

Quantum walks and exact RG in de Sitter space

Daniel Green and Kshitij Gupta

Department of Physics, University of California at San Diego, La Jolla, California 92093, USA



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The local physics of light scalar fields in de Sitter space is well described by classical random walks, as expressed through the framework of stochastic inflation. Recent work has clarified how this formalism arises from quantum field theory and the renormalization group (RG), allowing for corrections to this formalism to be determined order by order. Yet, this description is incomplete. For example, the quantum dynamics of these fields are expected to become important when determining the tail of the probability distribution for the fluctuations. In this paper, we develop the understanding of fields in de Sitter as a quantum walk in order to bridge the gap between the classical and quantum description. We use the framework of exact RG to calculate the evolution equation for the reduced density matrix of the long wavelength fields. This master equation provides the direct map from light fields to models of quantum walks. We show how to reduce the master equation to stochastic inflation, and provide a new understanding of how the higher-order corrections arise. In the process, we demonstrate that divergences and secular growth in de Sitter, for both light and heavy fields, can be absorbed by (dynamical) renormalization.

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I. INTRODUCTION

A complete quantum description of our Universe remains one of the key ambitions of theoretical physics [1]. It requires reckoning with some of the most basic conceptual questions about the origin of the Universe, the predictions of quantum gravity, and the interpretation of quantum mechanics in closed systems.

The dynamics of light scalar fields present a controlled opportunity to address some of these essential questions. This first challenge is that massless scalar fields do not have a de Sitter (dS) invariant vacuum state [2]. In addition, perturbative corrections exhibit secular growth that causes the perturbative expansion to break down [3,4]. It was suggested long ago [5,6] that these features of fields in de Sitter should be interpreted as consequences of local field values undergoing a classical random walk. In this language, known as stochastic inflation, the absence of a de Sitter invariant state and the secular growth become familiar aspects of random walks without a potential [7].

Recently, it has been shown that the random walk description of light scalars can be recovered in quantum field theory (QFT) on a fixed dS background using renormalization group (RG) techniques [8–18]. In the process, many of these longstanding issues are resolved at both a technical

and conceptual level. From the RG point of view, the theory evolves away from the trivial fixed point to a nontrivial self-similar description that is equivalent to a random walk. At the same time, this description must have a limited regime of validity; a classical random walk generates entropy which cannot be the case in a closed quantum system. The stochastic formalism is therefore limited to subsystems within de Sitter space [19–21], although that aspect is not always transparent in the derivation. A complete understanding of light fields in de Sitter should furnish an explanation for the origin and breakdown of this description.

A more complete quantum stochastic description of the long modes in de Sitter is nearly as old [22–26] as the original, classical description. Rather than calculating a probability distribution in the field basis, one calculates the evolution for the reduced density matrix of the long wavelength fields. Tracing over the short wavelength modes gives rise to an evolution equation. When expressed in terms of the Wigner quasiprobability distribution, this evolution remains in a random walk form in “phase space.” This approach has been revived more recently in [27–31], motivated in part by recent progress understanding the quantum nature of de Sitter space.

Despite this long history, the quantum description of light fields remains underdeveloped. For one, the stochastic evolution equation appears even in trivial free field theories, where the short and long wavelength models are uncorrelated. In this regard, it cannot be the case that the random walk arises only from tracing over short wavelength modes. Instead, one is performing an additional trace, at least

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implicitly, to arrive at a Fokker-Planck equation. Although one recovers the expected evolution to leading order, the quantum mechanical nature of this evolution equation depends on this procedure. Many of these subtleties arise even in flat space, and are a consequence of working with EFT observables within the in-in formalism [32–35]. These challenges make it impossible for current techniques to determine the tail of the local one-point distribution [36,37], or whether the universe is eternally inflating [38]. These types of questions require a nonperturbative understanding of the quantum evolution of the density matrix.

In this paper, we will use the framework of quantum walks [39] to better understand the quantum nature of the long wavelength fluctuations in de Sitter. A walker can be coupled to N -quantum coins (or N -harmonic oscillators) so that the location of the walker is entangled with the space of the coins. Tracing over the coins gives rise to different types of stochastic behavior, including both purely classical and uniquely quantum walks, depending on how the systems are coupled. The coins alone may be in a product state so that no entanglement entropy is generated until they are coupled to the walker. In fact, we will show that the origin of stochastic inflation is most easily derived by introducing a local physical measurement similar to the walker.

We will then use the framework of exact RG [40] to show that light fields in de Sitter space obey the equations of a general quantum walk. We will show how integrating out a shell of momentum gives rise to an evolution equation for the reduced density matrix, generalizing previous results from flat space [33,41] (due in part to momentum space entanglement [42–46]). This master equation for the evolution of the density matrix only gives rise to the physics of stochastic inflation after we sum over Fourier modes (tracing over information at separated points). After deriving an exact equation for the evolution, we will show that this equation can be approximated by the original equations of stochastic inflation and its generalization to the density matrix and Wigner distribution. The resulting evolution equations are most intuitive in the Wigner representation, where we find an evolution equation qualitatively similar to the Caldeira-Leggett model [47] and quantum Brownian motion [48,49], but that differs in several key respects.

As a byproduct of our exact RG analysis, we demonstrate that the divergences in cosmology can be absorbed into the definition of the action and operators in precise analogy with RG in flat space. Unlike flat space, the RG transformation involves time evolution giving rise to dynamical RG flow. This confirms the intuition that large logs found in perturbation theory can, and should be, resummed using RG techniques [50,51].

A. Summary of key results

The main result of this paper is the derivation of the exact evolution of the reduced density matrix of a light scalar

field on a dS background, Eq. (3.45). We further trace over the Fourier modes of the field to arrive at a master equation for the evolution of the density matrix for the measurement of the field at a single point, Eq. (4.14). When expressed in terms of the Wigner distribution, this takes the form of a random walk in phase space, Eq. (4.16).

When our results are reduced to the diagonal elements, we reproduce the results of stochastic inflation at leading order [5–7]. Moreover, corrections to the Fokker-Planck equation arise from the self-consistent elimination of the off-diagonal terms in the Wigner distribution and match the known corrections at this order [10,14,52]. In this sense, our results are consistent with known results from stochastic inflation and QFT on a fixed dS background. However, the formalism is nonperturbative and applies to the entire density matrix, opening the possibility of calculating statistical quantities that are not accessible using previous techniques.

The organization of this paper is as follows: In Sec. II, we discuss simple models of quantum walks that capture both conventional classical random walks and uniquely quantum contributions. In Sec. III, we derive the exact RG equations for light fields in dS. We then show how they combine with time evolution to produce Lindbladian evolution of the density matrix. In Sec. IV, we show how our exact RG reproduces the equations for quantum walks. Then, we derive stochastic inflation and its next-to-leading order (NLO) corrections from our master equation, and discuss how these techniques can be applied at next-to-next-to-leading order (NNLO) and beyond. In Sec. V, we demonstrate that dynamical RG in dS resums all UV logs. We conclude in Sec. VI.

II. QUANTUM WALKS FOR COSMOLOGY

The evolution of a scalar field in de Sitter has long been argued to (locally) resemble a random walk [5,6]. The resulting stochastic approach is classical in nature, leaving quantum mechanics responsible only for the microscopic origin of the noise. The universe, of course, remains quantum mechanical and ultimately the classical random walk must be replaced with a quantum generalization. In this section, we will present some simple quantum walks that will serve as useful intuition for de Sitter space. The models in this section are closely related to the literature on quantum Brownian motion [47–49]; however, it is important that the quantum walks relevant to cosmology are not equivalent to these models and we will emphasize the points of departure.

A. Density matrix of a Gaussian walk

We will imagine our system is split into a massive field whose wave function in the position basis $\Psi(X)$ encodes the random fluctuations of a N harmonic oscillators in their ground states,

$$\psi(\vec{x}_i) = \frac{1}{(2\pi\sigma^2)^{1/4}} \exp(-x_i^2/(4\sigma^2)). \quad (2.1)$$

In this regard, the location of heavy field X defines a pointer basis [53] for the measurement of the oscillators. We can implement a classical Gaussian random walk using this quantum system as follows: for $N = 0$, the initial state of the walker is definitely at the origin, $\Psi(X) = \delta(X)$. We then act on $|\Psi\rangle$ by a shift proportional to x_i for each $i = 1, \dots, N$:

$$|\Psi, N\rangle = e^{i\hat{P}\sum_i \hat{x}_i} |X = 0\rangle \otimes \prod_i |\psi_i\rangle. \quad (2.2)$$

The reduced density matrix for the walker after coupling to N oscillators is given by

$$\begin{aligned} \text{Tr}_{\psi_i} \rho &= \rho_N(X, X') \\ &= \left(\prod_i \int dx_i |\psi(x_i)|^2 \right) \left| X = \sum_i x_i \right\rangle \left\langle X' = \sum_i x_i \right| \end{aligned} \quad (2.3)$$

$$= \delta(X - X') \frac{1}{\sqrt{2\pi\sigma^2 N}} \exp(-X^2/(2\sigma^2 N)). \quad (2.4)$$

The second line can be determined by solving Eq. (2.2) in the momentum basis of the walker. This is mathematically equivalent to a common derivation of probability distribution for N steps of a classical random walk using the Fourier transform. Indeed, this is a classical random walk for all practical purposes.

To gain intuition for our evolution equations, it is useful to derive the master equation for the evolution of the density matrix when increasing the number of harmonic oscillators from N to $N + 1$,

$$\rho_{N+1}(X, X') = \text{Tr}_{x_{N+1}} [e^{i\hat{P}\hat{x}_{N+1}} \rho_N \otimes |\psi\rangle\langle\psi| e^{-i\hat{P}\hat{x}_{N+1}}]. \quad (2.5)$$

We can use this expression to derive an equation for the time evolution in terms of $T = Ndt$ in the continuum limit $N \rightarrow \infty$ and $dt \rightarrow 0$ while holding T and $\sigma^2 N = \alpha T$ finite. Taking the trace in the x_{N+1} -position basis and treating $x_{N+1} = \mathcal{O}(dt^{1/2})$ we can expand both sides to $\mathcal{O}(dt)$ to find

$$\frac{\partial}{\partial T} \rho = -\frac{\alpha}{2} (\hat{P}^2 \rho + \rho \hat{P}^2 - 2\hat{P} \rho \hat{P}). \quad (2.6)$$

In the X -position basis $\rho(X, X', T)$ this becomes the differential equation

$$\frac{\partial}{\partial T} \rho(X, X', T) = \frac{\alpha}{2} \left(\frac{\partial^2}{\partial X^2} + \frac{\partial^2}{\partial X'^2} + 2 \frac{\partial^2}{\partial X \partial X'} \right) \rho(X, X', T). \quad (2.7)$$

Naively, one might expect the diagonal elements $X = X'$ to obey the classical heat equation, but this is not the case.

In fact, the density matrix is actually singular if we take $X = X'$, which suggests that we need to be more careful solving for the density matrix, even when we expect it to be diagonal.

The master equation for our walk is easier to solve in the momentum representation of the density matrix, $X \rightarrow P$ and $X' \rightarrow P'$, so that our evolution is given by

$$\frac{\partial}{\partial T} \rho(P, P', T) = -\frac{\alpha}{2} (P - P')^2 \rho(P, P', T). \quad (2.8)$$

We can directly solve this first-order ordinary differential equation (ODE) to find

$$\rho(P, P', T) = \rho(P, P', 0) \exp\left(-\frac{\alpha}{2} (P - P')^2 T\right). \quad (2.9)$$

We can find the solution in the position basis by Fourier transform,

$$\begin{aligned} \rho(X, X', T) &= \int \frac{dP dP'}{(2\pi)^2} e^{iPX} e^{-iP'X'} \rho(P, P', 0) \\ &\quad \times \exp\left(-\frac{\alpha}{2} (P - P')^2 T\right). \end{aligned} \quad (2.10)$$

If we impose that $\rho(P, P', 0) = 1$, directly integrating reproduces Eq. (2.4).

We can gain a lot of useful intuition for these walks rewriting our results in terms of the Wigner distribution function,¹

$$W(x, p) = \int d\Delta x e^{-ip\Delta x} \rho\left(X = x + \frac{\Delta x}{2}, X' = x - \frac{\Delta x}{2}\right). \quad (2.11)$$

By direct calculation, our solution in Eq. (2.4) takes a familiar form,

$$W(x, p, T) = \exp(-x^2/(2\alpha T)). \quad (2.12)$$

We can bring this even closer to the classical description by transforming the master equation for the density matrix, Eq. (2.7), to an evolution equation for Wigner distribution. Specifically, if we change variables to $X, X' = x \pm \Delta x$ and using the chain rule

$$\frac{\partial}{\partial x} = \frac{\partial}{\partial X} + \frac{\partial}{\partial X'}. \quad (2.13)$$

¹We can also define the Wigner distribution starting from momentum representation of the density matrix, $\rho(P, P', T)$, by taking the inverse Fourier transform of $P - P' \rightarrow x$, while the “momentum” of the Wigner representation becomes the other linear combination $P + P' \rightarrow 2p$.

This is the linear combination of derivatives in the evolution of the density matrix, so that

$$\frac{\partial}{\partial T} W(x, p, T) = \frac{\alpha}{2} \frac{\partial^2}{\partial x^2} W(x, p, T). \quad (2.14)$$

In this precise sense, we see that the Wigner distribution obeys the heat equation for a Gaussian quantum walk in the position basis. This is a strong hint that the Wigner distribution provides a more natural starting point for discussing the evolution of these kinds of quantum walks.

None of the conclusions in this section should be surprising, as we have simply translated a purely classical random walk into quantum language. However, in anticipating the possibility of a more fundamentally quantum system, it is useful to look at mathematically equivalent (and still effectively classical walks) in other representations. For example, suppose our walk was defined in momentum space so that one step of the walk shifts,

$$|\Psi, N\rangle = e^{i\hat{X} \sum_i \hat{x}_i} |P=0\rangle \times \prod_i |\psi_i\rangle. \quad (2.15)$$

Following the above logic with $\hat{P} \rightarrow \hat{X}$, the density matrix obeys the master equation in the position basis

$$\frac{\partial}{\partial T} \rho(X, X', T) = -\frac{\alpha}{2} (X - X')^2 \rho(P, P', T). \quad (2.16)$$

Of course, in terms of the Wigner distribution, using $X - X' = 2\Delta x$, this is just the heat equation,

$$\frac{\partial}{\partial T} W(x, p, T) = \frac{\alpha}{2} \frac{\partial^2}{\partial p^2} W(x, p, T). \quad (2.17)$$

In terms of Wigner distribution, both types of walks look purely classical. In contrast, recognizing a classical walk in terms of the density matrix directly requires picking the right variables from the outset.

