Quantization of the ModMax oscillator

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We quantize the ModMax oscillator, which is the dimensional reduction of the modified Maxwell theory to one spacetime dimension. We show that the propagator of the ModMax oscillator satisfies a differential equation related to the Laplace equation in cylindrical coordinates, and we obtain expressions for the classical and quantum partition functions of the theory. To do this, we develop general results for deformations of quantum-mechanical theories by functions of conserved charges. We show that canonical quantization and path integral quantization of such deformed theories are equivalent only if one uses the phase space path integral; this gives a precise quantum analog of the statement that classical deformations of the Lagrangian are equivalent to those of the Hamiltonian.

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I. INTRODUCTION

Historically, quantum field theories first arose via the quantization of classical field theories. Following the modern usage of the term [1], we understand a quantum field theory (QFT) to mean any model that is compatible with certain physical principles including quantum mechanics, locality, and Lorentz invariance on a fixed (d + 1)-dimensional spacetime manifold. When d = 0, the Lorentz structure becomes essentially trivial and one has an ordinary theory of quantum mechanics.

However, we still lack a systematic understanding of the process of quantization for several reasons. One reason is that it is not known how to uniquely quantize a general classical theory, except in the case of theories which can be brought into a conventional form with a quadratic kinetic term. A famous example is the Nambu-Goto action of string theory; rather than attempting to quantize this theory directly, one first rewrites it in the classically equivalent form of the Polyakov action, which can then be quantized because the theory is quadratic in derivatives. A second reason is that not all quantum field theories admit classical limits. This means that certain QFTs cannot ever be

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understood by quantization of a classical theory, defined. for instance. by a Lagrangian or Hamiltonian (indeed, many QFTs are non-Lagrangian and thus do not even admit such a description). Because of these observations, it is sometimes said that "quantization is not a functor."

To better understand quantization and the space of QFTs, it seems that one must develop new tools. One such tool is to describe new quantum field theories using controlled deformations of old ones. An example of such a deformation, which has generated considerable interest in the past several years, is the $T\bar{T}$ deformation of two-dimensional QFTs. The $T\bar{T}$ operator refers to the coincident point limit

$$\mathcal{O}_{T\bar{T}}(x) = \lim_{y \to x} (T^{\mu\nu}(x)T_{\mu\nu}(y) - T^{\mu}_{\ \mu}(x)T^{\nu}_{\ \nu}(y)), \quad (1.1)$$

which was shown in [2] to define a local operator in any translation-invariant 2d QFT.

Using any such 2d QFT as a seed theory, one can define a family of theories, labeled by a flow parameter λ , which arise from deforming the seed theory by $T\bar{T}$. At the classical level, we think of this parametrized family of actions as solving the flow equation

$$\frac{\partial S_{\lambda}}{\partial \lambda} = \frac{1}{2} \int d^2 x (T^{(\lambda)\mu\nu} T^{(\lambda)}_{\mu\nu} - (T^{(\lambda)\mu}{}_{\mu})^2), \qquad (1.2)$$

where $T^{(\lambda)}_{\mu\nu}$ is the stress tensor computed from S_{λ} ,

$$T_{\mu\nu} = \frac{-2}{\sqrt{-g}} \frac{\delta S_{\lambda}}{\delta g^{\mu\nu}}.$$
 (1.3)

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However, the interpretation of the differential equation (1.2) for the classical Lagrangian can be somewhat subtle. Because $\mathcal{O}_{T\bar{T}}$ exists in the spectrum of local operators in the seed theory, deforming by this operator should lead to a well-defined quantum theory. One could ask whether this quantum theory corresponds to the quantization, in some appropriate sense, of the classical action S_{λ} .

To address this question, one can use an alternative characterization of the quantum theory obtained by a $T\bar{T}$ deformation. For instance, it is known [3,4] that the S-matrix for scattering in a $T\bar{T}$ -deformed OFT is obtained by dressing the S-matrix of the undeformed theory with a momentum-dependent phase known as a Castillejo-Dalitz-Dyson factor [5]. This gives an independent description of scattering in the quantum theory, which can be compared to predictions from quantization of the solution to (1.2). At the one-loop level, one must add specific counterterms when renormalizing the classical Lagrangian in order to reproduce the expected behavior of the $T\bar{T}$ -deformed S-matrix [6–8].¹ This suggests that the action which solves Eq. (1.2) does not, by itself, contain enough data to quantize and obtain the correct $T\bar{T}$ -deformed QFT at the quantum level; additional information from the S-matrix characterization is needed.

Another piece of evidence for this perspective comes from the observation that one can use different notions of the energy-momentum tensor in defining the flow (1.2). For instance, the Noether stress tensor is defined as the conserved current associated with spatial translations, while the Hilbert stress tensor (1.3) is defined as the variation of the action with respect to the metric; these two notions do not agree in general, and one can also consider other stress tensors that are related to these two by improvement transformations. For theories involving fermions, a direct quantization of the classical actions that solve the flow equations (1.2) driven by different definitions of the stress tensor leads to inequivalent Hilbert spaces [9,10].

We will interpret these observations by taking the following perspective. Although the solution to the classical $T\bar{T}$ flow equation (1.2) is useful, and often gives interesting hints about the nature of a $T\bar{T}$ -deformed QFT, the process of quantizing this deformed Lagrangian can be ambiguous. Indeed, as we have stressed above, quantization is not a functor: except in simple cases such as free theories, we do not understand a unique and systematic prescription for turning classical theories into quantum ones. Rather, what we *mean* by the quantum theory of a $T\bar{T}$ -deformed seed is determined by other characterizations such as the *S*-matrix or torus partition function [11–13]. These independent pieces of data should be viewed as picking out the correct

prescription for performing the quantization of the $T\bar{T}$ deformed Lagrangian. This is in analogy with the viewpoint that the proper quantization prescription for the Nambu-Goto string is the one which proceeds by first rewriting the theory in Polyakov form and then quantizing using the path integral.

It is natural to ask whether adopting this perspective offers us insights into the quantization of other models. Recently, a family of related theories that exhibit nonanalytic square-root structures in their Lagrangians has been introduced, all of which satisfy some classical flow equation similar to (1.2). We will now take a detour to describe some purely classical aspects of this collection of theories before returning to issues of quantization.

The first member of this class to be introduced was a four-dimensional gauge theory known as the modified Maxwell or ModMax model [14], which is described by the action

$$S_{\text{ModMax}}(\gamma) = \frac{1}{4} \int d^4x \left(-\cosh(\gamma) F^{\mu\nu} F_{\mu\nu} + \sinh(\gamma) \sqrt{(F^{\mu\nu} F_{\mu\nu})^2 + (F^{\mu\nu} \tilde{F}_{\mu\nu})^2} \right), \quad (1.4)$$

where $F_{\mu\nu}$ is the field strength of the Abelian gauge field A_{μ} and $\tilde{F}^{\mu\nu} = \frac{1}{2} \epsilon^{\mu\nu\rho\tau} F_{\rho\tau}$ is its Hodge dual. When $\gamma = 0$, the action (1.4) reduces to that of the usual Maxwell theory.

As a classical theory, the ModMax model (1.4) exhibits several intriguing properties. It is the unique conformally invariant and electric-magnetic duality-invariant extension of the Maxwell theory. It also satisfies a flow equation driven by a function of the energy-momentum tensor [15,16], namely

$$\frac{\partial S_{\text{ModMax}}(\gamma)}{\partial \gamma} = \frac{1}{2} \int d^4 x \sqrt{T^{(\gamma)\mu\nu} T^{(\gamma)}_{\mu\nu}}, \qquad (1.5)$$

where $T_{\mu\nu}^{(\gamma)}$ is the stress tensor of the ModMax theory (1.4) at parameter γ . Unlike the flow (1.2) for the Lagrangian of a $T\bar{T}$ -deformed 2*d* QFT, the operator on the right side of (1.5) is classically marginal. Note that, since the ModMax theory is conformally invariant and thus its stress tensor has vanishing trace, the operator driving the flow (1.5) need not have any dependence on $T^{(\gamma)\mu}{}_{\mu}$. However, it is convenient to define another combination that does involve the trace and that reduces to (1.5) in the conformal limit. For a theory in *D* spacetime dimensions with energy-momentum tensor $T_{\mu\nu}$, let

$$\mathcal{R}^{(D)} = \sqrt{\frac{1}{D} T^{\mu\nu} T_{\mu\nu} - \frac{1}{D^2} (T^{\mu}{}_{\mu})^2}.$$
 (1.6)

¹Because the $T\bar{T}$ operator is irrelevant in the sense of the renormalization group, this behavior is partly expected. Adding a generic irrelevant operator will typically activate infinitely many counterterms; the surprise is that the irrelevant $T\bar{T}$ deformation does not lead to a loss of analytic control.

