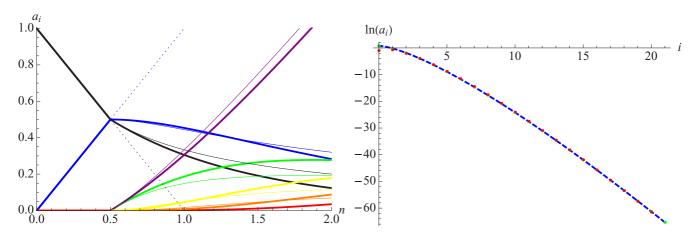


FIG. 11. Left: MF phase diagram of the weighted Manna model (wMM) (thick lines), compared to the MF phase diagram of the standard Manna model (thin lines). Right: The distribution of a_i for the wMM as a function of i (red points), for n = 0.8. The blue dashed line and green points are a Poisson distribution with $2.286e^{-0.989i}/i!$.

APPENDIX D: MF SOLUTIONS FOR THE WEIGHTED MANNA MODEL AND THE ABELIAN SANDPILE MODEL

1. MF solution for the weighted Manna model

The MF rate equations for the weighted Manna model are

$$\partial_t a_i = -(i-1)a_i \Theta(i \ge 2) + (i+1)a_{i+2}$$

$$+2 \left[\sum_{j \ge 2} a_j (j-1) \right] (a_{i-1} - a_i).$$
 (D1)

To give an analytic solution of this set of equations is left as a challenge to the reader. On the left of Fig. 11, we show the result of a numerical integration of the rate equations, starting from a Poisson distribution with on average n grains per site. As can be seen from the right of Fig. 11, the tail of the ensuing distribution remarkably no longer is an exponential, but close to a Poissonian. This can be seen from Eq. (D1): Setting

$$a_i \to \frac{a_2(2\rho)^{i-2}}{i!},\tag{D2}$$

the leading term in i is canceled. Indeed, the numerical solution confirms that this is approximately true; more precisely, $\frac{a_i 2\rho}{(i+1)a_{i+1}} \rightarrow 1$ for i large. (In practice, one can get the first $20 \dots 50$ coefficients, depending on n.)

2. MF solution for the 2D-ASM model

As a further excursion, consider the 2-dimensional Abelian sandpile model (ASM), which moves four grains to the four nearest neighbors, if the height at a given sites reaches or exceeds four. Its MF rate equations are

$$\partial_t a_i = -a_i \Theta(i \geqslant 4) + a_{i+4} + 4 \left[\sum_{j \geqslant 4} a_j \right] (a_{i-1} - a_i).$$
 (D3)

Using $\sum_i a_i = 1$, one can eliminate the infinite sum in the square bracket. Let us again suppose geometric progression, this time for i > 4,

$$a_i = a_4 \kappa^{i-4}, \quad i > 4. \tag{D4}$$

Solving as for the Manna model the 5 first equations, as well as the constraints (172) and (173), we obtain a nontrivial solution, with non-negative coefficients, for $n \ge \frac{3}{2}$,

$$a_0 = \frac{1}{4n - 2},\tag{D5}$$

$$a_1 = \frac{2(n-1)}{(1-2n)^2},$$
 (D6)

$$a_2 = \frac{12(n-2)n+13}{2(2n-1)^3},$$
 (D7)

$$a_i = \frac{4(n-1)[4(n-2)n+5]}{(1-2n)^4} \left[\frac{2n-3}{2n-1}\right]^{i-3}, \quad i \geqslant 3.$$
 (D8)

This is depicted in Fig. 12. For $n < \frac{3}{2}$, any set of positive a_i , such that $a_i = 0$ for i > 3, and which satisfy the constraints $\sum_i a_i = 1$ and $\sum_i i a_i = n$, is possible. Which solution is picked is given by the initial conditions. Integrating the rate equations (D3) starting from a Poissonian distribution with expectation n leads to the solid lines drawn in Fig. 12. This reproduces the solution (D5)–(D8) for $n > n_c \approx 1.542$. We have checked these analytical predictions by a direct numerical solution; see dots in Fig. 12.