B. Quantum Brownian motion with a potential

Now we would like to introduce additional forces on our quantum system. We start from the same Gaussian walk for a quantum particle via N harmonic oscillators in a bath, but now we will additionally immerse the walker, X , in a deterministic potential $V(X)$. As a result, we have the combined Hamiltonian of the walker and the harmonic oscillators as

$$H = \frac{P^2}{2M} + V(X) + P \sum_i x_i. \quad (2.18)$$

We can understand this Hamiltonian using the operator evolution equations, or the classical equations of motion,

$$\dot{X} = \frac{P}{M} + x_i \quad \dot{P} = -V'(X). \quad (2.19)$$

This model is very similar to the Caldeira-Leggett model [47] but we notice the noise directly in the position variable, rather than a random force that appears in the $\dot{P}$ equation of motion. If $X \rightarrow x$ and $P \rightarrow p$ were purely classical statistical variables, we expect the classical phase space probability distribution $W_c(x, p, T)$ to follow the Fokker-Planck equation

$$\begin{aligned} \frac{d}{dT} W_c(x, p, T) = & \alpha^2 \frac{\partial^2}{\partial x^2} W_c(x, p, T) - \frac{p}{M} \frac{\partial}{\partial x} W_c(x, p, T) \\ & + V'(x) \frac{\partial}{\partial p} W_c(x, p, T). \end{aligned} \quad (2.20)$$

Of course, X and P are operators and require a quantum mechanical equivalent of this evolution equation. After tracing over the harmonic oscillators, x_i , the evolution of the walker will be described by the evolution of the reduced density matrix, ρ_R . We can write this evolution using the Lindbladian formalism, so that

$$\dot{\rho}_R = i[H, \rho] + \sum_{\mu} \left(L_{\mu} \rho L_{\mu}^{\dagger} - \frac{1}{2} \{L_{\mu}^{\dagger} L_{\mu}, \rho_R\} \right), \quad (2.21)$$

where L_{μ} are the jump operators given by

$$L_{\mu} = iP_i \langle \mu | \hat{x} | 0 \rangle_i. \quad (2.22)$$

The states $|\mu\rangle_i$ and $|0\rangle_i$ are states of the i th harmonic oscillator that we are tracing over at time t . The jump operator is nonzero only for the first excited state,

$$L_{\mu=1} = iP\alpha \quad \alpha = \langle 1 | \hat{x} | 0 \rangle = \sqrt{\frac{1}{2m\omega}} = \sigma, \quad (2.23)$$

and the Lindbladian evolution is therefore

$$\begin{aligned} \dot{\rho}_R = & -i[H, \rho] - \alpha^2 (p - p')^2 \rho \\ = & -i \left(\frac{p^2}{2M} - \frac{p'^2}{2M} \right) \rho - i(V(x) - V(x')) \rho_R \\ & - \alpha^2 (p - p')^2 \rho_R. \end{aligned} \quad (2.24)$$

To compare to the classical probability distribution, we transform to the Wigner representation

$$\begin{aligned} \frac{\partial W}{\partial t} = & \alpha^2 \frac{\partial^2 W}{\partial x^2} - i \left(V \left(x + \frac{i}{2} \partial_p \right) - V \left(x - \frac{i}{2} \partial_p \right) \right) W \\ & - \frac{p}{M} \frac{\partial W}{\partial x}. \end{aligned} \quad (2.25)$$

Expanding around the extremum where ∂_p is small, we get that

$$\begin{aligned} & V\left(x + \frac{i}{2}\partial_p\right) - V\left(x - \frac{i}{2}\partial_p\right) \\ &= i\frac{\partial}{\partial p}V'(x) - \frac{i}{24}\frac{\partial^3}{\partial p^3}V'''(x) + \dots \end{aligned} \quad (2.26)$$

and thus we have

$$\frac{\partial W}{\partial t} = \alpha^2 \frac{\partial^2 W}{\partial x^2} - \frac{p}{M} \frac{\partial W}{\partial x} + V'(x) \frac{\partial W}{\partial p} - \frac{V'''(x)}{24} \frac{\partial^3 W}{\partial p^3} + \dots \quad (2.27)$$

Comparing this to the classical phase space distribution (2.20), we see that we get higher order derivatives in p purely due to a quantum effect. These terms are suppressed by powers of $\hbar$.

C. A walk for stochastic inflation

Having introduced quantum walks with a potential, the final ingredient required for building intuition for cosmology is introduction of friction for the walker, X . This is straightforward in the Langevin description of a classical walk but is much more involved for a quantum walk. However, we have seen that quantum walks end up looking like classical walks in phase space when using the Wigner distribution. We will therefore consider a classical walk with friction that will reproduce key features of our quantum walk for the dS density matrix.

In the presence of friction, the general Langevin equation for a classical walker takes the form

$$\dot{x} = \frac{p}{m} + \xi_1 \quad \dot{p} = -\gamma p - V'(x) + \xi_2, \quad (2.28)$$

where $\gamma > 0$ controls the friction. For most of Newtonian physics applications, one considers Langevin equations with $\xi_1 = 0$. Such a system of equations describes random forces due to thermal baths, and the equilibrium probability distribution is the Maxwell-Boltzmann distribution

$$W_{\text{eq}}(x, p) \propto \exp\left(-\beta\left(\frac{p^2}{2m} + V(x)\right)\right). \quad (2.29)$$

This is not the kind of walk that describes fields in de Sitter. Instead, we should consider a Langevin system where the noise is only in x rather than in p , $\xi_1 \neq 0$ and $\xi_2 = 0$. In other words, we want to study the following walk:

$$\dot{x} = \frac{p}{m} + \sqrt{2D}\xi \quad \dot{p} = -\gamma p - V'(x), \quad (2.30)$$

which represents a particle in an external potential, with friction in p characterized by γ and kicks in x characterized by D . As we will see, even at $\hbar \rightarrow 0$, this equation reproduces key features that we will encounter in stochastic inflation. The phase space probability distribution, W_c , at

leading order, evolves as

$$\frac{dW_c}{dt} = D \frac{d^2 W_c}{dx^2} + \frac{d}{dp}(\gamma p W_c + V'(x) W_c) - \frac{p}{m} \frac{dW_c}{dx}. \quad (2.31)$$

In order to solve this system, we will consider the overdamped limit $\gamma \gg 1/t$, and will derive the solutions as a series in this limit. We assume that at leading order, the background value x is fixed while p fluctuates to come to equilibrium around a given value of x . The expectation value of p is described by

$$\frac{d}{dt}\langle p \rangle_x = -\gamma \langle p \rangle_x - V'(x). \quad (2.32)$$

The solution is

$$\begin{aligned} \langle p \rangle_x &= e^{-\gamma t} \langle p(0) \rangle_x - \int_0^t ds e^{\gamma(s-t)} V'(x) \\ &= e^{-\gamma t} \langle p(0) \rangle_x - \frac{V'(x)}{\gamma} (1 - e^{-\gamma t}), \end{aligned} \quad (2.33)$$

where we are treating $V(x)$ as time independent at leading order. We see that in the overdamped limit, $\langle p \rangle_x \rightarrow -V'(x)/\gamma$. Now, we can solve for the probability distribution of x . Integrating over p and substituting $\langle p \rangle_x$, we get the Fokker-Planck equation for x ,

$$\frac{dP(x)}{dt} = D \frac{d^2 P(x)}{dx^2} + \frac{d}{dx} \left(\frac{V'(x)}{m\gamma} P(x) \right), \quad (2.34)$$

and hence the equilibrium distribution for x given by

$$P_{\text{eq}}(x) = A e^{-V(x)/m\gamma D}. \quad (2.35)$$

We would also like to construct the equilibrium solution for the full phase space distribution, i.e., the equilibrium solution for (2.31). However, unlike the case for the Gibbs distribution, we do not have an analytic solution for generic $V(x)$. However, a closed form solution does exist for quadratic $V(x)$. Writing the potential as $V(x) = \frac{1}{2} k x^2$, we see that the equilibrium probability distribution is given by

$$W_{\text{eq}} = \alpha \exp(Ax^2 + Bxp + Cp^2), \quad (2.36)$$

where

$$A = -\frac{\gamma}{2D}, \quad B = -\frac{\gamma^2}{kD}, \quad C = -\frac{\gamma}{2mkD} - \frac{\gamma^3}{2k^2D}. \quad (2.37)$$

Explicitly, we have

$$W_{\text{eq}} \propto \exp \left(-\frac{\gamma}{2D} x^2 - \frac{\gamma^2}{Dk} xp - \frac{\gamma}{2kD} \left(\frac{1}{m} + \frac{\gamma^2}{k} \right) p^2 \right). \quad (2.38)$$

On integrating out p , we get

$$P(x) = A \exp \left(-\frac{\frac{1}{2}k\gamma}{kD + D\gamma^2 m} x^2 \right). \quad (2.39)$$

Writing this in the form $P \propto \exp(-V_{\text{eff}}(x)/m\gamma D)$, we see that the effective potential is given by

$$V_{\text{eff}} = \frac{1}{2} kx^2 - \frac{1}{2} \frac{k^2 x^2}{m\gamma^2} + \dots \quad (2.40)$$

where the expansion is in $k/m\gamma^2 \ll 1$. We see that even in the case of a Gaussian theory, the momentum will induce a nontrivial effect on the nonperturbative probability distribution of the position, and in particular on the tails. Finally, we note that our equilibrium solution is quite different from the thermal Gibbs distribution, which cleanly factorizes out as $\exp(-\beta(p^2/2m + V(x)))$. We will comment more on this in the next sections.

1. Corrections to the effective potential

Having provided the leading-order solution for the evolution equation for $P(x)$, we want to understand the next-order correction. We can do this in the Langevin picture,

$$\dot{p} = -\gamma p - V'(x), \quad \dot{x} = \frac{p}{m} + \sqrt{2D}\xi(t). \quad (2.41)$$

Formally, we can solve this via

$$\begin{aligned} p(t) &= - \int ds e^{-\gamma(t-s)} V'(x(s)) \\ &= - \int ds e^{-\gamma(t-s)} V'(x(t)) \\ &\quad - \int ds e^{-\gamma(t-s)} (s-t) V''(x(t)) \dot{x}(t) + \dots \end{aligned} \quad (2.42)$$

Using the Langevin equations to substitute $\dot{x}$ in terms of p , we get

$$p = -\frac{V'(x)}{\gamma} - \frac{V'(x)V''(x)}{m\gamma^3} + \frac{V''(x)}{\gamma^2} \sqrt{2D}\xi(t) + \dots \quad (2.43)$$

So, substituting this back into the equations for x , we get

$$\dot{x} = \sqrt{2D}\xi(t) \left(1 + \frac{V''(x)}{m\gamma^2} \right) - \left(\frac{V'(x)}{m\gamma} + \frac{V'(x)V''(x)}{m^2\gamma^3} \right). \quad (2.44)$$

Using the Stratonovich convention, we see that this translates into the Fokker-Planck equation,

$$\begin{aligned} \frac{\partial P}{\partial t} &= D \frac{\partial}{\partial x} \left(\left(1 + \frac{V''}{m\gamma^2} \right) \frac{\partial}{\partial x} \left(1 + \frac{V''}{m\gamma^2} \right) P(x) \right) \\ &\quad + \frac{\partial}{\partial x} \left(\frac{V'(x)}{\gamma} + \frac{V'(x)V''(x)}{m\gamma^3} \right) P. \end{aligned} \quad (2.45)$$

For the harmonic case, we get that

$$\frac{\partial P}{\partial t} = D \left(1 + \frac{k}{m\gamma^2} \right)^2 \frac{\partial^2 P}{\partial x^2} + \frac{\partial}{\partial x} \left[\left(\frac{kx}{m\gamma} + \frac{k^2 x}{m^2 \gamma^3} \right) P \right]. \quad (2.46)$$

For a Fokker-Planck equation of the form (2.34), the equilibrium solution is given by

$$P_{\text{eq}} \propto \exp \left(-\frac{V_{\text{eff}}}{m\gamma D} \right). \quad (2.47)$$

Thus, for the harmonic case, since we have that our effective diffusion in the Fokker-Planck equation is $D' = D(1 + \frac{k}{m\gamma^2})^2$, we get that V_{eff} is given by

$$V_{\text{eff}} = \frac{1}{2} kx^2 \left(1 + \frac{k}{m\gamma^2} \right) \left(1 - 2 \frac{k}{m\gamma^2} \right) = \frac{1}{2} kx^2 \left(1 - \frac{k}{m\gamma^2} \right), \quad (2.48)$$

which matches the harmonic effective potential in (2.40). In Sec. IV, we will use a morally equivalent procedure to derive similar corrections to stochastic inflation.

2. Comments on the equilibrium distribution

While we have a stationary (time-independent) solution, it is crucially different from the Gibbs distribution, which exhibits detailed balance. Detailed balance tells us that there are no probability currents, and that forward and backward transition rates are equally likely

$$\begin{aligned} \frac{dP_i}{dt} &= \sum W(j \rightarrow i) P_j - \sum W(i \rightarrow j) P_i = 0 \\ W(j \rightarrow i) P_j &= W(i \rightarrow j) P_i \quad \forall i, j. \end{aligned} \quad (2.49)$$

This is the statement of microscopic time reversibility. However, for our stationary solution, time reversibility and hence detailed balance is manifestly broken, and we have nonzero currents which tell us the preferential probability flow. Writing

$$\frac{dP}{dt} = -\nabla \cdot J = -(\partial_x J_x + \partial_p J_p), \quad (2.50)$$

we get

$$J_x = \frac{p}{m} W - D \frac{\partial W}{\partial x}, \quad J_p = -\gamma p W - V'(x) W. \quad (2.51)$$

Evaluating the currents for our equilibrium solution in Eq. (2.38), we get

$$J_x = \left(\gamma x + \frac{\gamma^2 p}{k} + \frac{p}{m} \right) W_{\text{eq}}, \quad J_p = -\gamma \left(p + \frac{V'(x)}{\gamma} \right) W_{\text{eq}}. \quad (2.52)$$

We can rewrite this, in our small parameter limit $k \ll m\gamma^2$, as

$$J_x = \frac{\gamma^2}{k} \left(p + \frac{V'(x)}{\gamma} \right) W_{\text{eq}}, \quad J_p = -\gamma \left(p + \frac{V'(x)}{\gamma} \right) W_{\text{eq}}. \quad (2.53)$$

We thus see that whenever we are off the momentum saddle point axis $p = -V'(x)/\gamma$, we get a nonzero current. However, on integrating out p , we constrain ourselves to flow along this axis. Thus, the reduced one-dimensional Fokker-Planck equation

$$\frac{dP}{dt} = D \frac{d^2 P(x)}{dx^2} + \frac{d}{dx} \left(\frac{V'(x)}{m\gamma} P(x) \right) \quad (2.54)$$

does satisfy detailed balance at equilibrium, as expected. Thus, stochastic inflation represented in the flat Friedmann-Robertson-Walker (FRW) slicing has an arrow of time given by nonzero currents when the observer has access to the full phase space distribution.

III. EXACT RG AND THE DENSITY MATRIX

We would now like to take our intuition from quantum walks and apply it to the reduced density matrix of the long wavelength modes in de Sitter. Specifically, we will be constructing a density matrix

$$\rho_R[\varphi_+(\vec{x}), \varphi_-(\vec{x}), t] = \text{Tr}_\Phi \rho[\phi_+, \phi_-], \quad \phi_\pm = \varphi_\pm + \Phi, \quad (3.1)$$

where $\varphi(\vec{x}, t)$ is a local field defined in terms of the Fourier modes $\phi(\vec{k}, t)$ with $k < aH$ and $\Phi(\vec{x}, t)$ is the local field composed of the hard modes with $k \geq aH$ that we want to trace over.

We have intuition from RG flow and EFT that suggests that the evolution of the reduced density matrix can be understood as the exact RG for the density matrix $\rho[\phi_+, \phi_-]$ after integrating out a shell of modes with $k \in [\Lambda, \Lambda + \delta\Lambda]$

where $\Lambda > aH$. For purely long wavelength observables, we are going to trace over all the short modes anyway and ρ_R is independent of $\Lambda \gg aH$, and plays the same role as the Euclidean partition function of an exact RG in flat space [40]. However, because of the expansion of the Universe, modes that are in $\rho_R(t)$ would not have been in $\rho_R(t')$ for $t' \ll t$. Therefore, understanding dynamical RG (RG plus time evolution) introduces subtleties beyond the usual exact RG. Nevertheless, we will find that by following analogous exact RG treatment for the density matrix in flat space [33,41] (see also Refs. [54,55]), we can derive the evolution of the density matrix in dS under time evolution. Our approach is related to previous work deriving the stochastic formalism using functional/exact RG [56–59].

A. Setup

We will start from the density matrix associated with the UV field ϕ , $\rho[\phi_+, \phi_-]$. This field is defined up to some physical cutoff Λ_0 . Our interest is not the UV density matrix, but instead the reduced density matrix of the long distance modes. We can do this at the level of the path integral by integrating out shells of momentum down to an effective cutoff $aH < \Lambda < \Lambda_0$. The density matrix of the IR field, φ , can be defined as a path integral over the contour $\gamma \in (-\infty(1-i\epsilon), t_0] \cup [t_0, -\infty(1+i\epsilon))$ with the boundary condition $\varphi(\vec{x}, t_\pm \rightarrow t_0) = \varphi_\pm(\vec{x})$ where $t_+ \in (-\infty(1-i\epsilon), t_0]$ and $t_- \in [t_0, -\infty(1+i\epsilon))$. We then express the density matrix in terms of the action as

$$\begin{aligned} \rho_R[\varphi_+, \varphi_-] &= \int_\gamma D\varphi T \exp(iS_0 + iS_{\text{int}} + iS_\partial), \\ S_0[\varphi(\vec{x}, t), \Lambda] &\equiv \int_\gamma dt a^3(t) \int \frac{d^3 k}{(2\pi)^3} \frac{1}{2} \nabla_\mu \varphi(\vec{k}, t) K^{-1} \\ &\quad \times \left(\frac{k^2}{\Lambda^2} \right) \nabla^\mu \varphi(\vec{x}, t), \end{aligned} \quad (3.2)$$

where Λ is the scale where we integrate out the short modes to define an EFT. The UV density matrix, $\rho[\phi_+, \phi_-]$ is defined in the same way, but with $\Lambda = \Lambda_0$. Our goal is to show that in integrating out a shell $[\Lambda, \Lambda + \delta\Lambda]$, the form of the density matrix does not change.