In terms of the traceless part of the stress tensor, which we write as $\hat{T}_{\mu\nu} = T_{\mu\nu} - \frac{1}{D}g_{\mu\nu}T^{\rho}{}_{\rho}$, this operator is simply

$$\mathcal{R}^{(D)} = \frac{1}{\sqrt{D}} \sqrt{\hat{T}^{\mu\nu} \hat{T}_{\mu\nu}}.$$
 (1.7)

Including this dependence on the trace allows us to extend certain flow equations to nonconformal theories. For instance, there is a two-parameter family of ModMax-Born-Infeld theories labeled by couplings (λ, γ) , which reduces to (1.4) when $\lambda = 0$ and to the Born-Infeld theory when $\gamma = 0$. This family satisfies two commuting classical flow equations, one driven by a four-dimensional version of the

 $T\bar{T}$ operator and one driven by the operator $\mathcal{R}^{(4)}$ [17–19]. The operator $\mathcal{R}^{(3)}$ also appears in the flow equation that deforms the 3*d* Maxwell Lagrangian into the Born-Infeld theory in three dimensions [20].

When D = 2, the combination $\mathcal{R}^{(2)}$ is the root- $T\bar{T}$ operator introduced in [21].² Applying this deformation to the seed theory that describes N massless free scalar fields ϕ^i produces a second example of a nonanalytic classical Lagrangian. The resulting deformed theory can also be obtained from the 4*d* modified Maxwell theory by dimensional reduction [16], so we will sometimes refer to this model as the modified scalar theory. This model is described by the action

$$S_{\text{Modified Scalar}}(\gamma) = \frac{1}{2} \int d^2 x \left(\cosh(\gamma) \partial_\mu \phi^i \partial^\mu \phi^i + \sinh(\gamma) \sqrt{2 \partial_\mu \phi^i \partial_\nu \phi^i \partial^\mu \phi^j - (\partial_\mu \phi^i \partial^\mu \phi^i)^2} \right), \tag{1.8}$$

where the index i = 1, ..., N labels the N scalars. This action satisfies the flow equation

$$\frac{\partial S_{\text{Modified Scalar}}(\gamma)}{\partial \gamma} = \mathcal{R}^{(2)}, \qquad (1.9)$$

as shown in [21,26]. Like the 4*d* ModMax theory, the modified scalar theory is classically conformally invariant, and thus the trace of its stress tensor vanishes, so only the term $T^{\mu\nu}T_{\mu\nu}$ appearing under the square root in $\mathcal{R}^{(2)}$ gives a nonzero contribution.

The two-dimensional root- $T\bar{T}$ deformation that gives rise to the theory (1.8) appears to share some of the interesting properties of the $T\bar{T}$ deformation, such as preserving classical integrability in several examples [27]. However, unlike the case of $T\bar{T}$, it is not yet known how to define the root- $T\bar{T}$ deformation at the quantum level and obtain flow equations for quantities such as the *S*-matrix; proposed flow equations for the finite-volume spectrum and torus partition function of a root- $T\bar{T}$ deformed conformal field theory (CFT) were given in [28] and supported using evidence from holography, but there is no general proof of these results. This presents an obstruction to carrying out the procedure that we have described above—namely identifying the correct prescription for the quantization of these models using some additional input—for root- $T\bar{T}$ deformed models such as (1.4) and (1.8).

In this work, we take up the task of studying the quantization of such nonanalytic models in a simplified setting where one *can* carry out this program explicitly, namely in the arena of (0 + 1)-dimensional theories. By performing a particular dimensional reduction described in [29], either the ModMax theory or its modified scalar analog can be reduced to a 1*d* model that describes a harmonic oscillator with a nonanalytic interaction term. We refer to this system, which was first studied in [30], as the ModMax oscillator. The simplest version of this theory features two position variables x(t), y(t), and is described by the Lagrangian

$$L_{\text{ModMax oscillator}}(\gamma) = \frac{1}{2} \int dt \left(\cosh(\gamma) (\dot{x}^2 + \dot{y}^2 - x^2 - y^2) + \sinh(\gamma) \sqrt{((\dot{x} + y)^2 + (x - \dot{y})^2)((\dot{x} - y)^2 + (x + \dot{y})^2)} \right).$$
(1.10)

When $\gamma = 0$, this theory reduces to a two-dimensional isotropic harmonic oscillator with unit mass and frequency. This gives the third example of a nonanalytic theory.³

At first glance, it is not obvious that the full Lagrangian (1.10) at finite γ will be amenable to exact quantization because of the velocity-dependent square-root interaction. However, one might become more optimistic about the prospects of quantization after observing that this Lagrangian obeys a root- $T\bar{T}$ -like flow equation,

²We refer the reader to [22–25] for other work related to the root- $T\bar{T}$ operator.

³There are several other examples of related nonanalytic theories, which we will not discuss in detail here: a supersymmetric extension of ModMax [31,32], a 6d ModMax-like tensor theory [33], and a 4d duality-invariant supersymmetric theory that is referred to as the MadMax sigma model [34].

$$\frac{\partial L_{\text{ModMax oscillator}}(\gamma)}{\partial \gamma} = \sqrt{E_{\gamma}^2 - J_{\gamma}^2}, \qquad (1.11)$$

where E_{γ} and J_{γ} are the energy and angular momentum, respectively, of the theory (1.10) at parameter γ . This is the dimensional reduction of the flow equations driven by $\mathcal{R}^{(4)}$ and $\mathcal{R}^{(2)}$ that are obeyed by the ModMax and modified scalar models, respectively.

Because the ModMax oscillator can be described as a deformation of the harmonic oscillator by conserved charges, this suggests that one might perform canonical quantization of this theory using the prescription described in [35,36] for quantum-mechanical deformations by functions of the Hamiltonian. That is, we first choose a basis of simultaneous eigenfunctions of the Hamiltonian and total angular momentum operators in the undeformed harmonic oscillator theory. We then declare that the eigenfunctions of the harmonic oscillator, but with energy eigenvalues that have been shifted by the square-root combination appearing in (1.11).

This prescription gives a simple and elegant way to define a quantum theory of the ModMax oscillator. However, as we have emphasized, quantization is not a functor: it is not clear that this is the only prescription, or even the correct one. For instance, because this quantization scheme is so simple, one might expect that it is also possible to quantize (1.10) using the path integral formulation and get equivalent results. However, it is generally very difficult to perform the path integral for any theory that is not quadratic in derivatives. One of our goals in this work is to perform a detailed comparison of quantization prescriptions for the ModMax oscillator and check that they agree.

We will show that any deformation of a 1*d* theory by conserved charges induces a flow equation for the propagator, or partition function, of the theory. Remarkably, for the case of the deformation (1.11), this flow equation is the Laplace equation in cylindrical coordinates. We will see that this flow equation for the partition function is consistent with, and can be derived from, either the canonical quantization prescription or the phase space path integral formulation. This strategy of reformulating the quantization of a deformed theory in terms of a flow equation for the partition function, or some other quantity, might be useful for cases where one cannot quantize the classical Lagrangian or Hamiltonian directly. This is the main motivation for the present work: we perform a detailed analysis of the quantization of this simple 1d model in the hope that some of the insights from studying this problem may be useful in understanding the quantization of other nonanalytic models, such as the ModMax theory itself.⁴

The layout of this paper is as follows: In Sec. II, we study general classical deformations in a class of 1*d* theories and prove basic results such as the equivalence of deformations in the Lagrangian and Hamiltonian formulations. In Sec. III, we consider deformations of quantum-mechanical theories by conserved charges using both canonical quantization and the path integral formalism, and we derive flow equations for quantities such as the propagator and partition function. In Sec. IV, we apply the machinery developed in previous sections to the theory of the ModMax oscillator; understanding the quantum-mechanical properties of this model is the main goal of the present work. In Sec. V, we summarize our results and identify some directions for future research. A first-order check of the equivalence of Lagrangian and Hamiltonian flows appears in the Appendix.

II. CLASSICAL DEFORMATIONS

In this section, we will consider deformations of (0 + 1)-dimensional theories at the classical level. We focus on theories that describe the dynamics of a collection of real bosons $x^i(t)$, i = 1, ..., N. The generalization to theories with fermions $\psi^i(t)$ or more general degrees of freedom is straightforward, although we will not consider such cases here.⁵

We view such a theory as being defined by either a Lagrangian or a Hamiltonian,

$$L(x^{i}, \dot{x}^{i})$$
 or $H(x^{i}, p^{i})$, (2.1)

where p^i is the momentum that is canonically conjugate to the variable x^i ,

$$p^i = \frac{\partial L}{\partial \dot{x}^i}.$$
 (2.2)

Throughout this work, we assume that the indices *i*, *j*, etc., that label positions and momenta are raised or lowered with the trivial Euclidean metric δ_{ij} . We therefore do not distinguish between upstairs and downstairs indices.

We will often write expressions such as (2.1) in which the argument of a function of positions, velocities, or momenta carries an index such as *i*. In such expressions, the index *i* is not meant to be a free index, but is merely shorthand to indicate dependence on all of the corresponding variables as *i* runs from 1 to *N*. Explicitly,

$$L(x^{i}, \dot{x}^{i}) = L(x^{1}, \dots, x^{N}, \dot{x}^{1}, \dots, \dot{x}^{N}).$$
(2.3)

⁴See Ref. [37] for a discussion of equivalent classical forms of ModMax that may be useful for quantization.

⁵A convenient way to incorporate fermions is to define flow equations in superspace. These manifestly supersymmetric flows have been extensively studied; see Ref. [38] and references therein for a review of such deformations in field theory or [29,39] for the corresponding flows in 1*d* theories.