Note that the same analysis can be done for an ASM on a lattice of coordination number n; the most interesting such case is n = 3, e.g., the 2-dimensional honeycomb lattice. The MF equations of motion then read

$$\partial_t a_i = -a_i \Theta(i \geqslant n) + a_{i+n} + n \left[\sum_{j \geqslant n} a_j \right] (a_{i-1} - a_i).$$
 (D9)

APPENDIX E: THE TOPPLE-AWAY MANNA MODEL, ITS CONVERGENCE TO MF, AND LIMIT OF LARGE DIMENSIONS

The question arises as to how one can go from the standard Manna model with nearest-neighbor topplings to its mean-field variant. There is indeed a series of models which does this interpolation. To this aim, define the following:

The topple-away Manna model of range r. If on a site (i, j) there are two grains, do twice: Randomly pick a site (i', j') with

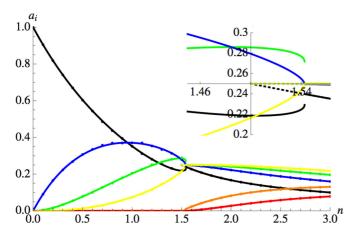


FIG. 12. Main plot: The phase diagram for the MF version of the 2d-ASM. The branch starting at n=1.5 and $a_{i,i\leqslant 3}=\frac{1}{4}$ is the MF solution, Eqs. (D5) and following. The branch starting at n=0 was obtained by a direct numerical integration of the rate equation (D3), setting to zero all $a_{i,i>100}$, and waiting until a steady state is reached. The dots are the result of a direct MC simulation of the MF-ASM model, starting from a Poisson distribution; we randomly distributed n times size 2 grains on a lattice of size 150. Note that Eq. (D3) allows us to predict the state reached by Monte Carlo. Also note that the branch with $a_{i,i\geqslant 4}=0$, which exists up to $n\leqslant 3$, remains attractive beyond $n=\frac{3}{2}$ (see inset, solid lines). The dashed lines are the solution (D5) and following.

 $|i-i'| \le r, |j-j'| \le r$, and move a grain from site (i,j) to site (i',j'). (We exclude the origin; two grains may end up on the same site.)

We can show via a numerical simulation that this variant converges to the MF Manna model for $r \to \infty$; in practice for r = 15, one is already very close. In Fig. 13 we show how n_c (below which there is no active phase) decreases with the range r; in fact, we find numerically that

$$n_c \approx \frac{1}{2} + \frac{0.1612}{r}.$$
 (E1)

In Fig. 14 (solid lines) we show how for r = 6 the phase diagram is already close to the MF model (thin lines). Intuitively, this is not surprising: When increasing the number of sites onto which an active site can depose its grains during a

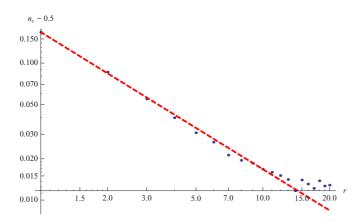


FIG. 13. A log-log plot of $n_c - 0.5$ as a function of the range r. The shown line is a power-law fit $n_c - 0.5 = 0.168069/r$.

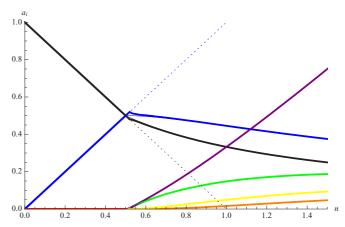


FIG. 14. The phase diagram for r = 6. The transition is at $n_c = 0.517$. Figure to be compared to Fig. 8. The thin underlying lines are the MF solution, the thin dotted lines the unstable solution.

toppling event, these "neighboring" sites are better and better described by a MF approach.

It is also intuitively clear that the same phenomenon takes place in high dimension d. Indeed, for $d = \infty$, each active grain has 2d neighbors, and supposing the state can locally be described by the number of grains and the activity, the assumptions in the derivation of MF theory become valid in the large-d limit.

APPENDIX F: CHANGING VARIABLES IN THE CSPI

In Ref. [17] the authors propose an alternative exact path-integral representation for stochastic systems, and show its equivalence to the CSPI formalism. Their idea is to start from a rate equation for the particle number ρ on a site. Suppose in a time step δt the particle number changes with probability $p(\rho)\delta t$ by a fixed amount $\delta \rho$. To write down the path integral for a time slice δt , the change $\rho(t+\delta t)=\rho(t)+\delta\rho$ is enforced by an integral over the auxiliary field ρ^* , $\int_{\rho^*}e^{-\rho^*[\rho(t+\delta t)-\rho(t)-\delta\rho]}$. Averaging this quantity over the random process results in the action for a time slice,

$$\int_{\rho^*} e^{-\delta \mathcal{S}[\rho^*, \rho]} = \int_{\rho^*} e^{-\rho^* [\rho(t+\delta t) - \rho(t)]} \langle e^{\rho^* \delta \rho} \rangle$$