We will define ∇_μ as the covariant derivative and $\nabla^2 = \nabla_\mu \nabla^\mu$. To simplify notation, when acting on Fourier modes, it is implied that we also Fourier transformed the derivative, so that (for example)

$$\nabla_\mu \nabla^\mu f(\vec{k}, t) \equiv \left(\frac{\partial^2}{\partial t^2} + 3H \frac{\partial}{\partial t} + \frac{k^2}{a^2} \right) f(\vec{k}, t). \quad (3.3)$$

The additional terms in the action S_I and S_∂ are the bulk interaction and boundary terms respectively. The reduced density matrix will then be traced over all modes with $k > aH$ so we will have $\Lambda \geq aH(t)$.

For the density matrix, a boundary term is always required in order to have a well-defined action principle. At the quadratic level, this implies

$$S_{0,\partial} \supset - \int d^3x a^3(t) \frac{1}{2} (\varphi_+ K^{-1} \dot{\varphi}(t_{0,+}) - \varphi_- K^{-1} \dot{\varphi}(t_{0,-})), \quad (3.4)$$

where $t_{0,\pm}$ means we evaluate the derivative with respect to $t_{\pm}$ and then take $t_{\pm} \rightarrow t_0$.

B. Integrating out

Now we will proceed to integrate out a shell of momentum. Because we are performing the path integral over $t < t'$, any change to the density matrix as we change Λ must arise from fields at time t . We will show that all such effects can be captured by the boundary action, S_{∂} so that

$$\begin{aligned} \frac{\partial}{\partial s} \rho_R[\varphi_+, \varphi_-, t] &= i \int_{\gamma} D\varphi \left(\frac{\partial}{\partial s} S_0 + \frac{\partial}{\partial s} S_I + \frac{\partial}{\partial s} S_{\partial} \right) \\ &\times \exp(iS_0 + iS_{\text{int}} + iS_{\partial}) = 0, \end{aligned} \quad (3.5)$$

where $s \equiv \log \Lambda$. As in the case of exact RG, we are looking to define the change to the interaction and boundary terms that will make the vanishing of this expression manifest. At this stage, Eq. (3.5) is a conjecture, as we must demonstrate that any dependence of the density matrix on Λ can be absorbed into S_{∂} .

The key idea is to use our knowledge of the quadratic (bulk) action to simplify this expression. In order to simplify the notation, we will suppress the integral over Fourier modes and isolate the integration over time. Quantities that are integrated over same time variable we will write in the compact notation

$$\begin{aligned} f * g &\equiv \int dt f(t) g(t), \\ f * H * g &\equiv \int dt \int dt' f(t) H(t, t') g(t'). \end{aligned} \quad (3.6)$$

We will keep track of time integration in this way in order to isolate the effect of the nontrivial boundary conditions imposed at $t = t_{0,\pm}$.

1. Quadratic terms

Now we can use our knowledge of the free theory to write

$$\frac{\partial}{\partial s} S_0 = \frac{1}{2} \frac{\delta}{\delta \varphi} S_0 * G' * \frac{\delta}{\delta \varphi} S_0 + \delta S_{\partial,2}, \quad (3.7)$$

where we defined the Green's function $G(t, t')$ so that

$$\begin{aligned} G^{-1} &= K^{-1} \nabla_{\mu} \nabla^{\mu}, \quad G^{-1} G = a^{-3}(t) \delta(t - t') \delta(\vec{x} - \vec{x}'), \\ G' &= \frac{\partial}{\partial s} G = \left(\frac{\partial}{\partial s} K \right) G. \end{aligned} \quad (3.8)$$

The boundary term $\delta S_{\partial,2}$ is determined using integration by parts. Specifically, we can write

$$\frac{\partial}{\partial s} S_0 = a^3(t) \frac{K'}{K^2} \varphi \nabla^2 \varphi = (K^{-1} \varphi \nabla^2) * G' * (K^{-1} \nabla^2 \varphi), \quad (3.9)$$

using the definition of the Green's function. The correction to the quadratic boundary term comes from integrating Eq. (3.9) by parts until it takes the form of Eq. (3.7). Defining $\bar{S}_0 = S_0 + S_{0,\partial}$ and using

$$\frac{\delta}{\delta \varphi} \bar{S}_0 = K^{-1} a^3(t) \nabla^2 \varphi, \quad (3.10)$$

we can see that

$$\frac{\delta}{\delta \varphi} \bar{S}_0 * G' * \frac{\delta}{\delta \varphi} \bar{S}_0 = K^{-1} \varphi \nabla^2 * G' * K^{-1} \nabla^2 \varphi - K^{-1} (\nabla_0 \varphi G' - \varphi \nabla_0 G')|_{t_{0,-}}^{t_{0,+}} * \frac{\delta}{\delta \varphi} \bar{S}_0. \quad (3.11)$$

As written yet, this is not a boundary term because it still involves a time integral through the $\delta_{\varphi} \bar{S}_0$ term. Fortunately we can integrate the second term by parts, and after simplification we find that only a boundary term remains of the form $\delta S_{\partial,0} = \int d^3x a^6(t_0) K^{-2} \frac{1}{2} f(\varphi)$ where

$$\begin{aligned} f &= (\varphi_- - \varphi_+) (\nabla_t G'(t_+) \nabla_{t'} \varphi(t_-) - \nabla_t G'(t_-) \nabla_{t'} \varphi(t_+) - \nabla_t \nabla_{t'} G'(\varphi_- - \varphi_+)) \\ &\quad - (\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+)) (G'(\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+)) - \nabla_{t'} G'(t_+) \varphi_- + \nabla_{t'} G'(t_-) \varphi_+). \end{aligned} \quad (3.12)$$

In writing this expression, we anticipated that

$$\begin{aligned} iG'(t, t', \vec{p})|_{t=t'=t_0} &\equiv \Delta_0(p), \\ i\nabla_t \nabla_{t'} G(t, t', \vec{p})|_{t=t'=t_0} &\equiv \Delta_2(p) \end{aligned} \quad (3.13)$$

are independent of taking $t_{\pm} \rightarrow t_0$. To determine the remaining terms, we note that $G(t, t', \vec{p})$ is the path ordered Green's function along γ ,

$$\begin{aligned} G'(t, t', \vec{p}) &= i\langle T_{\gamma} \varphi(\vec{p}, t_1) \varphi(-\vec{p}, t_2) \rangle' \\ &= i \frac{dK}{ds} \frac{H^2}{2p^3} (1 - ip\tau_1)(1 + ip\tau_2) e^{ip(\tau_1 - \tau_2)}, \end{aligned} \quad (3.14)$$

where t_1 (t_2) is the early (late) time along the contour γ and τ_i is the conformal time associated with physical time t_i . As a result, we have, for the massless field

$$\Delta_0(p, t_0) = iG' = iG'(t_0, t_0, \vec{p}) = -\frac{\partial K}{\partial s} \frac{H^2}{2p^3} (1 + \tau_0^2 p^2), \quad (3.15)$$

$$\Delta_2(p, t_0) = \partial_t \partial_{t'} iG'(t_0, t_0, \vec{p}) = -\frac{\partial K}{\partial s} \frac{H^4}{2p^3} \tau_0^4 p^4. \quad (3.16)$$

The single derivatives depend on the order, and the order of limits that we take as $t, t' \rightarrow t_0$ matters, and we find

$$\begin{aligned} i\partial_t G'(t, t_0, \vec{p})|_{t=t_{\pm}} &= i\partial_{t'} G'(t_0, t', \vec{p})|_{t'=t_{\mp}} \\ &= -\frac{\partial K}{\partial s} \frac{H^3}{2p^3} (\tau_0^2 p^2) (1 \pm i\tau_0 p) \\ &\equiv \Delta_R \pm i\Delta_I, \end{aligned} \quad (3.17)$$

where

$$\begin{aligned} \Delta_R &= -K' \frac{H^3}{2p^3} (p^2 \tau_0^2), \\ i\Delta_I &= iK^{-1} K' \frac{H^3}{2p^3} (\tau_0^3 p^3) = -i \frac{K'}{2a^3}. \end{aligned} \quad (3.18)$$

The final result is that

$$\begin{aligned} if &= -\Delta_2(\varphi_- - \varphi_+)^2 - \Delta_0(\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+))^2 \\ &+ \Delta_R\{(\varphi_+ - \varphi_-), (\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+))\} \\ &+ i\Delta_I[(\varphi_- - \varphi_+)(\nabla_t \varphi_- + \nabla_t \varphi_+) \\ &- (\nabla_t \varphi_- - \nabla_t \varphi_+)(\varphi_+ + \varphi_-)]. \end{aligned} \quad (3.19)$$

These additional quadratic terms will play a significant role in stochastic inflation.

The corrections to the boundary action will be given two different interpretations but are equivalent. The first, in precise analogy with exact RG, is that we change the quadratic boundary action to absorb this term

$$\frac{\partial}{\partial s} S_{\partial} \supset -if. \quad (3.20)$$

This interpretation is most useful when discussing the resummation of logs in the sense of RG flow. Although this is a local boundary term, it is unfortunately not equivalent to unitary evolution. Specifically, we have mixed $\varphi_+ - \varphi_-$ terms that we associate with open quantum systems. The second, and more typical, interpretation of the boundary terms is to treat them as a change to the density matrix,

$$\frac{\partial}{\partial s} \rho_R \supset if \rho_R. \quad (3.21)$$

Here if is a local superoperator in the $\varphi_{\pm}$ basis. This interpretation will be most useful for describing the density matrix during stochastic inflation.

Finally, the running of the required quadratic boundary terms of the action will also contribute a correction to the density matrix via

$$\frac{\partial}{\partial s} S_{0,\partial} = a^3(t) \frac{K^{-2}}{2} \left(\frac{\partial K}{\partial s} \right) (\varphi_+ \nabla_t \varphi_+ - \varphi_- \nabla_t \varphi_-). \quad (3.22)$$

Using the observation that

$$\Delta_I = \frac{H^3 \tau_0^3}{2} \frac{\partial K}{\partial s} = -\frac{\partial K}{\partial s} \frac{1}{2a^3}, \quad (3.23)$$

we can rewrite

$$i \frac{\partial}{\partial s} S_{0,\partial} = a^6(t) \Delta_I K^{-2} (\varphi_+ \nabla_t \varphi_+ - \varphi_- \nabla_t \varphi_-). \quad (3.24)$$

Since $[\varphi, \nabla_t \varphi] \neq 0$ we need to prescribe a time ordering. The correct time ordering is the symmetrized one, which gives

$$\begin{aligned} i \frac{\partial}{\partial s} S_{0,\partial} &= a^6(t) \Delta_I \frac{K^{-2}}{2} (\varphi_+ \nabla_t \varphi_+ + \nabla_t \varphi_+ \varphi_+ \\ &- \varphi_- \nabla_t \varphi_- - \nabla_t \varphi_- \varphi_-). \end{aligned} \quad (3.25)$$

Now we can combine all the terms to find

$$\begin{aligned} \left[\frac{1}{2} if + iK^2 \frac{\partial}{\partial s} S_{0,\partial} \right] &= -\frac{1}{2} \Delta_2(\varphi_- - \varphi_+)^2 - \frac{1}{2} \Delta_0(\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+))^2 \\ &+ \frac{1}{2} \Delta_R\{(\varphi_+ - \varphi_-), (\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+))\} - i\Delta_I(\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+))(\varphi_+ + \varphi_-). \end{aligned}$$

Overall, we find that the evolution of the density matrix due to the quadratic terms takes the form

$$\frac{\partial}{\partial s} \rho_R \supset i a^6 K^{-2} \frac{f}{2} \rho_R + \frac{\partial S_{0,\partial}}{\partial s}, \quad (3.26)$$

where we have suppressed the implicit momentum integral.

2. Interaction terms

After determining the quadratic boundary terms, we then proceed to use the Schwinger-Dyson (SD) equations to eliminate S_0 . Specifically, at the operator level, the SD equations reflect our ability to integrate by parts so that

$$\begin{aligned} i \int_{\gamma} D\varphi \left(\mathcal{O} \frac{\delta}{\delta \varphi} S_0 \right) e^{i(S_0 + S_I)} \\ = - \int_{\gamma} D\varphi \left(i \mathcal{O} \frac{\delta}{\delta \varphi} S_I + \frac{\delta}{\delta \varphi} \mathcal{O} \right) e^{i(S_0 + S_I)}, \end{aligned} \quad (3.27)$$

where $S_I = S_{\text{int}} + S_{\partial}$. This equation holds for $\varphi(t \neq t_0)$ because we are integrating over all fields up to the boundary conditions at t_0 . One should worry about the potential contributions at $t = t_0$; fortunately these are manifestly pure boundary terms and can be determined directly.

At this stage, there is no practical difference between the dS and the flat space in-in or density matrix calculations [33,41], other than the precise form of G . Therefore, we apply the SD equations repeatedly, to write

$$\begin{aligned} \frac{1}{2} \frac{\delta S_0}{\delta \varphi} \star G' \star \frac{\delta S_0}{\delta \varphi} \\ = - \frac{1}{2} \int_{\gamma} dt \left(\frac{\delta^2 S_0}{\delta \varphi \delta \varphi} G'(t, t) \right) + \frac{1}{2} \frac{\delta S_0}{\delta \varphi} \star G' \star \frac{\delta S_I}{\delta \varphi} \\ = \frac{1}{2} \int_{\gamma} dt \left(\frac{\delta^2 S_I}{\delta \varphi \delta \varphi} G'(t, t) \right) - \frac{1}{2} \frac{\delta S_I}{\delta \varphi} \star G' \star \frac{\delta S_I}{\delta \varphi} \\ - \frac{1}{2} \int_{\gamma} dt \left(\frac{\delta^2 S_0}{\delta \varphi \delta \varphi} G'(t, t) \right). \end{aligned} \quad (3.28)$$

The term on the last line is a constant and just contributes to the overall normalization of ρ (which is fixed, in principle, by the trace condition, $\text{Tr} \rho = 1$). What remains are the two (well-known) contributions to the effective action from integrating out a shell of momentum, namely

$$\frac{\partial}{\partial s} S_I = - \frac{1}{2} \int_{\gamma} dt \left(\frac{\delta^2 S_I}{\delta \varphi \delta \varphi} G'(t, t) \right) + \frac{1}{2} \frac{\delta S_I}{\delta \varphi} \star G' \star \frac{\delta S_I}{\delta \varphi}. \quad (3.29)$$

The first term is purely a local function of t because there is a single integral and therefore contributes only to S_{int} . The second term includes both local and boundary terms. Specifically, because $G'(t, t)$ is exponentially suppressed

$$\frac{d}{d \log \Lambda} S_I = \left(\text{purple circle } t \text{ --- } \text{yellow circle } t' + \text{blue circle with self-loop} \right)$$

$\frac{\delta}{\delta \varphi} S_I$ $\frac{\delta}{\delta \varphi} S_I$ $\frac{\delta^2}{\delta \varphi^2} S_I$

FIG. 1. The equation for the evolution of the effective action under dynamical RG. The key difference between this diagrammatic representation and conventional exact RG is that the times t and t' are not necessarily on the same branch of the Schwinger-Keldysh contour, giving rise to additional boundary terms.

in $t - t'$ (or, as we will see, $t - t_0$ or $t' - t_0$), we can Taylor expand and integrate to generate local operators.

However, since the local action can only involve operators in one of the $\gamma_+ = t_+ \in (-\infty(1 - i\epsilon), t_0]$ or $\gamma_- = t_- \in [t_0, -\infty(1 + i\epsilon)]$ regions, we may still have terms mixing t_+ and t_- that do not arise in the bulk action, see Fig. 1 for illustration. Therefore, defining

$$S_{\text{int},\pm} = \int_{\gamma_{\pm}} dt_{\pm} \mathcal{L}, \quad (3.30)$$

the correction to the bulk (local) action is

$$\begin{aligned} \frac{\partial}{\partial s} S_{\text{int},\pm} = - \frac{1}{2} \int_{\gamma} dt \left(\frac{\delta^2 S_{\text{int},\pm}}{\delta \varphi \delta \varphi} G'(t_{\pm}, t_{\pm}) \right) \\ + \frac{1}{2} \frac{\delta S_{\text{int},\pm}}{\delta \varphi} \star G'_{\pm,\pm} \star \frac{\delta S_{\text{int},\pm}}{\delta \varphi}. \end{aligned} \quad (3.31)$$

Allowing for a boundary action $S_{\partial}(\varphi_+, \varphi_-)$, we then find boundary terms

$$\frac{\partial}{\partial s} S_{\partial,\text{int}} = \frac{\delta S_{\text{int},+}}{\delta \varphi} \star G'_{+,-} \star \frac{\delta S_{\text{int},-}}{\delta \varphi}. \quad (3.32)$$

It will generally be the case that $G_{+,-}$ is exponentially suppressed in $t_{\pm} - t_0$ and therefore

$$\frac{\partial}{\partial s} S_{\partial,\text{int}} = \int d^3x \sum_{i,j} \mathcal{O}_{+,i} \mathcal{O}_{-,j}, \quad (3.33)$$

where $\mathcal{O}_{\pm}$ are local operators of $\varphi(t_{\pm})$ defined at $t_{\pm} \rightarrow t_0$ respectively that can be derived from the Taylor expansion of Eq. (3.32) at $t_+ = t_- = t_0$. We see this term necessarily mixes $\varphi(t_{\pm})$ and therefore is not associated with conventional Hamiltonian evolution. However, because this is a pure boundary term, it can be interpreted as a redefinition of the operators at t_0 .