A. Flow equations for Lagrangians and Hamiltonians

Our first goal is to study deformations, or flows, in the space of classical theories, which are defined as follows: Let $\mathcal{O}(x^i, \dot{x}^i; \lambda)$ be some function of the coordinates x^i and their time derivatives, which may also depend on a real variable λ . The differential equation

$$\frac{\partial L}{\partial \lambda} = \mathcal{O}(x^i, \dot{x}^i; \lambda) \tag{2.4}$$

defines a one-parameter family of Lagrangians $L(x^i, \dot{x}^i; \lambda)$. We refer to Eq. (2.4) as a flow equation, and we say that the function \mathcal{O} is the operator which drives the flow.⁶

A theory may equivalently be described in the Hamiltonian formulation by the function

$$H(x^{i}, p^{i}) = p^{j} \dot{x}^{j} - L(x^{i}, \dot{x}^{i}), \qquad (2.5)$$

which is the Legendre transform of $L(x^i, \dot{x}^i)$. In Eq. (2.5), one must view all instances of the velocities $\dot{x}^i = \dot{x}^i(p^j)$ as being implicitly defined in terms of the conjugate momenta.

In analogy with the Lagrangian flow (2.4), we may consider a differential equation

$$\frac{\partial H}{\partial \lambda} = \mathfrak{D}(x^i, p^i; \lambda). \tag{2.6}$$

We use the symbol $\mathfrak{D}(x^i, p^i)$, which is a function of positions and canonical momenta, to distinguish it from the function $\mathcal{O}(x^i, \dot{x}^i)$ that defines the flow equation for the Lagrangian.

Because we are interested in both Lagrangian and Hamiltonian descriptions of a physical system, it is natural to ask how the flow equations (2.4) and (2.6) are related. A general deformation will modify both the Lagrangian L or Hamiltonian H and the relationship between the velocities \dot{x}^i and the conjugate momentum p^i . Because such a deformation has two effects, one should check explicitly whether the diagram



commutes, and if so, under what conditions.

Fortunately, it turns out that the Lagrangian flow equation (2.4) and Hamiltonian flow equation (2.6) lead to deformed quantities L_{λ} and H_{λ} , which are related by a Legendre transform, as expected, so long as the operators \mathcal{O} and \mathfrak{D} are related in the appropriate way. This result is stated more precisely in the following theorem.

Theorem 1. Let $L_0(x_i, \dot{x}_i)$ be a Lagrangian for a collection of coordinates $x_i(t)$, and let $H_0(x_i, p_i)$ be the corresponding Hamiltonian. Given a function $\mathcal{O}(x^i, \dot{x}^i; \lambda)$, consider a one-parameter family of Lagrangians L_{λ} that satisfy the differential equations

$$\frac{\partial L_{\lambda}}{\partial \lambda} = \mathcal{O}(x^i, \dot{x}^i; \lambda), \qquad (2.8)$$

with the initial conditions $L_{\lambda} \to L_0$ as $\lambda \to 0$. Then the Hamiltonian associated with L_{λ} ,

$$H_{\lambda}(x^{i}, p^{i}) = p^{j} \dot{x}^{j} - L_{\lambda}(x^{i}, \dot{x}^{i}), \qquad (2.9)$$

satisfies the flow equation

$$\frac{\partial H}{\partial \lambda} = \mathfrak{D}(x^i, p^i; \lambda), \qquad (2.10)$$

where the function \mathfrak{O} is defined by

$$\mathfrak{O}(x^i, p^i; \lambda) = -\mathcal{O}(x^i, \dot{x}^i(p^j; \lambda); \lambda), \qquad (2.11)$$

and where $\dot{x}^i(p^j; \lambda)$ represents the functional dependence between the velocities and conjugate momenta in the theory L_{λ} .

Conversely, given a function $\mathfrak{D}(x^i, p^i; \lambda)$, consider the family of Hamiltonians H_{λ} that obey

$$\frac{\partial H_{\lambda}}{\partial \lambda} = \mathfrak{O}(x^i, p^i; \lambda), \qquad (2.12)$$

with initial condition $H_{\lambda} \to H_0$ as $\lambda \to 0$. The Lagrangians L_{λ} associated with H_{λ} ,

$$L_{\lambda}(x^{i}, \dot{x}^{i}; \lambda) = p^{j} \dot{x}^{j} - H_{\lambda}(x^{i}, \dot{x}^{i}; \lambda), \qquad (2.13)$$

satisfy the flow equation (2.8), where the operator O is defined by

$$\mathcal{O}(x^i, \dot{x}^i; \lambda) = -\mathfrak{O}(x^i, p^i(\dot{x}^j; \lambda); \lambda), \qquad (2.14)$$

and where $p^i(\dot{x}^j; \lambda)$ represents the functional dependence between the conjugate momenta and velocities in the theory H_{λ} .

The interpretation of this theorem is that the diagram (2.7) commutes, so long as the Lagrangian deformation \mathcal{O} and the Hamiltonian deformation \mathfrak{D} are correctly related using the constraint between the velocities and conjugate momenta in the theory at finite λ . Note that this is not the

⁶The use of the term "operator" is motivated by similar flow equations in field theories, such as the deformation of a 2*d* QFT by the $T\bar{T}$ operator. In the present context, O is not a true quantum-mechanical operator acting on a Hilbert space, but merely a classical function of positions and velocities.

same as using the relationship between p^i and \dot{x}^i in the undeformed theories L_0 and H_0 . Unsurprisingly, if one uses the relation between \dot{x}^i and p^i , which is valid in the seed theories, the corresponding flows commute only to leading order in the deformation parameter λ . This was proven for field theories describing a single field ϕ and conjugate momentum $\pi = \frac{\partial \mathcal{L}}{\partial \phi}$ in Appendix A of [40]. To make the present work self-contained, we include the analog of their leading-order proof for (0 + 1)-dimensional theories of N positions $x^i(t)$ in our Appendix.

Proof. We first prove the forward direction, beginning with the flow equation (2.8) for the Lagrangian. We view the velocities \dot{x}^i as independent variables, while at each value of λ , the conjugate momenta $p^i(\dot{x}^i; \lambda)$ are determined via the relation

$$p^{i}(\dot{x}^{i};\lambda) = \frac{\partial L_{\lambda}}{\partial \dot{x}^{i}}.$$
 (2.15)

The Legendre transform that defines the Hamiltonian, written in a way which emphasizes the λ dependence, is

$$H_{\lambda}(x^{i}, p^{i}(\dot{x}^{j}; \lambda)) = p^{j}(\dot{x}^{k}; \lambda)\dot{x}^{j} - L_{\lambda}(x^{i}, \dot{x}^{i}).$$
(2.16)

We differentiate both sides of (2.16) with respect to λ to find

$$\frac{\partial H_{\lambda}}{\partial p^{i}}\frac{\partial p^{i}}{\partial \lambda} + \frac{\partial H_{\lambda}}{\partial \lambda} = \frac{\partial p^{j}}{\partial \lambda}\dot{x}^{j} - \frac{\partial L_{\lambda}}{\partial \lambda}.$$
 (2.17)

By the Hamilton equations of motion, we have

$$\frac{\partial H_{\lambda}}{\partial p^{i}} = \dot{x}^{i}, \qquad (2.18)$$

and thus the terms $\frac{\partial p^j}{\partial \lambda} \dot{x}^j$ on both sides of Eq. (2.17) cancel, leaving

$$\frac{\partial H_{\lambda}}{\partial \lambda} = -\frac{\partial L_{\lambda}}{\partial \lambda} = -\mathcal{O}(x^{i}, \dot{x}^{i}(p^{j}; \lambda); \lambda).$$
(2.19)

The object on the right side of (2.19) is precisely the operator $\mathfrak{D}(x^i, p^i; \lambda)$. This completes the first half of the proof.

Now we show the reverse direction. Suppose that the Hamiltonian obeys the flow equation (2.12). We now view the p^i as independent variables while the \dot{x}^i are fixed as

$$\dot{x}^{i}(p^{i};\lambda) = \frac{\partial H_{\lambda}}{\partial p^{i}}.$$
(2.20)

Again making all functional dependence explicit, the Legendre transform that defines the Lagrangian is

$$L_{\lambda}(x^{i}, \dot{x}^{i}(p^{j}; \lambda)) = p^{j} \dot{x}^{j}(p^{i}; \lambda) - H_{\lambda}(x^{i}, p^{i}).$$
(2.21)

Differentiating with respect to λ gives

$$\frac{\partial L_{\lambda}}{\partial \dot{x}^{i}} \frac{\partial \dot{x}^{i}}{\partial \lambda} + \frac{\partial L_{\lambda}}{\partial \lambda} = p^{j} \frac{\partial \dot{x}^{j}}{\partial \lambda} - \frac{\partial H_{\lambda}}{\partial \lambda}.$$
 (2.22)

After using the relation

$$\frac{\partial L_{\lambda}}{\partial \dot{x}^{i}} = p^{i}, \qquad (2.23)$$

we see that the terms $p^j \frac{\partial \dot{x}^j}{\partial \lambda}$ cancel on either side of Eq. (2.22), and we are left with

$$\frac{\partial L_{\lambda}}{\partial \lambda} = -\frac{\partial H_{\lambda}}{\partial \lambda} = -\mathfrak{D}(x^{i}, p^{i}(\dot{x}^{j}; \lambda); \lambda), \qquad (2.24)$$

which establishes the converse.