$$= \int_{\rho^*} e^{-\rho^* [\rho(t+\delta t) - \rho(t)]}$$

$$\times [1 + p(\rho)\delta t (e^{\rho^* \delta \rho} - 1) + \cdots]. \tag{F1}$$

Taking δt small, and concatenating many such slices, yields the path integral $\int \mathcal{D}[\rho]\mathcal{D}[\rho^*]e^{-S[\rho^*,\rho]}$, with action

$$S[\rho^*, \rho] = \int_t \rho^*(t) \partial_t \rho(t) - \mathcal{H}_{\rho}[\rho^*(t), \rho(t)], \quad (F2)$$

$$\mathcal{H}_{\rho}[\rho^*, \rho] = p(\rho)(e^{\rho^*\delta\rho} - 1). \tag{F3}$$

⁵The reader of Ref. [17] may benefit from the following remarks: (i) The sum in Eq. (2) contains both the pairs (i, j) and (j, i). (ii) The unit vector \mathbf{e}_{ij} points from j to i. (iii) There is no factor of 2 in the definition of γ after Eq. (3). (iv) In Eq. (6) μ and ν are exchanged.

Consider now the examples discussed in Sec. II B. There was pair annihilation with probability $p(\rho) = \frac{\nu}{2}\rho(\rho-1)$ and $\delta\rho = -2$, decay with rate $\mu\rho$ and $\delta\rho = -1$, and finally particle creation (branching) with rate $\kappa\rho$ and $\delta\rho = 1$. Promoting the fields to operators $\rho \to \hat{\rho}$ and $\rho^* \to \hat{\rho}^{\dagger}$, the corresponding Hamiltonian reads

$$\mathcal{H}_{\rho} = \frac{\nu}{2} [e^{-2\hat{\rho}^{\dagger}} - 1] \hat{\rho} (\hat{\rho} - 1) + \mu [e^{-\hat{\rho}^{\dagger}} - 1] \hat{\rho} + \kappa [e^{\hat{\rho}^{\dagger}} - 1] \hat{\rho}.$$
(F4)

We recall that as in Sec. IIJ the Hamiltonian is written in normal-ordered form, which is required in order to pass from the path integral to the operator formalism, and back. By construction, the commutation relations are canonical,

$$[\hat{\rho}, \hat{\rho}^{\dagger}] = 1. \tag{F5}$$

Let us compare (F4) to our Hamiltonian (15) for the same processes, reproduced here for convenience,

$$\mathcal{H} = \frac{\nu}{2} [\hat{a}^{\dagger} \hat{a}^2 - (\hat{a}^{\dagger})^2 \hat{a}^2] + \mu [\hat{a} - \hat{a}^{\dagger} \hat{a}] + \kappa [(\hat{a}^{\dagger})^2 \hat{a} - \hat{a}^{\dagger} \hat{a}].$$
(F6)

Demanding that $\mathcal{H} \stackrel{!}{=} \mathcal{H}_{\rho}$ uniquely identifies, already for $\nu = 0$, the operator identities

$$\hat{a}^{\dagger} = e^{\hat{\rho}^{\dagger}}, \quad \hat{a} = e^{-\hat{\rho}^{\dagger}}\hat{\rho}.$$
 (F7)

A few checks are in order: First, the commutation relations $[\hat{a}, \hat{a}^{\dagger}] = 1$ and $[\hat{\rho}, \hat{\rho}^{\dagger}] = 1$ are compatible. Since the transformation (F7) is canonical, i.e., has Jacobian one, the path integrals also map onto each other.

Finally, the term proportional to ν is correctly reproduced; it is the only term which requires the commutation relations, and for which normal ordering is important.

We finally note that the Hamiltonian (46) for hopping, in the continuous limit, $\mathcal{H} = -D\nabla \hat{a}^{\dagger} \nabla \hat{a}$, maps onto

$$\mathcal{H}_{\rho} = \int -D\nabla \hat{\rho}^{\dagger} \nabla \hat{\rho} + D(\nabla \hat{\rho}^{\dagger})^{2} \hat{\rho}.$$
 (F8)

This expression can directly be derived from the equivalent of Eq. (F1) for diffusion [17]. Decoupling the term quadratic in $\hat{\rho}^{\dagger}$ with a noise yields Eq. (157). That these two equations are equivalent is nontrivial: the formalism used to derive Eq. (157) works with coarse-grained fields, whereas $\hat{\rho}$ above is a microscopic field. The reason for this equivalence is that when coarse-graining Eq. (157), its linear structure allows passing coarse graining through to the fields.

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