Like the quadratic boundary terms another way to interpret mixed boundary terms is through the evolution of the density matrix. Specifically, S_{∂} acts as a superoperator on density matrix and therefore can be incorporated through the master equation

$$\frac{\partial}{\partial s} \rho_R \supset \sum_{i,j} \int d^3x \mathcal{O}_{+,i} \rho_R \mathcal{O}_{-,j} \equiv \mathbb{S}_{\text{int}} \rho_R, \quad (3.34)$$

where $\mathbb{S}_{\text{int}}$ is the superoperator in the Kraus representation. In this representation, all corrections to the effective action in S_{int} can be derived from any observable as they do not depend on t_0 or the fact that this is an in-in correlator. All terms specific to the in-in correlator at t_0 contribute to the evolution of the density matrix.

While the presence of the additional boundary terms will be important for cosmological applications, such as

stochastic inflation, it is important to note that they are needed in flat space when matching correlators [33]. The fact that they do not appear in matching S-matrix elements is a consequence of the fact that they can represent field redefinitions between the UV and IR descriptions.

3. Full density matrix RG

To summarize the exact RG calculation, integrating out a shell of momenta introduces a differential change to the reduced density matrix of the long wavelength modes,

$$\begin{aligned} \frac{\partial \rho_R}{\partial s} &= \int \frac{d^3p}{(2\pi)^3} a^6(t) i K^{-2} \left[\frac{1}{2} f(\varphi) \rho_R + \frac{\partial S_{0,\partial}}{\partial s} \right] \\ &= \int \frac{d^3p}{(2\pi)^3} a^6(t) K^{-2} \left[-\frac{1}{2} \Delta_2 (\varphi_- - \varphi_+)^2 - \frac{1}{2} \Delta_0 (\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+))^2 \right. \\ &\quad \left. + \frac{1}{2} \Delta_R \{ (\varphi_+ - \varphi_-), (\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+)) \} - i \Delta_I (\nabla_t \varphi(t_-) - \nabla_t \varphi(t_+)) (\varphi_+ + \varphi_-) \right] \rho_R + \mathbb{S}_{\text{int}} \rho_R, \end{aligned} \quad (3.35)$$

where $\mathbb{S}_{\text{int}}$ are higher order local terms defined in Eq. (3.34). We can further simplify the quadratic terms in this equation by recognizing that

$$a^3(t) K^{-1} \nabla_t \varphi(t_+) \rho_R = -i \frac{\delta}{\delta \varphi(t_+)} \rho_R \quad a^3(t) K^{-1} \nabla_t \varphi(t_-) \rho_R = i \frac{\delta}{\delta \varphi(t_-)} \rho_R, \quad (3.36)$$

leading to

$$\begin{aligned} \frac{\partial \rho_R}{\partial s} &= \frac{1}{2} \int \frac{d^3p}{(2\pi)^3} \left[\Delta_0 \left(\frac{\delta}{\delta \varphi_+} + \frac{\delta}{\delta \varphi_-} \right)^2 - a^6 K^{-2} \Delta_2 (\varphi_- - \varphi_+)^2 \right. \\ &\quad \left. + i \frac{a^3}{K} \Delta_R \left\{ (\varphi_+ - \varphi_-), \left(\frac{\delta}{\delta \varphi_+} + \frac{\delta}{\delta \varphi_-} \right) \right\} + 2 \frac{a^3}{K} \Delta_I \left(\frac{\delta}{\delta \varphi_+} + \frac{\delta}{\delta \varphi_-} \right) (\varphi_+ + \varphi_-) \right] \rho_R + \mathbb{S}_{\text{int}} \rho_R. \end{aligned} \quad (3.37)$$

Here it will be important that this is a functional equation as we have suppressed the momentum, $\varphi = \varphi(\vec{p})$.

At this point, it is instructive to compare this to the random walks that we have built intuition for in Sec. II. Schematically, the quadratic terms (i.e. dropping $\mathbb{S}_{\text{int}}$) in our evolution equation are of the form

$$\begin{aligned} \frac{\partial \rho_R}{\partial s} &= \Delta_0 \left(\frac{\partial}{\partial x} + \frac{\partial}{\partial x'} \right)^2 \rho_R + \Delta_2 (x - x')^2 \rho - i \Delta_R \left(\frac{\partial}{\partial x} + \frac{\partial}{\partial x'} \right) (x - x') \rho_R \\ &\quad + \frac{\Delta_I}{2} \left(\frac{\partial}{\partial x} + \frac{\partial}{\partial x'} \right) (x + x') \rho_R. \end{aligned} \quad (3.38)$$

As we commented on earlier, it is easier to gain intuition for random walks in the Wigner phase space formalism. Transforming the above equation to the Wigner representation, we get

$$\frac{\partial W}{\partial t} = \Delta_0 \frac{\partial^2}{\partial x^2} W + \Delta_2 \frac{\partial^2}{\partial p^2} W + \Delta_R \frac{\partial^2}{\partial x \partial p} W + \Delta_I \frac{\partial}{\partial x} (x W). \quad (3.39)$$

Looking at this equation, we can quickly identify what the different Δ represent. The Δ_0, Δ_2 terms represent the diffusion in x, p , while the Δ_R term tells us about the degree of correlation between the two sources of diffusion. The Δ_I term, represents a quantum drift and arises from the fact the conjugate momentum to φ depends on K , and hence Λ . All in all, we see clearly how on integrating out a shell of momenta, we recreate the key structure of random walks. However, it is essential for our

discussion that Eq. (3.37) is a functional equation and not an ODE [i.e. it is a differential equation for $\varphi(\vec{k})$]. There remains much work to translate the RG equation into the partial differential equations that describe stochastic inflation.

C. Combining time evolution and RG

The actual problem of interest is to integrate in a shell of comoving momentum $p \in [\Lambda, \Lambda + \Delta\Lambda]$ and evolve forward in time $a(t) \rightarrow a(t + \delta t) \approx a(t)(1 + H\delta t)$ so that the physical UV cutoff is unchanged,

$$\Lambda_{\text{phys}} = \frac{\Lambda}{a(t)} = \frac{\Lambda + \delta\Lambda}{a(t + \delta t)} \rightarrow \delta t \approx H^{-1} \frac{\delta\Lambda}{\Lambda}. \quad (3.40)$$

The nature of this transformation is illustrated in Fig. 2. Under time evolution the reduced density matrix evolves according to

$$\frac{d}{Hdt} \rho_R = \frac{\partial}{H\partial t} \rho_R + \frac{\partial}{\partial s} \rho_R, \quad (3.41)$$

where $s = \log \Lambda$ as before. The first term is just the Hamiltonian evolution. Assuming our scalar field is a canonical field evolving on a potential, the Hamiltonian for the action (3.2) with a potential is

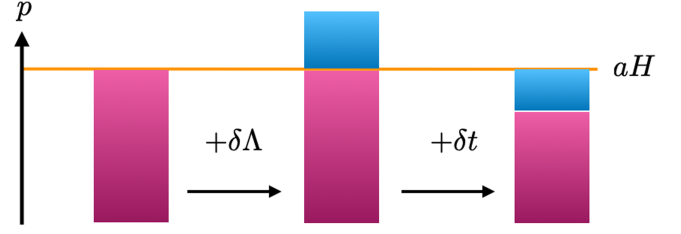


FIG. 2. The combination of RG and time evolution. The combined transformation holds the scale of the comoving cutoff, aH , fixed.

$$\hat{H} = \int \frac{d^3 p}{(2\pi)^3} \left[K \left(\frac{p^2}{\Lambda^2} \right) \left[\frac{1}{2a^3} \hat{\Pi}^2 \right] + K^{-1} \left(\frac{p^2}{\Lambda^2} \right) \left[\frac{a^3}{2} \frac{(\nabla \varphi)^2}{a^2} \right] + a^3 V(\varphi) \right] + \hat{H}_{\text{EFT}}, \quad (3.42)$$

where $\hat{H}_{\text{EFT}}$ includes any higher derivative terms. The Hamiltonian is responsible for time evolution in the usual way,

$$\frac{\partial}{\partial t} \rho_R = -i[\hat{H}, \rho_R]. \quad (3.43)$$

Writing this in the $\varphi_{\pm}$ basis, and dropping H_{EFT} for now, we find

$$\dot{\rho}_R = \int \frac{d^3 p}{(2\pi)^3} \left[\frac{iK}{2a^3} \left(\frac{\delta^2}{\delta\varphi_+^2} - \frac{\delta^2}{\delta\varphi_-^2} \right) - \frac{ia^3}{2K} \left(\frac{(\nabla\varphi_+)^2}{a^2} - \frac{(\nabla\varphi_-)^2}{a^2} \right) - ia^3(V(\varphi_+) - V(\varphi_-)) \right] \rho_R. \quad (3.44)$$

Now, we can combine one step of RG and one step of time evolution to get one step of dynamical RG,

$$\begin{aligned} \frac{d}{dt} \rho_R = & \frac{H}{2} \int \frac{d^3 p}{(2\pi)^3} \left[\Delta_0 \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right)^2 - \frac{a^6}{K^2} \Delta_2 (\varphi_- - \varphi_+)^2 \right. \\ & + i \frac{a^3}{K} \Delta_R \left\{ (\varphi_+ - \varphi_-), \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) \right\} + 2 \frac{a^3}{K} \Delta_I \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) (\varphi_+ + \varphi_-) \left. \right] \rho_R \\ & + \int \frac{d^3 p}{(2\pi)^3} \left[\frac{iK}{2a^3} \left(\frac{\delta^2}{\delta\varphi_+^2} - \frac{\delta^2}{\delta\varphi_-^2} \right) \rho_R - \frac{ia^3}{2K} \left(\frac{(\nabla\varphi_+)^2}{a^2} - \frac{(\nabla\varphi_-)^2}{a^2} \right) \rho_R \right. \\ & \left. - ia^3(V(\varphi_+) - V(\varphi_-)) \rho_R \right] - i[H_{\text{EFT}}, \rho_R] + \$_{\text{int}} \rho_R. \end{aligned} \quad (3.45)$$

For completeness, we reintroduced $\hat{H}_{\text{EFT}}$ arising from the higher derivative contributions to S_{eff} , and $\$_{\text{int}}$ is the interaction superoperator defined in Eq. (3.34). In practice, these terms will not play a meaningful role in stochastic inflation at leading or subleading order. Therefore, we will drop $\hat{H}_{\text{EFT}}$ and $\$_{\text{int}}$ from our analysis. The remaining terms play the dominant role in the exact RG equation for the density matrix that we will analyze in the majority of the paper.

D. From modes to local observables

One of the key issues in understanding the evolution of the density matrix is determining the operators that represent the long distance observations. To illustrate this issue and how it relates to the density matrix, let us first assume that we are in a free theory so that $S_{\text{int}} = 0$. Due to the boundary terms, the density matrix still evolves according to

$$\frac{\partial}{\partial s} \rho_R \supset \int \frac{d^3 p}{(2\pi)^3} \Delta_0(p) \left(\frac{\partial}{\partial\varphi_+(\vec{p})} + \frac{\partial}{\partial\varphi_-(-\vec{p})} \right) \left(\frac{\partial}{\partial\varphi_+(-\vec{p})} + \frac{\partial}{\partial\varphi_-(-\vec{p})} \right) \rho_R[\varphi_+, \varphi_-], \quad (3.46)$$

where φ_+ and φ_- are the complete set of Fourier modes with $p, q < \Lambda(t)$. However, since $\Delta_0(p)$ only has support over a range of momenta $p \in [\Lambda, \Lambda + \Delta\Lambda]$, the right-hand side is a diffusion term only for modes near the cutoff. Concretely, dropping the other terms, we would get a diffusion equation of the form

$$\begin{aligned} \frac{\partial}{H\partial t}\rho_R &= \delta(p - \Lambda(t)) \left(\frac{\partial}{\partial\varphi_+(\vec{p})} + \frac{\partial}{\partial\varphi_-(\vec{p})} \right) \\ &\times \left(\frac{\partial}{\partial\varphi_+(-\vec{p})} + \frac{\partial}{\partial\varphi_-(-\vec{p})} \right) \rho_R. \end{aligned} \quad (3.47)$$

Assuming the initial density matrix is a product between the UV and the IR modes, $\rho_0 = \rho_0(\varphi_+, \varphi_-) \otimes \rho_0(\Phi_+, \Phi_-)$, then after time $t > t_0$, we will have

$$\begin{aligned} \rho_R(\varphi_+, \varphi_-) &= \text{Tr}_{\Phi} \rho(\phi_+ = \varphi_+ + \Phi, \phi_- = \varphi_- + \Phi) \\ &= \rho_0(\varphi_+, \varphi_-). \end{aligned} \quad (3.48)$$

In other words, our diffusion equation does not couple modes and therefore cannot change the state of the long wavelength modes, including generating a mixed state from a pure state. This is true of all of the quadratic boundary terms. Of course, this is a trivial statement in terms of the wave function for a free theory,

$$\begin{aligned} \Psi[\phi, t] &= Z \exp \left(i \int d^3k \frac{(aH)^3}{2H^2} \left(\frac{k^2}{(aH)^2} + i \frac{k^3}{(aH)^3} \right) \right. \\ &\quad \times \phi(-\vec{k})\phi(\vec{k}) \Big) \\ &= \prod_k \psi[\phi(\vec{k})]. \end{aligned} \quad (3.49)$$

Here we are making the well-known observation that in a free translationally invariant theory, the Fourier modes are all independent and the wave function is a product state over these modes. As a result, tracing over any set of UV Fourier modes cannot create a mixed state for the remaining IR modes.

The description in terms of stochastic inflation and the resulting mixed state for the universe is therefore not just the result of integrating out short distance physics but also limiting our measurements to a local region. A common approach is to trace over the region outside of a sphere [60], giving rise to nontrivial density matrix due to entanglement in space.

In contrast, we will focus on local measurements of the long-wavelength field $\varphi(\vec{x})$ and its conjugate momentum $\pi(\vec{x})$ and trace over information that leaves $\varphi(\vec{x})$ [or $\pi(\vec{x})$] fixed. Using translation invariance, we can choose this point to be $\vec{x} = 0$ and

$$\varphi(\vec{x} = 0) = \bar{\varphi} = \int \frac{d^3k}{(2\pi)^3} \varphi(\vec{k}), \quad (3.50)$$

$$\pi(\vec{x} = 0) = \bar{\pi} = \frac{1}{N_k} \int \frac{d^3k}{(2\pi)^3} K \left(\frac{k^2}{\Lambda^2} \right) \pi(\vec{k}), \quad (3.51)$$

where $N_k = \int \frac{d^3k}{(2\pi)^3} K$ is the number of modes, that we introduced so that $[\bar{\varphi}, \bar{\pi}] = i$. We note that these are two local operators we can measure with a local detector. We should think of this in exact analogy with our quantum walk in Eq. (2.3); here $\prod_{\vec{k}} \psi[\phi(\vec{k})]$ is a product state of harmonic oscillators. It is only when we coupled $\sum \varphi(\vec{k})$ to an observer (X in the case of the quantum walk) and trace over the harmonic oscillators that we see the mixed state emerge.

Rather than explicitly calculating the trace over $\varphi(\vec{p})$ that leaves $\varphi(0)$ fixed, we can calculate the generating functional for φ ,

$$Z[J_{\pm}, t] = \text{Tr}[e^{J_+ \bar{\varphi}} e^{J_- \bar{\pi}} \rho_R(\varphi_+(\vec{x}), \varphi_-(\vec{x}))]. \quad (3.52)$$

Here it is important that we are working in the Schrödinger picture, so that all the time evolution is in the density matrix. The time evolution of this partition function is therefore given by

$$\frac{d}{dt} Z[J_{\pm}, t] = \text{Tr}[e^{J_+ \bar{\varphi}} e^{J_- \bar{\pi}} \frac{d}{dt} \rho_R(\varphi_+(\vec{x}), \varphi_-(\vec{x}))]. \quad (3.53)$$

This representation is useful because it immediately connects many of the previous perspectives on this system. In short, by allowing $J_{\pm} = ik_{\pm}$, we can write $Z[J_{\pm}, t]$ as the density matrix, Wigner distribution, or other quantities associated with the variables $\bar{\varphi}$ and $\bar{\pi}$. See Appendix A for derivations.