Note that we have stated this theorem and its proof for (0 + 1)-dimensional theories. However, one can repeat this argument almost verbatim, making the replacements

$$\begin{aligned} x^{i} &\to \phi^{i}, \qquad p^{i} \to \pi^{i}, \qquad L(x^{i}, \dot{x}^{i}) \to \mathcal{L}(\phi^{i}, \dot{\phi}^{i}), \\ H(x^{i}, p^{i}) \to \mathcal{H}(\phi^{i}, \pi^{i}), \end{aligned} \tag{2.25}$$

to obtain the corresponding theorem and its proof in any *d*-dimensional quantum field theory for a collection of fields ϕ^i and their conjugate momenta $\pi^i = \frac{\partial \mathcal{L}}{\partial \dot{\phi}^i}$.

1. Examples of classical Lagrangian and Hamiltonian deformations

It is instructive to see how Lagrangian and Hamiltonian flows are related in several examples, both when the assumptions of Theorem 1 hold and when they do not.

First, let us consider a nonexample of this theorem; we emphasize that this is only a toy problem meant for purposes of illustration, and not a flow of primary interest in this work. We begin from an undeformed theory that has only a free kinetic term:

$$L_0 = \frac{1}{2}m\dot{x}^2, \qquad H_0 = \frac{p^2}{2m}.$$
 (2.26)

The relationship between the undeformed velocity and its conjugate momentum is simply $p = m\dot{x}$. Now consider the pair of flows

$$\frac{\partial L_{\lambda}}{\partial \lambda} = \mathcal{O}(x, \dot{x}) = m\dot{x}, \qquad \frac{\partial H_{\lambda}}{\partial \lambda} = \mathfrak{O}(x, p) = -p, \quad (2.27)$$

with solutions

$$L_{\lambda} = \frac{1}{2}m\dot{x}^2 + \lambda m\dot{x}, \qquad H_{\lambda} = \frac{p^2}{2m} - \lambda p. \qquad (2.28)$$

The Lagrangian L_{λ} and Hamiltonian H_{λ} are not related by a Legendre transformation. The conjugate momentum evaluated using the Lagrangian L_{λ} is

$$p_{\lambda} = \frac{\partial L_{\lambda}}{\partial \dot{x}} = m \dot{x} + \lambda m, \qquad (2.29)$$

and thus the Legendre transform of L_{λ} , which we call \tilde{H}_{λ} , is

$$\tilde{H}_{\lambda} = p_{\lambda}\dot{x} - L_{\lambda} = \frac{p^2}{2m} - \lambda p + \frac{1}{2}m\lambda^2.$$
(2.30)

In this simple example, the difference between the Legendre transform \tilde{H}_{λ} and the Hamiltonian H_{λ} is only a constant term, but nonetheless the two quantities do not agree. However, the difference $\tilde{H}_{\lambda} - H_{\lambda}$ is of order λ^2 . Thus H_{λ} and \tilde{H}_{λ} agree to leading order in λ , which is consistent with the argument in the Appendix. In this example, the reason that the flows agree only at leading order is because the operators $\mathcal{O} = m\dot{x}$ and $\mathfrak{D} = -p$ only satisfy the required constraint $\mathcal{O} = -\mathfrak{D}$ if we use the relation between the velocity and conjugate momentum in the undeformed theories.

Next let us consider a modification of the above flow:

$$\frac{\partial L_{\lambda}}{\partial \lambda} = \mathcal{O}(x, \dot{x}) = \frac{\partial L_{\lambda}}{\partial \dot{x}}, \quad \frac{\partial H_{\lambda}}{\partial \lambda} = \mathfrak{O}(x, p) = -p. \quad (2.31)$$

To lowest order around the initial condition (2.26), the deforming operators of equation (2.31) agree with those in (2.27) since $\frac{\partial L_0}{\partial \dot{x}} = m\dot{x}$. Beyond first order, however, the flow defined by (2.31) has been altered in order to satisfy the assumptions of Theorem 1. We will solve these differential equations with initial conditions that are arbitrary functions of velocities or momenta, $L_0(\dot{x})$ and $H_0(p)$ [although note that one must specialize to the case of $L_0(\dot{x}) = \frac{1}{2}m\dot{x}^2$ for comparison to (2.28)]. One finds

$$L_{\lambda}(\dot{x}) = L_0(\dot{x} + \lambda), \qquad H_{\lambda}(p) = H_0 - \lambda p. \quad (2.32)$$

The Lagrangian L_{λ} and Hamiltonian H_{λ} of (2.32) are indeed related by a Legendre transform⁷ to all orders in λ , as guaranteed by Theorem 1.

Let us consider one more, slightly less trivial, example. Again beginning from an arbitrary velocity-dependent seed Lagrangian $L_0(\dot{x})$ and Hamiltonian $H_0(p)$, consider the flow equations

$$\frac{\partial L_{\lambda}}{\partial \lambda} = \left(L_{\lambda} - \dot{x} \frac{\partial L_{\lambda}}{\partial \dot{x}} \right)^2, \qquad \frac{\partial H_{\lambda}}{\partial \lambda} = -H^2.$$
(2.33)

This deformation was first considered in Appendix A of [35]. The Hamiltonian flow equation can be solved for any initial condition $H_0(x, p)$:

$$H_{\lambda} = \frac{H_0}{1 + \lambda H_0}.$$
 (2.34)

However, the Lagrangian flow equation is more complicated. If the seed theory is $L_0 = \dot{x}^2$, the solution is given in terms of a hypergeometric function:

$$L_{\lambda} = \frac{3}{4\lambda} \left({}_{3}F_{2} \left[-\frac{1}{2}, -\frac{1}{4}, \frac{1}{4}; \frac{1}{3}, \frac{2}{3}; \frac{256}{27} \lambda \dot{x}^{2} \right] - 1 \right).$$
(2.35)

This same hypergeometric function has appeared in several contexts related to classical deformations by conserved charges, including the $T\bar{T}$ deformation of the 2*d* Maxwell theory [17] and Yang-Mills [41]. Despite the complicated form of L_{λ} , one can check that it is indeed related to the simpler function H_{λ} of (2.34) by a Legendre transform when $H_0 = \frac{1}{4}p^2$. This is again required by the general argument of Theorem 1.

B. Deformations by conserved charges

The two examples (2.31) and (2.33) considered in the previous section are especially natural because they correspond to deformations of the theory by conserved quantities. Indeed, in Eq. (2.31) we deform the Lagrangian or Hamiltonian by the conjugate momentum p, which is conserved because the cyclic coordinate x does not appear in the Lagrangians $L_{\lambda}(\dot{x})$. Likewise, Eq. (2.33) is a deformation involving the Noether charge associated with time translation symmetry, which is the total energy of the system.

Such deformations by conserved charges are convenient to work with, since they give us a straightforward way to satisfy the conditions of Theorem 1. This is because a conserved quantity, associated with a particular symmetry, can easily be characterized in either the Lagrangian or the Hamiltonian formalism. In the former case we use Noether's theorem. If the Lagrangian is shifted by a total time derivative under the action of a symmetry generator δ ,

$$\delta L = \frac{df}{dt},\tag{2.36}$$

then the corresponding Noether charge

$$Q = \frac{\partial L}{\partial \dot{x}^i} \delta x^i + f \tag{2.37}$$

obeys $\frac{dQ}{dt} = 0$ when the equations of motion are satisfied. Likewise, in the Hamiltonian formulation—assuming that the charge Q does not depend explicitly on time—the relation

⁷This is an elementary property of the Legendre transform under translations. Let $f^*(p)$ represent the Legendre transform of a function f(x). If f(x) = g(x + y), then $f^*(p) = g^*(p) - py$.

$$\frac{dQ}{dt} = \{Q, H\} = 0$$
 (2.38)

holds on-shell.

Such conserved quantities typically have a clear physical interpretation, such as an energy or angular momentum, which makes them easy to describe either as functions of x^i and \dot{x}^i or as functions of x^i and p^i . This is in contrast to a deformation of the Lagrangian by an arbitrary combination $\mathcal{O}(x^i, \dot{x}^i)$ of kinematical variables, for which one would have to explicitly work out the dependence $\dot{x}^i(p^j)$ in order to find the corresponding Hamiltonian deformation $\mathfrak{O}(x^i, p^i)$. Because of the naturalness of deformations by conserved charges, and their relationship to interesting higher-dimensional deformations such as $T\bar{T}$ and root- $T\bar{T}$, we will focus on this class of flows in the remainder of this work.

More precisely, what we mean by a deformation by conserved charges is the following. Suppose that a seed theory L_0 has a collection of symmetries generated by variations δ_a and which are associated with conserved charges Q_a , for a = 1, ..., M, according to Eq. (2.37). We will always use early Latin indices such as a, b, c to label charges Q_a and middle Latin indices such as i, j, k to refer to coordinates x^i . We would like to define a flow equation of the form

$$\frac{\partial L_{\lambda}}{\partial \lambda} = f\left(Q_1^{(\lambda)}(x^i, \dot{x}^i), \dots, Q_M^{(\lambda)}(x^i, \dot{x}^i)\right), \quad (2.39)$$

where $Q_a^{(\lambda)}$ is the Noether charge associated with the symmetry δ_a in the theory L_{λ} . The corresponding Hamiltonian flow is

$$\frac{\partial H_{\lambda}}{\partial \lambda} = -f\left(Q_1^{(\lambda)}(x^i, p^i), \dots, Q_M^{(\lambda)}(x^i, p^i)\right).$$
(2.40)

In this equation, the Hamiltonian charges $Q_a^{(\lambda)}(x^i, p^i)$ are obtained by expressing the corresponding Lagrangian charges $Q_a^{(\lambda)}(x^i, \dot{x}^i)$ in terms of conjugate momenta, using the relation between p^i and \dot{x}^i in the theory L_{λ} .