We can now plug in the evolution equation for the density matrix (3.45), and simplify to get the evolution equation for Z . The details of the calculation can be found in Appendix B. For the massless field, we end up with

$$\begin{aligned} \frac{dZ[J_{\pm}]}{dt} &= \frac{1}{2} \int \frac{d^3p}{(2\pi)^3} \left[\left(\frac{H^3}{2p^3} J_+^2 - a^2(t_0) H \frac{p}{2} \frac{J_-^2}{N_k^2} - \frac{aH^2}{p} \frac{J_- J_+}{N_k} \right) \frac{\partial K}{\partial s} \right] Z[J_{\pm}] \\ &\quad + \frac{N_k}{a^3} J_+ \left(-\frac{iJ_+}{2} + \frac{\partial}{\partial J_-} \right) Z[J_{\pm}] \\ &\quad - \frac{J_-}{N_k} a^3 \tilde{V}'(-iJ_+) Z[J_{\pm}] + \int \frac{d^3p}{(2\pi)^3} \left(K \frac{aH^2}{2p} \frac{J_+ J_-}{N_k} - \frac{a^2 H p}{2} K \frac{J_-^2}{N_k^2} \right) Z[J_{\pm}]. \end{aligned} \quad (3.54)$$

While this equation still looks like a mess, it is a nice basis independent way of writing down the evolution of observables we measure in stochastic inflation. We will convert this equation to more well-known formulas in the next section.

IV. LIGHT FIELDS AND STOCHASTIC INFLATION

The description of long modes in cosmology has long been understood to mirror a random walk. In particular, as modes cross the horizon, they become classical in the sense that for a given Fourier mode, the commutator vanishes as $a \rightarrow \infty$,

$$[\varphi(\vec{k}), \varphi(\vec{k}')] = iH^2(aH)^{-3}\delta(\vec{k} + \vec{k}'). \quad (4.1)$$

Therefore, in an approximate sense, we must be able to treat $\varphi(\vec{k})$ as a classical statistical variable when $k \ll aH$. This intuition can be made precise using soft de Sitter effective theory [13], where quantum mechanical effects are irrelevant by power counting.

If we start from a free theory, where each Fourier mode is independent, then $\varphi(\vec{k})$ is like a i.i.d. random variable. Naturally, we would expect that the variable,

$$\bar{\varphi} = \varphi(\vec{x} = 0) = \int \frac{d^3k}{(2\pi)^3} \varphi(\vec{k}), \quad (4.2)$$

will behave like the location of a random walker. The goal of this section is to show precisely how this intuition arises within a fully quantum mechanical description, in analogy with the quantum walks from Sec. II.

Physically, we should interpret $\bar{\varphi}$ as the local measurement of the background value of the field at the point $\vec{x} = 0$. For example, we could introduce a local observer with a detector whose value is measured by $\hat{X}$ in a pointer state [53],

$$|X, \varphi; t\rangle = e^{i\hat{P}\bar{\varphi}}|0\rangle \times |\Psi[\phi]\rangle. \quad (4.3)$$

Tracing over the short distance modes then gives our random walk

$$\rho(X, X') = \text{Tr}[e^{i\hat{P}\bar{\varphi}}|0\rangle\langle 0|e^{-i\hat{P}\bar{\varphi}} \otimes \rho_R(\varphi_+, \varphi_-)]. \quad (4.4)$$

Of course, local measurements are also sensitive to short wavelength. However, we can isolate only the constant long wavelength modes using a local detector model [44] that can discriminate between time dependence or wavelength. The need for the local measurement is in line with recent work that shows the need for including the observer in order to make sense of physics in dS [19–21, 61–64]. This connection between the density matrix and a local observer is illustrated in Fig. 3.

A. Leading order

The general strategy we will use in this section is to write the evolution equation in terms of the Wigner functional. As usual in the case of quantum random walks, the Wigner description most naturally makes contact with classical dynamical systems.

The advantage of defining the generating functional $Z[J]$ in Eq. (3.52) is that it is equivalent to many representations, including the Wigner functional. Specifically, as shown in Appendix A, we can write

$$\begin{aligned} W[\bar{\varphi}, \bar{p}] &= \int d\Delta\varphi e^{-i\bar{p}\Delta\varphi} \\ &\times \int \frac{dk}{(2\pi)} e^{i\frac{k\Delta\varphi}{2}} e^{-ik\bar{\varphi}} Z[J_+ = ik, J_- = i\Delta\varphi]. \end{aligned} \quad (4.5)$$

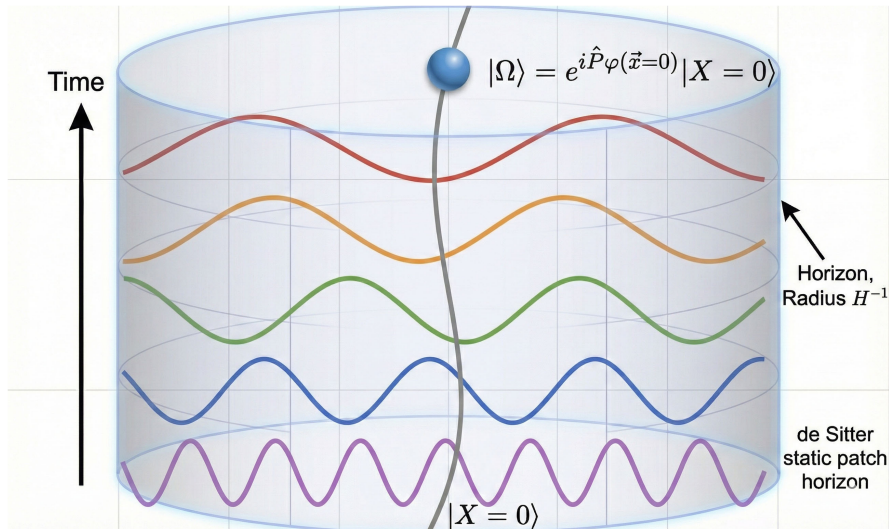


FIG. 3. Visualization of a local observer measuring $\bar{\varphi}$ with a detector along a worldline.

Integration by parts will allow us to convert the evolution equation for $Z[J_{\pm}]$ into an evolution equation for the Wigner distribution,

$$\begin{aligned} \frac{dZ[J_{\pm}]}{dt} = & \frac{1}{2} \int \frac{d^3 p}{(2\pi)^3} \left[\left(\frac{H^3}{2p^3} J_+^2 - a^2(t_0) H \frac{p(J_-)^2}{2N_k^2} - \frac{aH^2 J_- J_+}{p N_k} \right) \frac{\partial K}{\partial s} \right] Z[J_{\pm}] + \frac{N_k}{a^3} J_+ \left(-\frac{iJ_+}{2} + \frac{\partial}{\partial J_-} \right) Z[J_{\pm}] \\ & - \frac{J_-}{N_k} a^3 \tilde{V}'(-iJ_+) Z[J_{\pm}] + \int \frac{d^3 p}{(2\pi)^3} \left(K \frac{aH^2 J_+ J_-}{2p N_k} - \frac{a^2 H p}{2} K \frac{J_-^2}{N_k^2} \right) Z[J_{\pm}]. \end{aligned} \quad (4.6)$$

Since the Wigner distribution is closely related to the Fourier transform of $Z[J_{\pm}]$, the evolution equation from $Z[J_{\pm}] \rightarrow W[\bar{\varphi}, \bar{p}]$ can be determined using

$$J_+ = ik \rightarrow -\frac{\partial}{\partial \bar{\varphi}}, \quad J_- = i\Delta\varphi \rightarrow -\frac{\partial}{\partial \bar{p}}, \quad \frac{\partial}{\partial J_-} - i\frac{J_+}{2} \rightarrow \bar{p}. \quad (4.7)$$

From here, we can immediately recognize some of the familiar from describing a quantum walk (where we are rescaling $\bar{p} \rightarrow \bar{p}/N_k$),

$$J_+^2 \rightarrow \frac{\partial^2}{\partial \bar{\varphi}^2}, \quad J_- \tilde{V}'(-iJ_+) \rightarrow -\frac{1}{N_k} \frac{\partial}{\partial \bar{p}} V'(\bar{\varphi}), \quad J_+ \left(-\frac{iJ_+}{2} + \frac{\partial}{\partial J_-} \right) \rightarrow -N_k \frac{\partial}{\partial \bar{\varphi}} \bar{p}. \quad (4.8)$$

In contrast, the appearance of gradient terms, such as $a^{-2}(\nabla\varphi)^2$, have typically not arisen in studies of stochastic inflation and corrections thereof. Intuitively, although one might imagine that gradients are small corrections, for the modes at the horizon ($p \sim aH$), this is not necessarily the case.² Instead, as explained using SdSET, we should understand these terms as being purely scheme dependent and removable by a field redefinition. Concretely, in dimensional regularization, power-law divergences vanish, and therefore we would expect these terms to be absent as well.

The resolution here follows the same pattern as irrelevant operators in any RG. In the presence of a hard cutoff, one might generate mixing under RG between irrelevant and marginal operators, but these are removable by a change of basis. For illustration, suppose we have a scalar field theory in flat space with

$$S = \int d^4x \left(\frac{1}{2} \partial_\mu \phi \partial^\mu \phi - \frac{\lambda_4}{4!} \phi^4 - \frac{\lambda_6}{6!} \frac{1}{\Lambda^2} \phi^6 \right). \quad (4.9)$$

With a hard cutoff $\epsilon\Lambda$, the one-loop contribution to λ_4 β -function is

$$\frac{\partial}{\partial \log \Lambda} \lambda_4 \supset \frac{\lambda_6}{4\pi^2} \epsilon^2. \quad (4.10)$$

Here we have introduced ϵ to show how it depends on the value of the hard cutoff, but ultimately we should take $\epsilon \rightarrow 1$ as Λ is defined to be our RG scale. The key to eliminating this term is to realize that λ_6 has a classical β -function so that

$$\begin{aligned} \frac{\partial}{\partial \log \Lambda} \lambda_6 = 2\lambda_6 \rightarrow \frac{\partial}{\partial \log \Lambda} \left(\lambda_4 - \frac{\lambda_6}{8\pi^2} \epsilon^2 \right) & \equiv \frac{\partial}{\partial \log \Lambda} \tilde{\lambda}_4 \\ & = 0. \end{aligned} \quad (4.11)$$

We see that by redefining the coupling as $\tilde{\lambda}_4$, we can eliminate the apparent contribution to RG from λ_6 . Of course, when we regulate with dimensional regularization, the power-law divergence vanishes, and there is no need to redefine the coupling.

The evolution of the density matrix under dynamical RG is in precise analogy with the scalar field example. Using $N_k \approx (aH)^3$ so that $\dot{N}_k = 3HN_k$, we notice that

$$\begin{aligned} \frac{1}{H} \frac{\partial}{\partial t} \left[\left(-a^2(t_0) H \frac{p(J_-)^2}{2N_k^2} - \frac{aH^2 J_- J_+}{p N_k} \right) + \left(K \frac{aH^2 J_+ J_-}{2p N_k} - \frac{a^2 H p}{2} K \frac{J_-^2}{N_k^2} \right) \right] \\ = \left[\left(4a^2(t_0) H \frac{p(J_-)^2}{2N_k^2} + 2 \frac{aH^2 J_- J_+}{p N_k} \right) + \left(-2K \frac{aH^2 J_+ J_-}{2p N_k} + 4 \frac{a^2 H p}{2} K \frac{J_-^2}{N_k^2} \right) \right]. \end{aligned} \quad (4.12)$$

²Our analysis is most similar to the treatment of [10], which similarly found additional gradient terms. However, they imposed the UV cutoff $\Lambda = \epsilon aH$ such that these terms were suppressed by powers of ϵ . This is sufficient to postpone the issue, but at the cost of eliminating superhorizon modes with $\epsilon aH < p < aH$.

In contrast, the additional terms in Eq. (4.6) do not explicitly depend on time. We can therefore remove the terms in Eq. (4.12) entirely by a polynomial redefinition of $Z[J_{\pm}]$.

Removing the irrelevant³ terms in the equation by this redefinition, we are left with the evolution equation for the Wigner distribution

$$\frac{d}{dt}W[\bar{\varphi}, \bar{p}] = \frac{H^3}{8\pi^2} \frac{\partial^2}{\partial \bar{\varphi}^2} W[\bar{\varphi}, \bar{p}] - \frac{\bar{p}}{a^3} \frac{\partial}{\partial \bar{\varphi}} W + a^3 \frac{\partial}{\partial \bar{p}} V'(\bar{\varphi}) W. \quad (4.13)$$

It is instructive to compare the final Wigner evolution equation to the density matrix evolution that we derived from QFT, and see which terms survived. We started with density matrix evolution in (3.45), and see that the ultimate fate is that for local measurements, the following terms drop out due to being irrelevant:

$$\begin{aligned} \frac{d}{dt}\rho_R = & \frac{H}{2} \int \frac{d^3 p}{(2\pi)^3} \left[\Delta_0 \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right)^2 - \frac{a^6}{K^2} \Delta_2 (\varphi_- - \varphi_+)^2 \right. \\ & + i \frac{a^3}{K} \Delta_R \left\{ (\varphi_+ - \varphi_-), \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) \right\} + 2 \frac{a^3}{K} \Delta_I \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) (\varphi_+ + \varphi_-) \Big] \rho_R \\ & + \int \frac{d^3 p}{(2\pi)^3} \left[\frac{iK}{2a^3} \left(\frac{\delta^2}{\delta\varphi_+^2} - \frac{\delta^2}{\delta\varphi_-^2} \right) \rho_R - \frac{ia^3}{2K} \left(\frac{(\nabla\varphi_+)^2}{a^2} - \frac{(\nabla\varphi_-)^2}{a^2} \right) \rho_R - ia^3 (V(\varphi_+) - V(\varphi_-)) \rho_R \right]. \end{aligned} \quad (4.14)$$

Out of the remaining terms, we see

$$\begin{aligned} & \int \frac{d^3 p}{(2\pi)^3} \left[\Delta_0 \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right)^2 + 2 \frac{a^3}{K} \Delta_I \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) (\varphi_+ + \varphi_-) \right] \rho_R \\ & \rightarrow \frac{H^3}{8\pi^2} \frac{\partial^2 W}{\partial \bar{\varphi}^2} \int \frac{d^3 p}{(2\pi)^3} \left[\frac{iK}{2a^3} \left(\frac{\delta^2}{\delta\varphi_+^2} - \frac{\delta^2}{\delta\varphi_-^2} \right) \rho_R \right] \\ & \rightarrow -q \frac{\partial W}{\partial \bar{\varphi}} - \int \frac{d^3 p}{(2\pi)^3} ia^3 (V(\varphi_+) - V(\varphi_-)) \rho_R \rightarrow a^3 \frac{\partial}{\partial \bar{p}} V'(\bar{\varphi}) W. \end{aligned} \quad (4.15)$$

It is useful to remember that Δ_I contains the boundary term that determines the real part of the wave function, as shown around Eq. (3.23). As a result, it is directly related to the Gaussian fluctuations of each mode and thus will be crucial in the derivation of the Fokker-Planck equation. See Appendix B for details of this derivation.

Now, we can transform variables to $\bar{p} = a^3(t)q$ to get

$$\begin{aligned} \frac{d}{dt}W[\bar{\varphi}, q] = & \frac{H^3}{8\pi^2} \frac{\partial^2 W}{\partial \bar{\varphi}^2} + 3H \frac{\partial}{\partial q} (qW) \\ & - q \frac{\partial}{\partial \bar{\varphi}} W + \frac{\partial}{\partial q} V'(\bar{\varphi}) W. \end{aligned} \quad (4.16)$$

Having set up the full phase space evolution equation, we can now compare this to the walk introduced in Sec. II. We see that this reproduces the same exact differential

³Here we mean irrelevant in the technical RG sense, as follows from SdSET.

equation with $D = H^3/8\pi^2$, $\gamma = 3H$, $m = 1$.⁴ As explained, at leading order, we can integrate out q by setting $q = -V'(x)/3H$. We then get the well-known stochastic inflation result

$$\frac{dP}{dt} = \frac{H^3}{8\pi^2} \frac{\partial^2 P}{\partial \bar{\varphi}^2} + \frac{\partial}{\partial \bar{\varphi}} \left(\frac{V'(\bar{\varphi})}{3H} P(\bar{\varphi}) \right). \quad (4.17)$$

There are, of course, easier ways to arrive at this result. What is meaningful here is that the determination of this partial differential equation from the exact RG as a functional over all the Fourier modes is more nuanced

⁴We see that stochastic inflation is represented as a noisy x and deterministic p evolution. This agrees with previous results for the evolution of the density matrix [24,30], but contradicts some generalizations of stochastic inflation [65–67]. Generalizations exist, where even p can be noisy due to coupling to additional degrees of freedom. These are studied in the open EFT of inflation. See for instance [34] and references therein.

than one might have expected from just comparing Eqs. (3.38) and (3.39).