We must make an additional assumption in order to define such a flow: the variations δ_a must generate symmetries of the entire family of theories L_{λ} , rather than only for the seed theory L_0 . That is, we must assume that this deformation does not break any of the symmetries. This will be true if all of the charges Q_a are Poisson commuting,

$$\{Q_a, Q_b\} = 0. \tag{2.41}$$

This assumption is sufficient because, if Q_a is the Noether charge associated with a symmetry variation δ_a , then the Poisson bracket with Q_a generates the transformation of any function g as $\{Q_a, g\} = \delta_a g$. Thus the condition (2.41) implies that all of the charges Q_a are invariant under all of the symmetries generating these charges, and thus any deformation of a Lagrangian by a function of these charges will still enjoy the same symmetries.

For instance, we could consider a theory with two coordinates $x^1 = x$ and $x^2 = y$ and which has two conserved momenta p_x and p_y along with a conserved angular momentum J and a Hamiltonian H. In this case,

$$\{H, J\} = \{H, p_x\} = \{H, p_y\} = 0, \qquad (2.42)$$

but one has

$$\{J, p_x\} = p_y, \qquad \{J, p_y\} = -p_x.$$
 (2.43)

In this case, we cannot define a flow equation (2.39) if, for instance, the function f depends on J and p_x but not p_y . Such a deformation breaks the rotational symmetry between x and y, and thus J is no longer a conserved quantity in the deformed theory. However, we are free to construct a flow equation using the three quantities

$$Q_1 = H,$$
 $Q_2 = J,$ $Q_3 = p_x^2 + p_y^2,$ (2.44)

since

$$\{J, p_x^2 + p_y^2\} = 0, \qquad (2.45)$$

and thus these three charges Q_a Poisson commute.

The first examples of deformations by conserved charges, which we will also call $f(Q_a)$ flows, are those driven only by a function of the Hamiltonian H. This is the class of f(H) deformations considered in [35,36], and although this class of deformations has been studied in several previous works, we will review some results concerning these flows in our paper to illustrate our general formalism. The class of f(H) deformations includes the dimensional reduction of the $T\bar{T}$ deformation, in the special case where the 2d seed theory is conformally invariant. This dimensional reduction leads to the flow equation

$$\partial_{\lambda}H_{\lambda} = \frac{H_{\lambda}^2}{\frac{1}{2} - 2\lambda H_{\lambda}},\qquad(2.46)$$

which has the solution

$$H_{\lambda} = \frac{1}{4\lambda} \left(1 - \sqrt{1 - 8\lambda H_0} \right). \tag{2.47}$$

See Ref. [42] for the extension of these f(H) flows to the case of non-Hermitian Hamiltonian deformations. Another class of examples is of those that involve the product of the conserved energy E and a second charge Q, which were studied in [43] and which are similar to the $J\bar{T}$ deformations of 2d field theories.

However, in the present work we will be primarily interested in the class of $f(E, J^2)$ deformations studied in [29], which was motivated by the flow equation obeyed by the ModMax oscillator of [30]. These deformations can be applied to theories that enjoy an additional SO(N)symmetry which rotates the N coordinates x^i as

$$x^{i}(t) \rightarrow R^{i}{}_{j}x^{j}(t), \qquad R \in SO(N).$$
 (2.48)

The conserved currents associated with each of the rotation generators is a component of angular momentum,

$$J_{nm} = \frac{\partial L}{\partial \dot{x}^n} x_m - \frac{\partial L}{\partial \dot{x}^m} x_n.$$
(2.49)

We define the total angular momentum by

$$J^2 = J^{nm} J_{nm}.$$
 (2.50)

A general deformation by a function of both the energy and angular momentum, written in the Lagrangian formulation, is then

$$\frac{\partial L_{\lambda}}{\partial \lambda} = \mathcal{O}(E, J^2), \qquad (2.51)$$

for some function \mathcal{O} . We can, of course, write an equivalent flow equation for the Hamiltonian, $\partial_{\lambda}H_{\lambda} = \mathfrak{D}(E, J^2)$, where it is understood that E and J^2 are functions of the positions x^i and canonical momenta p^i in the Hamiltonian flow.

The main operator of interest in the present work is

$$\mathcal{R} = \sqrt{E(x^{i}, \dot{x}^{i})^{2} - J(x^{i}, \dot{x}^{i})^{2}}$$

or $\Re = \sqrt{H(x^{i}, p^{i})^{2} - J(x^{i}, p^{i})^{2}},$ (2.52)

where, following the conventions for \mathcal{O} and \mathfrak{D} , we use calligraphic letters to refer to operators that depend on configuration space variables and Fraktur symbols for functions of phase space coordinates. In physically interesting examples, such as the *N*-dimensional harmonic oscillator, one always has the bound $|E|^2 \ge |J|^2$ so that the argument of the square root in (2.52) is non-negative; we will assume this to be true in what follows. We will always use the symbol γ for the flow parameter when deforming by the operator \mathcal{R} , in contrast with λ , which we use as the parameter for a general deformation.

We refer to the combination (2.52), in either formulation, as the 1*d* root- $T\bar{T}$ operator. The reason for using this term is that, as shown in [29], this object arises from a certain dimensional reduction of the flow equation

$$\frac{\partial S_{\gamma}}{\partial \gamma} = \int d^2 x \sqrt{\frac{1}{2} T^{(\gamma)\mu\nu} T^{(\gamma)}_{\mu\nu} - \frac{1}{4} (T^{(\gamma)\mu}{}_{\mu})^2}, \qquad (2.53)$$

which defines the root- $T\bar{T}$ deformation of (1 + 1)dimensional field theories [21]. More precisely, given a field theory describing the dynamics of a collection of scalar fields $\phi^i(x, t)$ on a spatial circle $x \sim x + 2\pi R$, one can Fourier expand each scalar field as

$$\phi^{i}(x,t) = \sum_{n=-\infty}^{\infty} c_{n}^{i}(t) \exp\left(\frac{inx}{R}\right).$$
(2.54)

Truncating the theory to the dynamics of a single nonzero mode $c_m^i(t)$, m > 0, and integrating over the circle then yields a dimensionally reduced theory for the functions $c_m^i(t)$. Performing this reduction for the family of theories that arises from deforming a collection of free scalars by the root- $T\bar{T}$ operator (2.53) then yields a family of 1*d* theories which satisfy a flow equation driven by the combination (2.52). In particular, applying this 1*d* root- $T\bar{T}$ deformation to a seed theory of *N* bosons $x^i(t)$ subject to a harmonic oscillator potential yields the theory that we refer to as the ModMax oscillator. This will be the subject of Sec. IV.

To conclude this subsection, we point out that—although deformations by conserved charges are quite general—not all models of interest satisfy flow equations of this form. Interesting nonexamples include the Born oscillator and generalized Born oscillator, which were recently studied in [44,45]. A version of the Born oscillator for N coordinates x^i can be described by the Hamiltonian

$$H_{\lambda} = \frac{1}{\lambda} \left(\sqrt{(1 + \lambda p^i p^i)(1 + \lambda x^i x^i)} - 1 \right).$$
(2.55)

Despite its very symmetrical form, this Hamiltonian has the property that

$$\{H_{\lambda}, \partial_{\lambda} H_{\lambda}\} \neq 0, \tag{2.56}$$

and thus it cannot obey any flow equation of the form $\partial_{\lambda}H_{\lambda} = \mathfrak{D}(H_{\lambda}, Q_a).$

C. Flow of the classical partition function

We are ultimately interested in quantum observables, such as the propagator, for theories deformed by functions of conserved charges, which will be studied in Sec. III. The periodic Euclidean-time propagator reproduces the quantum thermal partition function, and the classical limit of this object is the ordinary classical partition function. It will therefore be useful to study the classical partition function in order to have a check against which to compare the results of Sec. III, since these quantities should agree in the limit $\hbar \rightarrow 0$. We will see that the flow equations satisfied by the classical and quantum partition functions under a general deformation by conserved charges are identical.

Consider a theory with Hamiltonian $H(x^i, p^i)$, for i = 1, ..., N, and a deformation by a function of conserved

charges Q_a , a = 1, ..., M, of the form in Eq. (2.40). One could also define $Q_0 = H$ to be the conserved charge associated with time translations, which would give M + 1 charges in total. We define the grand canonical partition function

$$\mathcal{Z}(\beta,\lambda,\mu_a) = \frac{1}{(2\pi\hbar)^N} \int dx^1 \cdots dx^N dp^1 \cdots dp^N$$
$$\times \exp\left(-\beta H_\lambda(x^i,p^i) + \sum_{a=1}^M \mu_a Q_a\right). \quad (2.57)$$

Here $\beta = \frac{1}{T}$ is the inverse temperature and μ_a is the chemical potential, or fugacity, for the conserved charge Q_a . Note that the sum over charges in (2.57) begins at a = 1 so that the Hamiltonian is not included, although we could treat the Hamiltonian symmetrically by defining $\mu_0 = -\beta$ and beginning the sum at a = 0. From now on, we will set $\hbar = 1$.