B. Next-to-leading order

Our goal in this section is to determine the leading corrections to the stochastic formalism. However, unlike perturbation theory around the vacuum, the fluctuations around a random walk have a nontrivial definition of leading order, next-to-leading order and so on, that requires explanation.

Stochastic inflation is conventionally defined in terms of $P(\bar{\varphi}, t)$, the probability distribution for the long wavelength field. In terms of the Wigner distribution, we get this probability by “marginalizing” over q :

$$P(\bar{\varphi}, t) = \int dq W(\bar{\varphi}, q, t). \quad (4.18)$$

If we start with our LO description of stochastic inflation, equilibrium probability distribution can be determined by demanding $\dot{P} = 0$ and yields

$$P_{\text{eq}}(\bar{\varphi}) = \int dq W_{\text{eq}}(\bar{\varphi}, q) = \tilde{A} \exp\left(-\frac{8\pi^2 V(\bar{\varphi})}{3H^4}\right). \quad (4.19)$$

Let us take the specific case where $V(\bar{\varphi}) = \frac{\lambda}{4!} \bar{\varphi}^4$. The equilibrium probability distribution allows for order one probabilities for any $|\bar{\varphi}| < \lambda^{-1/4} H$. One might be concerned that there is no reliable expansion in $\lambda \ll 1$ given that $\bar{\varphi} \rightarrow \infty$ as $\lambda \rightarrow 0$. Fortunately, the corrections come with positive powers of λ that a meaningful expansion exists. At leading order, the time evolution of the probability distribution, Eq. (4.17), scales as $\frac{d}{dt}P = \mathcal{O}(\lambda^{1/2})$ when $\varphi \sim \lambda^{-1}$. At next-to-leading order, the correction to the time evolution gives contributions to $\dot{P}$ that are $\mathcal{O}(\lambda)$, and are therefore suppressed by $\sqrt{\lambda} \ll 1$ as desired.

Previous work has calculated the NLO corrections to the stochastic formalism and to the evolution of the density matrix. At this order, corrections to the Fokker-Planck equation can be defined as a nontrivial⁵ correction to the effective potential [10,14,52]

$$V_{\text{eff,NLO}}(\bar{\varphi}) = \frac{\lambda}{4!} \bar{\varphi}^4 + \frac{1}{6} \frac{1}{3!} \frac{\lambda^2}{18H^2} \bar{\varphi}^6 + \mathcal{O}(\lambda^3 \bar{\varphi}^8). \quad (4.20)$$

Notice that the effective potential becomes a series expansion in $\phi^2 \lambda$ and therefore breaks down when $\phi \sim 1/\sqrt{\lambda}$.

⁵Here we are avoiding calling these corrections universal because they are sensitive to the split between the time-independent and time-dependent components of φ . Specifically, the correction below is correct for the definition used in [14,52] but appears to contradict [10]. The difference arises from a redefinition of the decaying mode and can be accounted for by matching [52].

The effective potential corrections at NLO and beyond turn out to be determined by a field redefinition in SdSET. Concretely, eliminating the decaying mode, which is the conjugate momentum, introduces these terms. It was argued in [36] that the breakdown of series expansion of V_{eff} arises because the conjugate momentum to ϕ cannot be ignored at large field values.

The role of the momentum in generating the effective potential is seen most directly if we interpret the evolution of the Wigner distribution as a Langevin system. This is in exact analogy to the walk we analyzed in Sec. II C. In that case, $\dot{\phi} = q + \xi$ where ξ is our noise and

$$\dot{q} = 3Hq + V'(\bar{\varphi}) \rightarrow q(t) = \int^t dt' V'(\bar{\varphi}(t')) e^{3H(t-t')}. \quad (4.21)$$

Now expanding $\varphi(t') = \varphi(t) + (t' - t)\dot{\varphi}(t) + \dots$ we have

$$\begin{aligned} q(t) &= -\frac{V'(\bar{\varphi})}{3H} + \dot{\varphi} V''(\bar{\varphi}) \int dt (t' - t) e^{3H(t-t')} \\ &= -\frac{V'(\bar{\varphi})}{3H} - \frac{1}{27H^3} V'(\bar{\varphi}) V''(\bar{\varphi}). \end{aligned} \quad (4.22)$$

For $V(\bar{\varphi}) = \frac{\lambda \bar{\varphi}^4}{4!}$, if we interpret $3Hq = -V'_{\text{eff}}$, then we can see that

$$V_{\text{eff}} = \frac{\lambda}{4!} \bar{\varphi}^4 + \frac{1}{6} \frac{1}{3!} \frac{\lambda^2}{18H^2} \bar{\varphi}^6 + \dots, \quad (4.23)$$

in exact agreement with the direct calculations in QFT.

Of course, we should be able to derive a correction to the Fokker-Planck equation from q without the need for a Langevin description,

$$\begin{aligned} \frac{d}{dt} W[\bar{\varphi}, q] &= \frac{H^3}{8\pi^2} \frac{\partial^2 W}{\partial \bar{\varphi}^2} + 3H \frac{\partial}{\partial q} (qW) \\ &\quad - q \frac{\partial}{\partial \bar{\varphi}} W + \frac{\partial}{\partial q} V'(\bar{\varphi}) W. \end{aligned} \quad (4.24)$$

However, since the evolution of q is purely deterministic at this order, we can use the intuition from the Langevin description to define a change of variables

$$Q = a^3(t) \left(q + \frac{1}{3H} V'_{\text{eff}}(\bar{\varphi}) \right) \quad (4.25)$$

so that

$$\frac{\partial}{\partial t} \rightarrow \frac{\partial Q}{\partial t} \frac{\partial}{\partial Q} + \frac{\partial}{\partial t} = 3H Q \frac{\partial}{\partial Q} + \frac{\partial}{\partial t}, \quad (4.26)$$

$$\frac{\partial}{\partial \bar{\varphi}} \rightarrow \frac{\partial Q}{\partial \bar{\varphi}} \frac{\partial}{\partial Q} + \frac{\partial}{\partial \bar{\varphi}} = \frac{a^3}{3H} V''_{\text{eff}}(\bar{\varphi}) \frac{\partial}{\partial Q} + \frac{\partial}{\partial \bar{\varphi}}, \quad (4.27)$$

$$\frac{\partial}{\partial q} \rightarrow \frac{\partial Q}{\partial q} \frac{\partial}{\partial Q} = a^3 \frac{\partial}{\partial Q}. \quad (4.28)$$

Making this substitution, we find

$$\begin{aligned} \frac{d}{dt} W[\bar{\varphi}, q] = & 3HW + \frac{H^3}{8\pi^2} \frac{\partial^2 W}{\partial \bar{\varphi}^2} + a^3 V''_{\text{eff}} \frac{H^2}{12\pi^2} \frac{\partial^2 W}{\partial Q \partial \bar{\varphi}} + a^6 \frac{H}{72\pi^2} (V''_{\text{eff}})^2 \frac{\partial^2 W}{\partial Q^2} \\ & + a^3 \frac{\partial}{\partial Q} \left(-V'_{\text{eff}} + \frac{1}{9H^2} V'_{\text{eff}} V''_{\text{eff}} + \frac{H^2}{24\pi^2} V'''_{\text{eff}} + V' \right) W \\ & - V''_{\text{eff}} \frac{\partial}{\partial Q} \left(\frac{Q}{3H} W \right) - \frac{Q}{a^3} \frac{\partial W}{\partial \bar{\varphi}} + \frac{1}{3H} \frac{\partial}{\partial \bar{\varphi}} (V'_{\text{eff}} W). \end{aligned} \quad (4.29)$$

In order to have a normalized distribution in these new variables, we now rescale the Wigner distribution

$$\tilde{W}[\bar{\varphi}, Q] = \frac{\partial q}{\partial Q} W[\bar{\varphi}, q] = a^3(t) W[\bar{\varphi}, q]. \quad (4.30)$$

We can remove the mean for Q by choosing V_{eff} (neglecting V'''_{eff} at the order we want to consider) so that

$$\begin{aligned} V'_{\text{eff}} \left(1 - \frac{1}{9H^2} V''_{\text{eff}} \right) &= V' \rightarrow V'_{\text{eff}}(\bar{\varphi}) \\ &\approx V'(\bar{\varphi}) + \frac{1}{9H^2} V'(\bar{\varphi}) V''(\bar{\varphi}). \end{aligned} \quad (4.31)$$

The final step to determining the evolution of the probability of $\bar{\varphi}$ is to expand in moments [30], defined by

$$W_N(\bar{\varphi}) = \int dQ \left(\frac{Q}{a^3} \right)^N \tilde{W}(\bar{\varphi}, Q), \quad P(\bar{\varphi}) \equiv W_0(\bar{\varphi}). \quad (4.32)$$

This is effectively the same procedure used to convert the Boltzmann equations to the fluid equation, as is commonly done for describing the CMB.⁶ After performing our change of variables, this procedure is straightforward and gives rise to the NLO evolution equation

$$\begin{aligned} \frac{dP}{dt} = & \frac{H^3}{8\pi^2} \frac{\partial^2 P}{\partial \bar{\varphi}^2} + \frac{\partial}{\partial \bar{\varphi}} \left(\frac{V'_{\text{eff}}(\bar{\varphi})}{3H} P(\bar{\varphi}) \right) \\ & + \frac{1}{3H} \frac{\partial}{\partial \bar{\varphi}} \left(V''_{\text{eff}} \frac{H^2}{12\pi^2} \frac{\partial}{\partial \bar{\varphi}} P(\bar{\varphi}) \right). \end{aligned} \quad (4.33)$$

The final term arises from solving for the evolution of W_1 at NLO and plugging it into the equation for W_0 . As a check, the equation agrees with the equivalent equations derived in Sec. II C. The details of this calculation are shown in

⁶See Ref. [68], for example, for a systematic expansion of the corrections to the tight coupling approximation similar to our approach.

Appendix C. In the $\lambda\bar{\varphi}^4$ theory, the final term is removable by a field redefinition [52] of $\bar{\varphi}$ and thus is not universal.

The essential takeaway from this calculation is that all of the information relevant to both the LO and NLO calculation of stochastic inflation is present in the master equation for the Wigner distribution without including higher-order corrections of S_{eff} in our exact Wilsonian scheme.

C. NNLO and beyond

When calculating the corrections to stochastic inflation from QFT, the first nontrivial correction arises at NNLO, which is $\mathcal{O}(\lambda^{3/2})$ in the Fokker-Planck equation. Specifically, at this order we generate a non-Gaussian noise term

$$\frac{dP}{dt} = \frac{H^3}{8\pi^2} \frac{\partial^2 P}{\partial \bar{\varphi}^2} + \frac{\partial}{\partial \bar{\varphi}} \left(\frac{V'_{\text{eff}}(\bar{\varphi})}{3H} P(\bar{\varphi}) \right) + \frac{d_0}{3!} \frac{\partial^3}{\partial \bar{\varphi}^3} (\bar{\varphi} P(\bar{\varphi})), \quad (4.34)$$

where $d_0 = \frac{\lambda}{192\pi^2}$ was calculated in [52]. However, this result has not been obtained in a fully consistent scheme that removes all the subdivergences. See Ref. [69] for progress in this direction.

An essential difference between the approach in this paper and [52] is that we are working in a Wilsonian EFT, rather than a continuum EFT [70]. In the Wilsonian description, we integrated out the short modes explicitly, giving rise to higher dimension interactions, including higher derivative terms, in S_{eff} . These terms do not arise in a continuum theory, as they would vanish in dimensional regularization. Nevertheless, the corrections to these terms in the Wilsonian description are needed to reproduce higher order contributions to the RG flow [40]. In this regard, the challenge of calculating NNLO corrections in our exact RG scheme is not unique to de Sitter, but a common aspect of Wilsonian versus continuum descriptions.

In addition to these higher order terms from S_{eff} , we expect additional terms from eliminating Q systematically. These terms are also generally scheme dependent and will

bring higher powers of λ . There is a unique contribution at NNLO that is both straightforward to calculate and could contribute to d_0 , that arises from

$$\frac{d}{dt} \tilde{W}(\bar{\varphi}, Q) \supset -\frac{1}{24} V'''(\bar{\varphi}) \frac{\partial^3}{\partial Q^3} \tilde{W}. \quad (4.35)$$

Solving for the higher moments of $\tilde{W}$ and reducing the equation to the solution of $P = W_0$, one finds the NNLO term

$$\frac{d}{dt} P \supset \frac{1}{648 H^3} \frac{\partial^3}{\partial \bar{\varphi}^3} (V'''(\bar{\varphi}) P). \quad (4.36)$$

See Appendix C for the derivation. Notice that using $V = \lambda \bar{\varphi}^4/4!$, this is a contribution to d_0 defined above but is not consistent with the two-loop calculation in [52]. Moreover, at NNNLO, we expect a contribution from a three-loop diagram that would give a four derivative correction proportional to λ . There is no such term that can arise from eliminating Q .

The fact that the higher order corrections appear to require a more detailed analysis of the RG evolution is consistent with the physical expectation that higher-derivatives terms arise from non-Gaussian noise. Specifically, in a Markovian random walk, these terms arise from non-Gaussian cumulants in the Kramers-Moyal expansion of the transition amplitudes [71]. It would therefore be surprising to fully capture this effect, which we associate with the behavior of modes at horizon crossing, purely from the elimination of the fast variables in the superhorizon description.

Technically speaking, SdSET offers a much more straightforward path to calculating the coefficients at NNLO and beyond. However the long-term goal of this program is to calculate the tail of the distribution of $\bar{\varphi}$ beyond the limits of SdSET, when $\lambda \bar{\varphi}^2 > 1$. This will likely⁷ require a hybrid approach of SdSET and exact RG. The tails are sensitive to UV physics and therefore cannot be calculated purely within an effective description [36,38] (see also Refs. [73,74] for specific models). Yet, for a theory like $\lambda \phi^4$, where the theory is renormalizable, one might expect to be able to describe the tail up to the Landau pole. This is reminiscent of the effective potential in flat space, which can be improved using RG to resum the log corrections. There is reason to suspect a similar approach could apply for large deviations in de Sitter, but we will leave this to future work.

⁷An alternate proposal for calculating the tail of the distribution was made in [72] directly in terms of the wave function. As the statistics of local operators are described by a density matrix, an effective description in terms of a pure state can miss essential contributions to these observables [33].

V. RENORMALIZATION IN dS

Although light fields are where the need for normalization is most acute, our understanding of QFT in dS requires RG for all masses and spins, just like flat space. Direct calculations of the spectral densities are naturally interpreted in the usual language of conformal invariance and anomalous dimensions [75–78]. Much like the case of physics in AdS, it is natural to expect the language of RG is useful for interpreting these results.

One strong piece of evidence that RG flow can be used directly in dS comes from conformal field theory: if a given theory is conformal in flat space, then the same theory in dS will exhibit the same list operators and scaling dimensions. This can be seen easily using Weyl transformation. However, at nontrivial fixed point, these dimensions arise from mixing between operators and/or anomalous dimensions generated by interactions and therefore the same RG resummation needed to describe the CFT in flat space must apply in dS as well [51].

In flat slicing of dS, the Weyl transformation that takes flat space correlators to cosmological correlators involves scaling by the scale factor $a(\tau)$. For a CFT, this requires that any anomalous dimensions are associated with time-dependent logs. The required matching between dS and flat space therefore strongly suggests that the RG resummation involves time-dependent logs in dS. In this sense, it is dynamical RG that should be on the same footing as conventional RG in flat space.⁸

In this section, we will demonstrate that dynamical RG in dS obeys these properties. We will show that the exact RG argument of Sec. III can be applied to massive fields without any meaningful changes to the evolution equation. Repeating the argument of Polchinski [40] then demonstrates that all the RG corrections can be absorbed into the definitions of couplings and local operators of the theory.

A. Scalar fields of any mass

Our starting point will be an interacting scalar field $\varphi(\vec{x})$ with a mass m such that the action and density matrix are defined by Eq. (3.2), now with

$$\begin{aligned} S_0[\varphi(\vec{x}, t), \Lambda] &= \int_{\gamma} dt a^3(t) \int \frac{d^3 k}{(2\pi)^3} \frac{1}{2} K^{-1} \left(\frac{k^2}{\Lambda^2} \right) (\nabla_{\mu} \varphi(\vec{k}, t) \nabla^{\mu} \varphi(\vec{x}, t) \\ &\quad - m^2 |\varphi(\vec{k}, t)|^2). \end{aligned} \quad (5.1)$$

⁸For principle series fields, some anomalous dimensions can be calculated by resumming bubble diagrams, more like a decay width [79,80]. These examples are compatible with the RG description, as shown in [81,82]. Yet, the conformal example [51] demonstrates that not all anomalous dimensions can be calculated by summing geometric series.

We may be concerned only with local operators constructed from $\varphi(\vec{x})$ and derivatives, $\mathcal{O}_n(\vec{x})$. Since we are interested in local correlation function in the in-in formalism (rather than the density matrix, following [33]), we will define a partition function

$$Z[\{J_n\}; t] = \text{Tr}_\phi \left[\exp \left(\sum_n \int d^3x J_n \mathcal{O}_n(\vec{x}) \right) \rho(\phi, \phi') \right], \quad (5.2)$$

where $\mathcal{O}_n(\vec{x})$ are local operators composed from the UV field ϕ in the Schrödinger representation. Because of the trace, we know the result must be independent of Λ , in contrast to the density matrix itself. Our goal is to show that the process of integrating out the short distance physics and evolving in time combine to give a local theory where the local action and operators are modified.

At this stage, we can repeat the argument from Sec. III with only minor modifications. Nothing about the procedure assumed the fields were massless, other than the specific form of the action and Green's function. However, we note that if we make the substitutions

$$\begin{aligned} \nabla^2 &\rightarrow \nabla^2 - m^2, & G^{-1} &= K^{-1}(\nabla^2 - m^2), \\ G^{-1}G &= a^{-3}\delta(t-t')\delta(\vec{x}-\vec{x}'), \end{aligned} \quad (5.3)$$

then an exact RG result is unchanged. The essential reason is that the mass term, or any potential $V(\phi)$, does not introduce any boundary terms when integrating by parts. The boundary terms are again defined by Δ_0 , Δ_2 , and Δ_{LR} with the Green's functions replaced by the massive ones, for example

$$\begin{aligned} \Delta_0(p, t_0) &= iG' = iG'(t_0, t_0, \vec{p}) \\ &= -\frac{\partial K}{\partial s} \frac{\pi}{4} (-\tau_0)^3 H_\nu^{(1)}(-p\tau_0) H_\nu^{(2)}(-p\tau_0), \end{aligned} \quad (5.4)$$

where τ_0 is the conformal time at physical time t_0 .

B. Diagrammatic understanding of exact RG

Given the setup, we can now apply exact RG to the dS generating functional directly. Our goal is to demonstrate that the change to $Z[\{J_n\}; t]$ from integrating out $p \in [\Lambda, \Lambda + \delta\Lambda]$ and shifting in time by $\delta t = H^{-1}\delta\Lambda/\Lambda$ can be completely absorbed into the definition of (a) the effective action S_{eff} and (b) the definition of the local operators $\mathcal{O}_n(\vec{x})$. This will ensure that our RG equations are differential equations with no explicit dependence on the UV physics and therefore can be integrated to determine correlation functions, just like the Callan-Symanzik equation for RG in flat space.

The first step is straightforward: we have already shown that integrating out a shell of momenta gives rise to the following change to the effective action, illustrated in Fig. 1,

$$\begin{aligned} \frac{\partial}{\partial s} S_{\text{int}} &= \int \frac{d^3p}{(2\pi)^3} \int_\gamma dt dt' \left(\frac{1}{2} \frac{\delta S_{\text{int}}}{\delta \varphi(t)} G(t, t') \frac{\delta S_{\text{int}}}{\delta \varphi(t')} \right. \\ &\quad \left. - \frac{1}{2} \left(\frac{\delta^2 S_{\text{int}}}{\delta \varphi(t) \delta \varphi(t')} G(t, t') \right) \right). \end{aligned} \quad (5.5)$$

The second term is manifestly local because

$$\frac{\delta^2 S_{\text{int}}}{\delta \varphi(t) \delta \varphi(t')} = \delta(t-t') \frac{\partial^2}{\partial^2 \varphi} \mathcal{L}_{\text{int}}(\varphi, t). \quad (5.6)$$

The first term, on the other hand, is bilocal in time and therefore technically nonlocal. However, with the $i\epsilon$ prescription, we know that

$$G(t, t') \propto e^{-\epsilon p|t-t'|}, \quad (5.7)$$

for t, t' on the same side of the Schwinger-Keldysh contour (i.e. $t, t' \in [-\infty(1-i\epsilon), t_0]$ or $t, t' \in [-\infty(1+i\epsilon), t_0]$). This ensures we can Taylor expand in $t-t'$ and integrate each term to find a finite result to all orders. In this sense, as far as the effective action is concerned, there is nothing about integrating out a shell of momentum in dS that is distinct from flat space.

The situation is similar from the contributions to the action from $t \in [-\infty(1-i\epsilon), t_0]$ and $t' \in [-\infty(1+i\epsilon), t_0]$ (or vice versa). In these cases, the Green's functions are again exponentially suppressed but now by

$$G(t = t_+, t' = t_-) \propto e^{-\epsilon p|t_+ - t_0| - \epsilon p|t_- - t_0|}. \quad (5.8)$$

These terms should therefore be expanded around $t_\pm = t_0$, giving rise to a local expansion in operators at the boundary. These terms are responsible for renormalizing the definition of local operators, or equivalently, the sources J_n .

The nontrivial changes that arise for in-in correlators in de Sitter space are the boundary terms at t_0 and the fact that $\Lambda(t)$ is combined with time evolution in the full RG transformation. Yet, it is straightforward to understand the impact of these terms. We writing our generating functional as

$$\begin{aligned} \frac{d}{dt} Z[\{J_n\}; t] &= \text{Tr}_\varphi \left[\exp \left(\sum_n \int d^3x J_n \mathcal{O}_n(\vec{x}) \right) \frac{d}{dt} \rho_R(\varphi, \varphi') \right]. \end{aligned} \quad (5.9)$$

Of course, this is identical to the derivation of the evolution in Eq. (3.54), so that schematically we have

$$\begin{aligned} \frac{d}{dt} Z[J_\pm, t] &= \text{Tr} \left[\exp \left(\sum_n \int d^3x J_n \mathcal{O}_n(\vec{x}) \right) \right. \\ &\quad \left. \times \left(-i[H_{\text{eff}}, \rho] + \sum_i L_i \rho L_i^\dagger - \frac{1}{2} \{L_i^\dagger L_i, \rho\} \right) \right]. \end{aligned} \quad (5.10)$$

Using the trace, we can exchange operators acting on ρ to acting on the local operator insertions. However, because these act locally on $\mathcal{O}_n(\vec{x})$ we can express time evolution as a local change to the operator (which would be the Heisenberg picture, in the absence of the L_i). Therefore, we can treat this apparent time evolution as the operator equation for

$$\frac{d}{dt}V[\{J_n\}, t] = i[H_{\text{eff}}, V[\{J_n\}, t]] + \sum_i L_i^\dagger V[\{J_n\}, t] L_i - \frac{1}{2}\{L_i^\dagger L_i, V[\{J_n\}, t]\}, \quad (5.11)$$

$$V[\{J_n\}, t] = \exp\left(\sum_n \int d^3x J_n \mathcal{O}_n(\vec{x})\right). \quad (5.12)$$

If $\mathcal{O}_n$ forms a complete basis of local operators, then the change to $V[\{J_n\}, t]$ is just a redefinition of J_n . This expression is purely local in time and therefore we can simply redefine $J_n \rightarrow J'_n$ to capture the time evolution of the system.

One might worry that the fact that $J_n(\vec{x})$ is a source that depends directly on position could introduce nonlocality in space. This does not occur because the nontrivial evolution of the operators arises from equal time commutators, and therefore are local in space by causality. This is manifest in the UV Hamiltonian description. In terms of the Wilsonian RG, the evolution of the operators arises from contribution that couple the two branches of the Schwinger-Kelgysch contour (t_+ and t_-) and are not equivalent to a local term in the action because of the path ordering of the exponential. Of course, the time ordering is only relevant for the terms that do not commute, which are again local in space at equal time. The same features appear in flat space where several examples have been worked out in detail [33].

The essential result here again follows from Eq. (3.5). The sources can be added to the boundary action

$$S_{\partial, \text{tot}} = S_{\partial} + \int d^3x \sum_n J_n \mathcal{O}_n(\vec{x}). \quad (5.13)$$

In these variables, our partition function is simply

$$Z[\{J_n\}; t] = \int_{\gamma} D\varphi T \exp(iS_0 + iS_{\text{int}} + iS_{\partial}), \quad (5.14)$$

and therefore obeys that same renormalization properties as the effective action. At this stage, the properties of the renormalization of de Sitter correlators and the flat space in-in correlators are identical, up to the definition of Green's functions and the fact that the de Sitter RG transformation includes a step of time evolution.

It is, perhaps, worth noting that the analogous proof of “renormalizability” in AdS remains incomplete [83]. In that case, we have the benefit of AdS/CFT which implies that the dual description is a conventional QFT and therefore the bulk of AdS should inherit what is known renormalization. Yet, some of the peculiar aspects of dynamical RG in

dS [82] have nearly identical avatars in AdS [84,85]. A proof in AdS would likely be instructive both for understanding holographic RG in AdS [86], and as another connection between physics in AdS and dS.

VI. CONCLUSIONS

Understanding the dynamics of light scalar fields on a fixed de Sitter background presents a surprising challenge. Perturbative calculations fail due to large time-dependent logs [4,87]. These logs have been understood and resummed using the ideas from classical random walks but leave the quantum nature of the fields obscure. Moreover, while understanding light scalars should be a stepping stone to quantum cosmology, the stochastic formalism bears little resemblance to the Euclidean path integral or wave function of the Universe [88]. Yet, both approaches yield the same equilibrium probability distribution [89], which demands a deeper explanation. A comparison of the density matrices of the no-boundary proposal [90–93] with the stochastic formalism offers the opportunity to make a deeper connection between these descriptions.

Understanding the de Sitter entropy [94], entanglement of inflationary modes, [60,95,96], and cosmology as an open quantum system [8,34] motivates the need to describe the evolution of the density matrix from a stochastic perspective. In this paper, we developed the tools of exact RG to provide a nonperturbative evolution equation for the long-wavelength density matrix. While elements of the stochastic description arise from tracing out subhorizon fluctuations, we also find that the description as a quantum walk only emerges when additionally tracing over information at separated points. This formalism clarifies and generalizes various aspects of stochastic inflation and should serve a powerful tool for exploring nonperturbative aspects of QFT in dS.

The density matrix in dS also played a significant role in prior work [22–26], but has largely been used for exploring the decoherence of fluctuations in cosmological backgrounds [97–103]. It is a practical fact that cosmological correlators are effectively classical, but the manner in which classicality emerges remains an area of active research. A potentially audacious theoretical and observational goal is to determine empirically that structure in the Universe arises from quantum fluctuations [104–108]. Such a determination would necessitate techniques beyond the Bell inequalities [105,109] and instead would have to isolate the quantum evolution during the inflationary epoch. A more precise understanding of the evolution of the density matrix in cosmology may serve as a setting stone to these ambitions as well.

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DATA AVAILABILITY

There are no publicly available research data or software supporting this manuscript. Requests for further information or data should be sent to the authors.

APPENDIX A: RELATIONSHIP BETWEEN GENERATING FUNCTIONALS AND PROBABILITIES

In the main text, we chose to derive the density matrix of the local long distance field φ by defining the generating functional

$$Z[J_{\pm}, t] = \text{Tr}[e^{J_+ \bar{\varphi}} e^{J_- \bar{\rho}} \rho_{\text{red}}(\varphi_+(\vec{x}), \varphi_-(\vec{x}))]. \quad (\text{A1})$$

While working with the generating function is a common framework, it does make the connection to previous work less transparent. The purpose of this section is to show this representation is identical to the previous approaches. To make the discussion easy to follow, we will work with a collection of harmonic oscillator wave functions $\psi_i(x)$ but we want to calculate a generating functional for $\bar{x} = \sum_k \hat{x}_k$ and $\bar{p} = N^{-1} \sum_k \hat{p}_k$, namely

$$Z[J_{\pm}] = \text{Tr}_k \left[e^{J_+ \bar{x}} e^{J_- \bar{p}} \prod_k |\psi_k\rangle \langle \psi_k| \right] = \text{Tr}[e^{J_+ \bar{x}} e^{J_- \bar{p}} \rho], \quad (\text{A2})$$

where $\rho = \prod_k \rho_k$ is the total density matrix for all k -harmonic oscillators.

First, we would like to compare this with our random walks discussed in Sec. II, where we defined a density matrix for a degree of freedom $\hat{X}$ via the interaction

$$\rho(X, X') = \text{Tr}_k [e^{i\hat{P}_X \bar{x}} |X=0\rangle \langle X'=0| e^{-i\hat{P}_{X'} \bar{x}} \otimes \rho], \quad (\text{A3})$$

where we took the trace in the $\hat{x}_k$ basis. Now we notice that if we work in the momentum basis $X \rightarrow P_X$, $|X=0\rangle \rightarrow$ constant and therefore

$$\begin{aligned} \rho(P_+, P_-) &= \text{Tr}_k [e^{iP_+ \bar{x}} \rho e^{-iP_- \bar{x}}] \\ &= Z[J_+ = i(P_+ - P_-), J_- = 0], \end{aligned} \quad (\text{A4})$$

where $P_{\pm}$ are just the momentum eigenvalues. We see that in the momentum basis, the density matrix is just the generating function with imaginary values of J . In fact, this also represents the measurements of a local observer along a worldline as in [19].

Another approach is to calculate the density matrix while imposing the constraint $\sum x_i = \bar{x}$ in the form of a δ -function,

$$\begin{aligned} &\rho(\bar{x} + \Delta x, \bar{x} - \Delta x) \\ &= \text{Tr}_k \left[\delta\left(\bar{x} + \Delta x - \sum_k x_k\right) \delta\left(\bar{x} - \Delta x - \sum_k x'_k\right) \right. \\ &\quad \left. \times \rho(\{x_i\}, \{x'_i\}) \right]. \end{aligned} \quad (\text{A5})$$

This was employed in [10] for $\Delta x = 0$ in order to determine the probability distribution in stochastic inflation for $\bar{\varphi} \rightarrow \bar{x}$. In order to derive such a result from the generating function, we first note that if we take $J_- = i2\Delta x$, then

$$Z[J_+, i\Delta x] = \text{Tr}[e^{J_+ \bar{x}} \rho(\{x + \Delta x/N\}, \{x' - \Delta x/N\})]. \quad (\text{A6})$$

Now defining $J_+ = ik$ and taking the Fourier transform we get

$$\int \frac{dk}{(2\pi)} e^{-ik\bar{x}} Z[J_+ = ik, J_- = i\Delta x] \quad (\text{A7})$$

$$= \text{Tr}_k \left[\delta\left(\bar{x} - \sum_k x_k\right) \rho(\{x_i + \Delta x/N\}, \{x_i - \Delta x/N\}) \right]. \quad (\text{A8})$$

Applying this formula back to the de Sitter context, if we take $\Delta x = 0$, $\bar{x} \rightarrow \bar{\varphi}$, and $x_k \rightarrow \varphi(\vec{k})$ we get

$$P(\bar{\varphi}) = \int D\varphi \delta\left(\bar{\varphi} - \int \frac{d^3 k}{(2\pi)^3} \varphi(\vec{k})\right) P[\varphi(\vec{k})] \quad (\text{A9})$$

in agreement with [10]. Finally, from this expression, we can derive the Wigner quasiprobability distribution for $\bar{x}$ and $\bar{x}'$, starting from

$$W[\bar{x}, \bar{p}] = \int d\Delta x e^{-i\bar{p}\Delta x} \rho\left(\bar{x} + \frac{\Delta x}{2}, \bar{x} - \frac{\Delta x}{2}\right). \quad (\text{A10})$$

We can transform $Z[J_{\pm}]$ as

$$\begin{aligned} \text{Tr}_k [e^{J_+ \bar{x}} e^{J_- \bar{p}} \rho] &= \text{Tr}_k [e^{\frac{iJ_+ J_-}{2}} e^{J_+ \bar{x}} e^{\frac{J_- \bar{p}}{2}} \rho e^{\frac{J_- \bar{p}}{2}}] \\ &= \text{Tr}_k \left[e^{\frac{iJ_+ J_-}{2}} e^{J_+ \bar{x}} \rho\left(x - \frac{iJ_-}{2N}, x' + \frac{iJ_-}{2N}\right) \right]. \end{aligned} \quad (\text{A11})$$

Then, taking $J_+ = ik$, $J_- = iy$, we get $W[\bar{x}, \bar{p}]$ via

$$W[\bar{x}, \bar{p}] = \int dy e^{-i\bar{p}y} \int \frac{dk}{(2\pi)} e^{i\frac{ky}{2}} e^{-ik\bar{x}} Z[J_+ = ik, J_- = iy]. \quad (\text{A12})$$

This, of course, is not limited to the Wigner distribution alone; any Fourier transform of the Wigner distribution can be derived in the same way.