It is convenient to define the expectation value of a function f in this ensemble as

$$\langle f(x^j, p^j) \rangle = \frac{1}{(2\pi)^N} \int dx^1 \cdots dx^N dp^1 \cdots dp^N f(x^j, p^j)$$
$$\times \exp\left(-\beta H_{\lambda}(x^i, p^i) + \sum_{b=1}^M \mu_b Q_b\right). \quad (2.58)$$

The derivative of the partition function with respect to λ is

$$\partial_{\lambda} \mathcal{Z} = -\beta \langle \partial_{\lambda} H_{\lambda} \rangle = -\beta \langle \mathfrak{O}(H_{\lambda}, Q_a) \rangle, \qquad (2.59)$$

and the derivatives with respect to the inverse temperature and chemical potentials are

$$\partial_{\beta} \mathcal{Z} = -\langle H_{\lambda} \rangle, \qquad \partial_{\mu_a} \mathcal{Z} = \langle Q_a \rangle.$$
 (2.60)

The partition function therefore obeys the flow equation

$$\partial_{\lambda} \mathcal{Z} = -\beta \mathfrak{D}[-\partial_{\beta}, \partial_{\mu_1}, \dots, \partial_{\mu_M}](\mathcal{Z}).$$
(2.61)

The right side of (2.61) is defined by expanding the operator \mathfrak{D} as a power series in each of its variables H_{λ} and Q_a , and then replacing each variable with the appropriate derivative. This is possible only if the deforming operator is an analytic function of its arguments, but we will see shortly how to extend this argument to some nonanalytic deformations.

For example, let us consider the flow equation for the one-dimensional $T\bar{T}$ deformation,

$$\partial_{\lambda}H_{\lambda} = \frac{H_{\lambda}^2}{\frac{1}{2} - 2\lambda H_{\lambda}} = 2H_{\lambda}^2 + 8\lambda H_{\lambda}^3 + 32\lambda^2 H_{\lambda}^4 + 128\lambda^3 H_{\lambda}^5 + \cdots.$$
(2.62)

We may schematically write the corresponding flow equation for the partition function as

$$\partial_{\lambda} \mathcal{Z} = -\beta \left(\frac{\partial_{\beta}^2}{\frac{1}{2} + 2\lambda \partial_{\beta}} \right) \mathcal{Z}, \qquad (2.63)$$

where the rational function of derivatives is again defined by the Taylor series expansion whose first few terms are shown in (2.62). We can invert this infinite series of derivatives to write the equivalent flow equation

$$\left(\frac{1}{2} + 2\lambda\partial_{\beta}\right)\left(\frac{1}{\beta}\partial_{\lambda}\mathcal{Z}\right) = -\partial_{\beta}^{2}\mathcal{Z}, \qquad (2.64)$$

which can be expressed as

$$\left(4\lambda\partial_{\lambda}\partial_{\beta}+2\beta\partial_{\beta}^{2}+\left(1-\frac{4\lambda}{\beta}\right)\partial_{\lambda}\right)\mathcal{Z}(\lambda,\beta)=0.$$
 (2.65)

This is the flow equation for the classical partition function of a theory deformed by the $1d T\bar{T}$ flow. It also turns out that the quantum partition function obeys the same flow equation.⁸ We will see in Sec. III that this is true more generally: the flow equations for the classical and quantum partition functions are always identical for deformations by any function of conserved charges.

An interesting feature of the differential equation (2.65) is that its solution can be written as an integral transform of the undeformed partition function $Z_0(\beta)$ at $\lambda = 0$. This can be understood as an analog of the solution to the heat equation, written as a convolution against the heat kernel, since (2.65) takes the form of a diffusion-type equation where the parameter λ plays the role of time. This integral kernel solution is

$$Z_{\lambda}(\beta) = \int_0^{\infty} d\beta' \frac{\beta}{\sqrt{-8\pi\lambda}\beta'^{3/2}} \exp\left(\frac{(\beta - \beta')^2}{8\lambda\beta'}\right) Z_0(\beta'),$$
(2.66)

which was obtained and studied in [36]. It is easy to check that the integral expression (2.66) automatically solves the differential equation (2.65).

We now turn our attention to deformations that take the form

$$\frac{\partial H_{\lambda}}{\partial \lambda} = \sqrt{f(H_{\lambda}, Q_a)}, \qquad (2.67)$$

where f is an analytic function. In the case where $f(H_{\lambda} = 0, Q_a = 0) = 0$, this deforming operator does not admit a Taylor series expansion because the square-root

⁸For the negative sign of the deformation parameter, this also matches the differential equation one obtains from studying Jackiw-Teitelboim gravity with a finite radial cutoff [46].

function is not analytic around 0. For such deformations, we cannot define the differential operator appearing on the right side of Eq. (2.61) via a series.

We can attempt to circumvent this difficulty in one of two ways. The first way is to attempt to define a fractional derivative by diagonalizing the differential operator. That is, if we can identify a complete basis of eigenfunctions ψ_n with the property

$$f(-\partial_{\beta},\partial_{\mu_a})\psi_n = \nu_n\psi_n, \qquad (2.68)$$

for some non-negative eigenvalues ν_n , then we simply define the fractional differential operator to act as

$$\sqrt{f(-\partial_{\beta},\partial_{\mu_a})}\psi_n = \sqrt{\nu_n}\psi_n.$$
(2.69)

We may then expand the partition function \mathcal{Z} in this basis as

$$\mathcal{Z} = \sum_{n} c_n \psi_n \tag{2.70}$$

and obtain a flow equation

$$\partial_{\lambda} \mathcal{Z} = -\beta \sum_{n} \sqrt{\nu_n} \, c_n \psi_n. \tag{2.71}$$

Although it should be possible, in principle, to carry out this procedure of defining the fractional derivative—at least in some examples—we will not pursue it further here.

Instead, we will attempt to remedy the nonanalyticity by taking a second derivative. In certain cases, this can convert a first-order flow equation driven by a square-root operator to a second-order flow equation driven by an analytic operator. For a general deformation by an operator $\mathfrak{D} = \mathfrak{D}(H_{\lambda}, Q_a)$, the second derivative of the partition function with respect to λ is

$$\partial_{\lambda}^{2} \mathcal{Z} = \left\langle -\beta \frac{\partial \mathfrak{D}}{\partial \lambda} + \beta^{2} \mathfrak{D}^{2} \right\rangle, \qquad (2.72)$$

and for an operator $\mathfrak{D} = \sqrt{f(H_{\lambda}, Q_a)}$ of the form in Eq. (2.67), this is

$$\partial_{\lambda}^{2} \mathcal{Z} = \left\langle \frac{-\beta}{2} \frac{\partial_{\lambda} f}{\sqrt{f}} + \beta^{2} f \right\rangle.$$
(2.73)

The second term now depends only on f but not its square root, so this term can be expressed in terms of a power series in derivatives of \mathcal{Z} with respect to β and the μ_a as before. The first term, however, still depends on \sqrt{f} and is not manifestly analytic for a deformation involving a generic function f. However, we will revisit this expression in Sec. IV for the special case $f(H_{\lambda}, Q_a) = \sqrt{H_{\lambda}^2 - J_{\lambda}^2}$ and the

III. QUANTUM DEFORMATIONS

solution such as (2.66), its general solution can be written

in terms of exponentials and Bessel functions.

In this section we turn to the deformation of quantummechanical theories by conserved charges in (0 + 1) spacetime dimensions. As we mentioned in the Introduction, there is no universal method for quantizing a general classical theory; see, for instance, [47] for a survey of quantization methods from a mathematical perspective.

Here we will focus on canonical quantization and path integral quantization. When discussing operators in the canonical formalism, we will use hats to distinguish them from the corresponding classical variables; for instance, we write

$$\hat{x}^i | \vec{x} \rangle = x^i | \vec{x} \rangle, \qquad \hat{p}^i | \vec{p} \rangle = p^i | \vec{p} \rangle.$$
 (3.1)

We will use vector symbols \vec{x} and \vec{p} to represent the collection of all components x^i and p^i , respectively, for i = 1, ..., N. The component indices i, j, etc., are not to be confused with the subscripts A, B, which we will introduce shortly and which refer to the initial and final configurations that determine the boundary conditions of the path integral.

One of the results of Sec. II is that deformations by conserved charges in the Hamiltonian and Lagrangian formulations are equivalent, if one uses the correct relationship between variables in the deforming operators. Because the quantum theory features either the Hamiltonian, in canonical quantization, or the Lagrangian, in conventional path integral quantization, a natural first question to ask is whether deformations by conserved charges in these two formalisms are again equivalent quantum mechanically.

For example, one might ask whether computing the propagator using the unitary time evolution operator $\hat{U}(t_B, t_A)$ associated with a given Hamiltonian $\hat{H}(t)$,

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A) = \langle \vec{x}_B | \hat{U}(t_B, t_A) | \vec{x}_A \rangle,$$
$$\frac{\partial \hat{U}(t, t')}{\partial t} \Big|_{t=t'} = -i\hat{H}(t), \qquad (3.2)$$

agrees with a definition using the Feynman path integral,

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A) = \int_{\vec{x}(t_A) = \vec{x}_A}^{\vec{x}(t_B) = \vec{x}_B} \mathcal{D}\vec{x} \exp\left(i \int_{t_A}^{t_B} dt L_0\right), \quad (3.3)$$

after deforming \hat{H}_0 by an operator $\hat{\mathfrak{D}}(\hat{Q}_a)$ in (3.2) and deforming L_0 by the corresponding operator $\mathcal{O}(Q_a) = -\mathfrak{D}(Q_a)$ in (3.3).

However, as we will review, this is not the right question to ask. The expression (3.3) for the propagator, in terms of a path integral over the position coordinates \vec{x} with the standard measure, is only valid for Lagrangians that are quadratic in derivatives. After deforming a seed theory by a function of conserved charges, the resulting deformed Lagrangian will often have more general dependence on the derivatives \dot{x}^i . In such situations, the conventional path integral (3.3) does not correctly compute the deformed propagator, and instead one must use a phase space path integral:

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A) = \int_{\vec{x}(t_A) = \vec{x}_A}^{\vec{x}(t_B) = \vec{x}_B} \mathcal{D}\vec{x} \int_{\vec{p}(t_A) = \vec{p}_A}^{\vec{p}(t_B) = \vec{p}_B} \mathcal{D}\vec{p} \\ \times \exp\left(i \int_{t_A}^{t_B} dt \left(p^i(t) \dot{x}_i(t) - H_0(\vec{p}, \vec{x})\right)\right),$$
(3.4)

The quantity appearing in the exponential of (3.4) is *not* the classical action, because the functions $p^i(t)$ in the first term are not the canonical momenta, but rather dummy functions that are path-integrated over. Thus, for a general non-quadratic action, the classical Lagrangian plays no role in the path integral quantization of the theory, and only the phase space path integral defined in (3.4) is important.

Given this observation, we should not ask whether deformations of the Hamiltonian and Lagrangian are equivalent in the quantum theory. Instead, we should ask whether deforming the Hamiltonian operator in the expression (3.2) is equivalent to deforming the classical function of phase space variables $H_0(\vec{p}, \vec{x})$ in (3.4). This will be the topic of Sec. III B. First, we will take a detour to review the phase space path integral.

A. Phase space path integral

In this section, we will review the path integral quantization of a general Hamiltonian

$$\hat{H}(\hat{x}^i, \hat{p}^i), \tag{3.5}$$

which need not be quadratic in the momenta \hat{p}^i that are conjugate to the position operators \hat{x}^i . Only in the case of this quadratic dependence on momenta does the general phase space path integral reduce to the ordinary Feynman path integral.⁹

To begin, we will assume that the Hamiltonian is an analytic function of the variables \hat{x}^i and \hat{p}^i . Later we will be

interested in deformations of such Hamiltonians by nonanalytic functions of charges, but for now we will require \hat{H} to admit an expansion

$$\hat{H}(t;\vec{x},\vec{p}) = \sum_{i,j,m,n} h_{ijmn}(t) (\hat{p}^i)^m (\hat{x}^j)^n.$$
(3.6)

One can always bring a general analytic Hamiltonian into this form by using the canonical commutation relation to move all position operators \hat{x}^i to the right of momentum operators \hat{p}^i . The form (3.6) of the Hamiltonian is said to be normal ordered. Taking the Hermitian conjugate of this expression reverses the order of the operators,

$$\hat{H}^{\dagger}(t;\vec{x},\vec{p}) = \sum_{i,j,m,n} h^*_{ijmn}(t) (\hat{x}^j)^n (\hat{p}^i)^m, \qquad (3.7)$$

and is thus antinormal ordered. We also assume that the Hamiltonian is Hermitian, so that $\hat{H} = \hat{H}^{\dagger}$. Our conventions for the position and momentum eigenstates are

$$\langle \vec{x} | \vec{p} \rangle = \exp\left(ix^i p_i\right),\tag{3.8}$$

and we take $\hbar = 1$.

We are interested in computing the propagator

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A) = \langle \vec{x}_B | \hat{U}(t_B, t_A) | \vec{x}_A \rangle, \qquad (3.9)$$

where the unitary time evolution operator \hat{U} is related to the Hamiltonian according to (3.2). Following the usual time-slicing procedure, we subdivide the time interval $T = t_B - t_A$ into a large number M + 1 of smaller intervals of length ϵ ,

$$t_0 = t_A, t_1 = t_A + \epsilon, t_2 = t_A + 2\epsilon, \dots, t_M = t_A + M\epsilon, t_{M+1} = t_B, (3.10)$$

where we have defined

$$\epsilon = \frac{T}{M+1}.\tag{3.11}$$

The time evolution operator decomposes into a product of operators $\hat{U}(t_{j+1}, t_j)$ over each of the smaller time intervals,

$$\hat{U}(t_B, t_A) = \prod_{j=0}^{M} \hat{U}(t_{j+1}, t_j).$$
(3.12)

Furthermore, we can use the completeness of the position eigenstates

$$\int d\vec{x} |\vec{x}\rangle \langle \vec{x}| = 1, \qquad (3.13)$$

⁹This is emphasized in some, but not all, textbooks. See, for instance, Sec. 2.2 of [48], Sec. 10.2 of [49], or Sec. 1.2 of [50].

where the integration measure $d\vec{x}$ is shorthand for $\prod_{i=1}^{N} dx^{i}$, to insert several resolutions of the identity and write

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A) = \int \left(\prod_{k=1}^M d\vec{x}_k\right) \left(\prod_{j=0}^M \langle \vec{x}_{j+1} | \hat{U}(t_{j+1}, t_j) | \vec{x}_j \rangle\right)$$
(3.14)

Here we have defined $\vec{x}_0 = \vec{x}_A$ and $\vec{x}_M = \vec{x}_B$. Note that the subscripts *j*, *k*, etc., on the position variables do not refer to the components x^i of the vector \vec{x} but rather to labels that index the different integration variables.

We can now focus on the propagator over one of the smaller time intervals of length ϵ . Over such an interval

from t to $t + \epsilon$, even if the Hamiltonian \hat{H} has explicit time dependence, we can approximate the unitary time evolution operator to $\mathcal{O}(\epsilon)$ as

$$\hat{U}(t+\epsilon,t) = e^{-i\hat{H}(t)\epsilon} + \mathcal{O}(\epsilon^2).$$
(3.15)

Inserting a complete set of momentum eigenstates using the completeness relation

$$\int d\vec{p} |\vec{p}\rangle \langle \vec{p}| = \frac{1}{(2\pi)^N}, \qquad (3.16)$$

we can then write a single factor in the integrand of (3.14) as

$$\langle \vec{x}_{j+1} | U(t+\epsilon,t) | \vec{x}_j \rangle = \int \frac{d\vec{p}}{(2\pi)^N} \langle \vec{x}_{j+1} | \hat{U}\left(t+\epsilon,t+\frac{\epsilon}{2}\right) | \vec{p} \rangle \langle \vec{p} | \hat{U}\left(t+\frac{\epsilon}{2},t\right) | \vec{x}_j \rangle.$$
(3.17)

It is convenient that we have two representations of the Hamiltonian (3.6) and (3.7), one with position operators to the right and one with momentum operators to the right. We evaluate the second matrix element in (3.17) to order ϵ using the normal-ordered form,

$$\langle \vec{p} | \hat{U} \left(t + \frac{\epsilon}{2}, t \right) | \vec{x}_j \rangle = \left\langle \vec{p} | 1 - \frac{i\epsilon}{2} \hat{H}(t) | \vec{x}_j \right\rangle + \mathcal{O}(\epsilon^2)$$

$$= \left\langle \vec{p} | \vec{x}_j \right\rangle - \frac{i\epsilon}{2} \left\langle \vec{p} | \sum h_{ikmn}(t) (\hat{p}^i)^m (\hat{x}^k)^n | \vec{x}_j \right\rangle + \mathcal{O}(\epsilon^2)$$

$$= e^{-i\vec{p}\cdot\vec{x}_j} - \frac{i\epsilon}{2} h(t; \vec{x}_j, \vec{p}) + \mathcal{O}(\epsilon^2),$$

$$(3.18)$$

where we use the symbol $h(\vec{x}_j, \vec{p})$ with no indices to refer to the normal-ordered Hamiltonian (3.6) with all operators replaced with classical variables.

Similarly, we use Hermiticity of \hat{H} along with the anti-normal-ordered form (3.7) for the Hamiltonian to evaluate the first matrix element appearing in (3.17),

$$\langle \vec{x}_{j+1} | \hat{U} \left(t + \epsilon, t + \frac{\epsilon}{2} \right) | \vec{p} \rangle = \left\langle \vec{p} | 1 - \frac{i\epsilon}{2} \hat{H}^{\dagger} \left(t + \frac{\epsilon}{2} \right) | \vec{x}_j \right\rangle + \mathcal{O}(\epsilon^2)$$

$$= \left\langle \vec{x}_{j+1} | \vec{p} \right\rangle - \frac{i\epsilon}{2} \left\langle \vec{x}_{j+1} | \sum h_{ikmn}^* \left(t + \frac{\epsilon}{2} \right) (\hat{x}^k)^n (\hat{p}^i)^m | \vec{p} \right\rangle + \mathcal{O}(\epsilon^2)$$

$$= e^{i\vec{x}_{j+1} \cdot \vec{p}} - \frac{i\epsilon}{2} h^* \left(t + \frac{\epsilon}{2} ; \vec{x}_{j+1}, \vec{p} \right) + \mathcal{O}(\epsilon^2), \qquad (3.19)$$

where we similarly write $h^*(t; \vec{x}_{j+1}, \vec{p})$ for the anti-normal-ordered Hamiltonian (3.6) with operators replaced by classical variables.