APPENDIX B: EVOLUTION EQUATION FOR THE GENERATING FUNCTIONAL

In this section, we want to simplify the evolution equation for the generating functional. We start with the definition

$$Z[J_{\pm}, t] = \text{Tr}[e^{J_+ \bar{\varphi}} e^{J_- \bar{\pi}} \rho_{\text{red}}(\varphi_+(\vec{x}), \varphi_-(\vec{x}))], \quad (\text{B1})$$

and hence we have

$$\frac{d}{dt} Z[J_{\pm}, t] = \text{Tr}[e^{J_+ \bar{\varphi}} e^{J_- \bar{\pi}} \frac{d}{dt} \rho_{\text{red}}(\varphi_+(\vec{x}), \varphi_-(\vec{x}))]. \quad (\text{B2})$$

At this stage, we can substitute the evolution equation for ρ_{red} using (3.45). Defining $\mathcal{O} = e^{J_+ \bar{\varphi}} e^{J_- \bar{\pi}}$, we have that

$$\mathcal{O}(\varphi_-(p), \varphi_+(p)) = \langle \varphi_- | e^{J_+ \bar{\varphi}} e^{J_- \bar{\pi}} | \varphi_+ \rangle = e^{J_+ \bar{\varphi}} \delta \left[\varphi_+(p) - \varphi_-(p) + \frac{iJ_- K(p^2/\Lambda^2)}{N_k} \right]. \quad (\text{B3})$$

Then, we equivalently have

$$\begin{aligned} Z[J_{\pm}, t] &= \int D\varphi_+ D\varphi_- \mathcal{O}(\varphi_-, \varphi_+) \rho(\varphi_+, \varphi_-) \\ &= \int D\varphi_+ D\varphi_- e^{J_+ \bar{\varphi}} \delta \left[\varphi_+(p) - \varphi_-(p) + \frac{iJ_- K(p^2/\Lambda^2)}{N_k} \right] \rho(\varphi_+, \varphi_-). \end{aligned} \quad (\text{B4})$$

Now, we can plug in our time derivative of ρ to get

$$\begin{aligned} \frac{dZ}{dt} &= \int D\varphi_+ D\varphi_- \mathcal{O} \int \frac{d^3 p}{(2\pi)^3} \left[\frac{H\Delta_0}{2} \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right)^2 - a^6(t_0) K^{-2} \frac{H\Delta_2}{2} (\varphi_- - \varphi_+)^2 \right. \\ &\quad \left. + ia^3(t_0) K^{-1} \frac{H\Delta_R}{2} \left\{ (\varphi_+ - \varphi_-), \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) \right\} + 2a^3 K^{-1} H\Delta_I \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) (\varphi_+ + \varphi_-) \right] \\ &\quad \left. + \frac{iK}{2a^3} \left(\frac{\delta^2}{\delta\varphi_+^2} - \frac{\delta^2}{\delta\varphi_-^2} \right) \rho - \frac{iK^{-1}a^3}{2} \left(\frac{(\nabla\varphi_+)^2}{a^2} - \frac{(\nabla\varphi_-)^2}{a^2} \right) \rho - ia^3 (V(\varphi_+) - V(\varphi_-)) \rho. \end{aligned} \quad (\text{B5})$$

We can integrate by parts until we simplify each term. All the derivatives will ultimately be put onto $\mathcal{O}$. The rules simplify to

$$\mathcal{O} \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) \rho \rightarrow -J_+ \mathcal{O} \rho \quad \mathcal{O}(\varphi_+ - \varphi_-) \rho \rightarrow -\frac{iJ_- K}{N_k} \mathcal{O} \rho \quad \mathcal{O} \left(\frac{\delta}{\delta\varphi_+} - \frac{\delta}{\delta\varphi_-} \right) \rho \rightarrow \left(2 \frac{N_k}{iK} \frac{\partial}{\partial J_-} \right) \mathcal{O} \rho. \quad (\text{B6})$$

We will also need to make a further simplification. In order to deal with the $\varphi_+ + \varphi_-$ terms, we will resort to perturbation theory and use the following leading order identity (for the massless case) as explained in Appendix B 1,

$$\begin{aligned} (\varphi_+ + \varphi_-) \rho &= -K \frac{H^2}{k^3} \left(\frac{\delta}{\delta\varphi_+} + \frac{\delta}{\delta\varphi_-} \right) \rho + \frac{i}{k\eta} (\varphi_+ - \varphi_-) \rho \\ \mathcal{O}(\varphi_+ + \varphi_-) \rho &\rightarrow K \frac{H^2}{k^3} J_+ \mathcal{O} \rho - aH \frac{J_- K}{N_k} \mathcal{O} \rho. \end{aligned} \quad (\text{B7})$$

Using these identities, we can completely simplify the above equation. We get

$$\begin{aligned} \frac{dZ}{dt} &= \frac{1}{2} \int \frac{d^3 p}{(2\pi)^3} \left[H\Delta_0 J_+^2 + a^6 H\Delta_2 \left(\frac{J_-}{N_k} \right)^2 - 2a^3 H\Delta_R \left(\frac{J_-}{N_k} J_+ \right) \right. \\ &\quad \left. - 2a^3 \Delta_I \frac{H^3}{p^3} J_+^2 + 2a^4 \frac{H^2}{p} \Delta_I \left(\frac{J_+ J_-}{N_k} \right) \right] K' Z + \frac{N_k}{a^3} J_+ \left(-\frac{iJ_+}{2} + \frac{\partial}{\partial J_-} \right) Z \\ &\quad - ia^3 \left(\tilde{V} \left(-iJ_+ - i\frac{J_-}{2N_k} \right) - \tilde{V} \left(-iJ_+ + i\frac{J_-}{2N_k} \right) \right) Z + \int \frac{d^3 p}{(2\pi)^3} \left(K \frac{aH^2}{2p} \frac{J_+ J_-}{N_k} - \frac{a^2 H p}{2} K \frac{J_-^2}{N_k^2} \right) Z. \end{aligned} \quad (\text{B8})$$

The terms with the p integral correspond to the shell of momentum just integrated in, while the local terms correspond to Hamiltonian time evolution. After using the form of the Δ for the massless theory, we get

$$\begin{aligned}
\frac{dZ[J_{\pm}]}{dt} = & \frac{1}{2} \int \frac{d^3 p}{(2\pi)^3} \left[\left(\frac{H^3}{2p^3} J_+^2 - a^2(t_0) H \frac{p}{2} \frac{(J_-)^2}{N_k^2} - \frac{aH^2}{p} \frac{J_- J_+}{N_k} \right) \frac{\partial K}{\partial s} \right] Z[J_{\pm}] \\
& + \frac{N_k}{a^3} J_+ \left(-\frac{iJ_+}{2} + \frac{\partial}{\partial J_-} \right) Z[J_{\pm}] - \frac{J_-}{N_k} a^3 \tilde{V}'(-iJ_+) Z[J_{\pm}] \\
& + \int \frac{d^3 p}{(2\pi)^3} \left(K \frac{aH^2}{2p} \frac{J_+ J_-}{N_k} - \frac{a^2 H p}{2} K \frac{J_-^2}{N_k^2} \right) Z[J_{\pm}].
\end{aligned} \tag{B9}$$

1. Identity useful for homogenization

The homogenization of terms of the form

$$\int \frac{d^3 p}{(2\pi)^3} \mathcal{O}(\varphi_-, \varphi_+) (\varphi_+ + \varphi_-) (p) \rho_R \tag{B10}$$

need to be dealt with perturbatively. The basic idea is that for harmonic oscillators, this represents a term of the form $x\psi \propto p\psi$ when ψ is a ground state. We can use a similar idea here, starting with the UV mode ϕ , that is defined for the free theory as

$$\phi(k) = \frac{H}{\sqrt{2k^3}} [(1 - ik\eta) e^{ik\eta} a_k + (1 + ik\eta) e^{-ik\eta} a_k^\dagger], \quad \Pi(k) = K^{-1} \frac{\sqrt{k}}{\sqrt{2H\eta}} [e^{ik\eta} a_k + e^{-ik\eta} a_k^\dagger].$$

Solving for a_k in the limit $k\eta \ll 1$, we get

$$\begin{aligned}
a_k &= i[\phi(k, \eta) v_k^* - K\Pi(k, \eta) u_k^*], \quad v_k = a^2 u_k', \quad u_k = \frac{H}{\sqrt{2k^3}} (1 + ik\eta) e^{-ik\eta}, \\
&= i \left[\frac{k^2}{H\sqrt{2k^3}\eta} (1 + ik\eta) \phi(k, \eta) - \frac{H}{\sqrt{2k^3}} K\Pi(k, \eta) \right] \\
&= \frac{i}{\sqrt{2k^3}} \left[\left(\frac{k^2}{H\eta} + \frac{ik^3}{H} \right) \phi - HK\Pi \right].
\end{aligned} \tag{B11}$$

The lowering operator annihilates the vacuum, giving

$$a_k \Psi_{\text{BD}} = 0 \Rightarrow \left(\frac{k^2}{H\eta} + \frac{ik^3}{H} \right) \phi(k) \Psi_{\text{BD}} = -iHK \frac{\delta}{\delta\phi(k)} \Psi_{\text{BD}},$$

as can be checked directly using

$$\Psi_{\text{BD}}[\phi, \eta] = \prod_{\vec{k}} \left(\frac{2k^3}{\pi} \right)^{1/4} \exp \left[\frac{iK^{-1}}{2H^2} \frac{k^2}{\eta(1 - ik\eta)} \phi(\vec{k}) \phi(-\vec{k}) \right] \frac{e^{-ik\eta/2}}{\sqrt{1 - ik\eta}}.$$

Thus, the density matrix satisfies

$$H^2 K \left(\frac{\delta}{\delta\phi_+} + \frac{\delta}{\delta\phi_-} \right) \rho = \left(\frac{ik^2}{\eta} (\phi_+ - \phi_-) - k^3 (\phi_+ + \phi_-) \right) \rho. \tag{B12}$$

This is a useful simplification. We will write this as

$$(\phi_+ + \phi_-) \rho = -\frac{H^2}{k^3} K \left(\frac{\delta}{\delta\phi_+} + \frac{\delta}{\delta\phi_-} \right) \rho + \frac{i}{k\eta} (\phi_+ - \phi_-) \rho. \tag{B13}$$

This is, of course, only true at the Gaussian level, which is sufficient for our purposes right now. Now, we can evaluate terms of the form

$$\int \frac{d^3 p}{(2\pi)^3} \mathcal{O}(\varphi_-, \varphi_+) (\varphi_+ + \varphi_-) (p) \rho_R \rightarrow \int \frac{d^3 p}{(2\pi)^3} \mathcal{O} \left(-\frac{H^2}{p^3} K \left(\frac{\delta}{\delta\phi_+} + \frac{\delta}{\delta\phi_-} \right) \rho + \frac{i}{p\eta} (\varphi_+ - \varphi_-) \rho_R \right) \tag{B14}$$

using the rules in (B6).

APPENDIX C: ELIMINATING MOMENTUM WITH MOMENTS

The type of random walk under consideration is generally referred to as a “fast-slow” system in the random walk literature. The fast variables have evolution with a time scale γ^{-1} where the slow system evolves on a timescale set by the potential V' and the amplitude of the noise, A . One then aims to systematically eliminate the fast variables, leaving a Markovian evolution equation for the slow variables as an expansion in γ^{-1} . There is a formal procedure [71] where the solutions are projected onto the solutions of the fast equations, giving a systematic

expansion in γ^{-1} . However, as we are only interested in calculating the NLO (and potentially NNLO) terms, this procedure is both overkill and somewhat ill-suited for the problem at hand.

The most straightforward approach to eliminating Q from the evolution of $W(\bar{\varphi}, Q)$, following [30], is to define the moments of the distribution

$$W_N(\bar{\varphi}) = \int dQ (a^{-3} Q)^N W(\bar{\varphi}, Q), \quad P() \equiv W_0(). \quad (C1)$$

We have changed variables such that the evolution equation is schematically of the form

$$\begin{aligned} \frac{d}{dt} W[\bar{\varphi}, Q] = & \frac{H^3}{8\pi^2} \frac{\partial^2 W}{\partial \varphi^2} + a^3 V''_{\text{eff}}(\bar{\varphi}) \frac{H^2}{12\pi^2} \frac{\partial^2 W}{\partial Q \partial \bar{\varphi}} + a^6 \frac{H}{72\pi^2} (V''_{\text{eff}}(\bar{\varphi}))^2 \frac{\partial^2 W}{\partial Q^2} - V''_{\text{eff}}(\bar{\varphi}) \frac{\partial}{\partial Q} \left(\frac{Q}{3H} W \right) \\ & - \frac{Q}{a^3} \frac{\partial W}{\partial \bar{\varphi}} + \frac{1}{3H} \frac{\partial}{\partial \bar{\varphi}} (V'_{\text{eff}}(\bar{\varphi}) W), \end{aligned} \quad (C2)$$

In terms of the moments, this equation becomes

$$\left(\frac{d}{dt} + 3HN \right) W_N[\bar{\varphi}] = \hat{L}_1 W_N - \frac{\partial W_{N+1}}{\partial \bar{\varphi}} + V''_{\text{eff}} \left(-N \frac{H^2}{12\pi^2} \frac{\partial}{\partial \bar{\varphi}} W_{N-1} + \frac{H}{72\pi^2} V''_{\text{eff}} N(N-1) \partial W_{N-2} \right), \quad (C3)$$

where $\hat{L}_1 W = \frac{\partial}{\partial \bar{\varphi}} \left(\frac{1}{3H} V'_{\text{eff}}(\bar{\varphi}) \right) W + \frac{H^3}{8\pi^2} \frac{\partial^2 W}{\partial \bar{\varphi}^2} + N \frac{V''_{\text{eff}}}{3H} W$. Therefore, we have

$$\frac{d}{dt} W_0 = \hat{L}_1 W_0 - \frac{\partial}{\partial \bar{\varphi}} W_1 \quad (C4)$$

$$\frac{d}{dt} W_1 + 3H W_1 = L_1 W_1 - V''_{\text{eff}} \frac{H^2}{12\pi^2} \frac{\partial}{\partial \bar{\varphi}} W_0 + \dots \quad (C5)$$

Here the ... are higher order in λ and can be dropped for the NLO terms. At this order, it is easy to solve for W_1 treating $3H = \gamma$ as the large parameter, so that

$$\frac{d}{dt} W_0 = \hat{L}_1 W_0 + \frac{1}{3H} \frac{\partial}{\partial \bar{\varphi}} V''_{\text{eff}} \frac{H^2}{12\pi^2} \frac{\partial}{\partial \bar{\varphi}} W_0. \quad (C6)$$

We see therefore both the correction to the effective potential and the noise at NLO.

At higher orders, such as NNLO, the natural question is whether it is eliminating the higher orders in W_N or nonlinear contributions to S_{eff} that will play the dominant role in determining the evolution equations. We will be specifically interested in whether the following NNLO term can explain the results of [52],

$$\frac{d}{dt} \tilde{W}(\bar{\varphi}, Q) \supset -\frac{1}{24} V'''(\varphi) a^9 \frac{\partial^3}{\partial Q^3} W. \quad (C7)$$

This gives rise to a new contribution to W_3 ,

$$\left(\frac{d}{dt} + 9H \right) W_3(\bar{\varphi}) \supset \frac{1}{4} V'''(\bar{\varphi}) W_0(\bar{\varphi}). \quad (C8)$$

Using Eq. (C3), we can solve for leading order in H^{-1} terms to solve for the contributions to W_2 and then W_1 , to find

$$W_2 \approx -\frac{1}{6H} \frac{\partial}{\partial \bar{\varphi}} W_3 = -\frac{1}{216H^2} \frac{\partial}{\partial \bar{\varphi}} V'''(\bar{\varphi}) W_0(\bar{\varphi}), \quad (C9)$$

$$W_1 \supset -\frac{1}{3H} \frac{\partial}{\partial \bar{\varphi}} W_2 = \frac{1}{648H^3} \frac{\partial^2}{\partial \bar{\varphi}^2} V'''(\bar{\varphi}) W_0(\bar{\varphi}), \quad (C10)$$

and therefore our correction to the evolution of $W_0 \equiv P$ becomes

$$\frac{d}{dt} P \supset -\frac{1}{648H^3} \frac{\partial^3}{\partial \bar{\varphi}^3} (V'''(\bar{\varphi}) P). \quad (C11)$$

This is of the correct form to describe our non-Gaussian noise term at NNLO but the coefficient does not appear to match the result from SdSET [52]. Here we have neglected many other NNLO terms, which are generally not universal, along with any corrections to S_{eff} in the interacting theory. This is therefore an incomplete calculation and shows the need for more efficient techniques for calculating at NNLO and beyond.

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