Using these results (3.18) and (3.19) for the matrix elements, we find

$$\begin{aligned} \langle \vec{x}_{j+1} | U(t+\epsilon,t) | \vec{x}_j \rangle &= \int \frac{d\vec{p}}{(2\pi)^N} \left(e^{i\vec{x}_{j+1}\cdot\vec{p}} - \frac{i\epsilon}{2} h^* \left(t + \frac{\epsilon}{2}; \vec{x}_{j+1}, \vec{p} \right) \right) \left(e^{-i\vec{p}\cdot\vec{x}_j} - \frac{i\epsilon}{2} h(t; \vec{x}_j, \vec{p}) \right) + \mathcal{O}(\epsilon^2) \\ &= \int \frac{d\vec{p}}{(2\pi)^N} e^{i(\vec{x}_{j+1}-\vec{x}_j)\cdot\vec{p}} \exp\left[-\frac{i\epsilon}{2} \left(h \left(t + \frac{\epsilon}{2}; \vec{x}_j, \vec{p} \right) + h^* \left(t + \frac{\epsilon}{2}; \vec{x}_{j+1}, \vec{p} \right) \right) \right] + \mathcal{O}(\epsilon^2). \end{aligned}$$
(3.20)

This discretization suggests that we should define the classical Hamiltonian

$$H(t; \vec{x}, \vec{p}) = \frac{1}{2} (h(t; \vec{x}, \vec{p}) + h^*(t; \vec{x}, \vec{p})) = \operatorname{Re}(h(t; \vec{x}, \vec{p})).$$
(3.21)

To leading order at small ϵ , the propagator over a small time interval ϵ is therefore

$$\langle \vec{x}_{j+1} | U(t+\epsilon,t) | \vec{x}_j \rangle = \int \frac{d\vec{p}}{(2\pi)^N} \exp[i(\vec{p}_j \cdot (\vec{x}_{j+1} - \vec{x}_j) - \epsilon H(t; \vec{x}_j, \vec{p}_j))].$$
(3.22)

The full time-sliced propagator (3.14) is obtained from the product of the individual factors (3.22) in the limit as $\epsilon \to 0$ and $M \to \infty$,

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A) = \lim_{\epsilon \to 0} \int \left(\prod_{k=1}^M d\vec{x}_k \right) \int \left(\prod_{l=0}^M \frac{d\vec{p}_l}{(2\pi)^N} \right)$$
$$\times \exp\left[i \sum_{j=0}^M (\vec{p}_j \cdot (\vec{x}_{j+1} - \vec{x}_j) - \epsilon H(t; \vec{x}_j, \vec{p}_j)) \right].$$
(3.23)

In the limit of small ϵ , the sum in the argument of the exponential becomes an integral:

$$\lim_{\epsilon \to 0} \sum_{j=0}^{M} \epsilon \left(\vec{p}_j \cdot \frac{(\vec{x}_{j+1} - \vec{x}_j)}{\epsilon} - H(t; \vec{x}_j, \vec{p}_j) \right)$$
$$= \int_{t_A}^{t_B} dt \left(p^i(t) \dot{x}_i(t) - H(t; \vec{x}(t), \vec{p}(t)) \right). \tag{3.24}$$

Here we have passed from discrete collections of \vec{x}_j , \vec{p}_j to continuous trajectories $x^i(t)$, $p^i(t)$. We conclude that the propagator for a general analytic Hamiltonian takes the form

$$K(\vec{x}_{B}, t_{B}; \vec{x}_{A}, t_{A}) = \int_{\vec{x}(t_{A}) = \vec{x}_{A}}^{\vec{x}(t_{B}) = \vec{x}_{A}} \mathcal{D}\vec{x} \int_{\vec{p}(t_{A}) = \vec{p}_{A}}^{\vec{p}(t_{B}) = \vec{p}_{B}} \mathcal{D}\vec{p} \\ \times \exp\left(i \int_{t_{A}}^{t_{B}} dt \left(p^{i}(t) \dot{x}_{i}(t) - H(\vec{p}, \vec{x})\right)\right),$$
(3.25)

where the path integral measures $D\vec{x}$ and $D\vec{p}$ are defined as the limits of the products in (3.23). To respect causality, we set the propagator equal to (3.25) when $t_B > t_A$ and set $K(\vec{x}_B, t_B; \vec{x}_A, t_A) = 0$ for $t_B < t_A$. Similarly, by performing this phase space path integral in Euclidean time with periodic boundary conditions, one can obtain a phase space integral expression for the finite-temperature partition function.

Again, it is important to emphasize that the integral (3.25) runs over all phase space paths $(\vec{x}(t), \vec{p}(t))$. For a generic path, there is no relationship between the coordinates and momenta; in particular, it is not the case that p^i is constrained to be equal to the canonical momentum which is conjugate to x^i . In the special case where the Hamiltonian is quadratic in the momenta, for instance, if

$$H = \frac{p^{i} p_{i}}{2m} + V(x), \qquad (3.26)$$

then the path integral over momenta can be evaluated, leaving a path integral over positions $x^i(t)$. Only when performing this evaluation is the expression $p^i(t)$ set equal to the conjugate momentum, which reproduces the usual Feynman path integral (3.3) that involves the Lagrangian. For nonquadratic Hamiltonians, however, no such reduction is possible, and we must use the more fundamental form (3.25).

B. Deformations in canonical quantization and path integral quantization

We now wish to study how observables in quantum mechanics are modified when the theory is deformed by conserved charges. Our goal is to show that the propagator, and hence the finite-temperature partition function, of a general theory satisfies a flow equation which is identical to the one derived in Sec. II C for the classical partition function. As a consistency check, we will also see that this flow equation for the propagator can be equivalently derived using either canonical methods or path integral methods.

1. Canonical analysis

First, let us see how to understand this flow equation in the canonical formalism. For the moment, we will specialize to the case of Hamiltonians which do not depend on time explicitly. In this case, the unitary time evolution operator can be written as

$$\hat{U}(t_B, t_A) = e^{-i\hat{H}(t_B - t_A)},$$
 (3.27)

and the propagator is

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A) = \langle \vec{x}_B | e^{-i\hat{H}(t_B - t_A)} | \vec{x}_A \rangle.$$
(3.28)

Suppose that the one-parameter family of Hamiltonian operators \hat{H}_{λ} satisfies a flow equation driven by a combination of conserved charge operators \hat{Q}_a ,

$$\frac{\partial \hat{H}_{\lambda}}{\partial \lambda} = \hat{\mathfrak{D}}(\hat{Q}_a). \tag{3.29}$$

If the operator $\hat{\Sigma}$ depends on charges besides the Hamiltonian, it will not be possible to derive a closed flow equation for the usual propagator (3.28). Instead, we must consider a more general propagator that includes sources for the various conserved charges Q_a . We therefore define the quantity

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A; \lambda; \mu_a) = \langle \vec{x}_B | \exp\left(-i\hat{H}_{\lambda}(t_B - t_A) + \sum_a \mu_a \hat{Q}_a\right) | \vec{x}_A \rangle, \qquad (3.30)$$

where μ_a are a collection of couplings that serve the same purpose as the chemical potentials in the analysis of the classical partition function in Sec. II C.

For simplicity, we define $T = t_B - t_A$, and we will suppress the arguments of the propagator in what follows. One has the relations

$$\partial_{\lambda}K = \left\langle \vec{x}_{B} \right| - iT\hat{\mathfrak{O}}e^{-i\hat{H}_{\lambda}T + \mu_{a}\hat{Q}_{a}} |\vec{x}_{A} \right\rangle,$$

$$\partial_{T}K = \left\langle \vec{x}_{B} \right| - i\hat{H}e^{-i\hat{H}_{\lambda}T + \mu_{a}\hat{Q}_{a}} |\vec{x}_{A} \right\rangle,$$

$$\partial_{\mu_{b}}K = \left\langle \vec{x}_{B} \right|\hat{Q}_{b}e^{-i\hat{H}_{\lambda}T + \mu_{a}\hat{Q}_{a}} |\vec{x}_{A} \right\rangle.$$
(3.31)

Summation is implied in the expression $\mu_a \hat{Q}_a = \sum_a \mu_a \hat{Q}_a$. Because the deforming operator $\mathfrak{O}(\hat{H}, \hat{Q}_a)$ is itself a function of the operators \hat{H} and \hat{Q}_a , we arrive at a differential equation

$$\partial_{\lambda}K = -iT\mathfrak{O}[i\partial_T, \partial_{\mu_a}]K. \tag{3.32}$$

This is identical to the flow equation (2.61) for the classical partition function after identifying $\beta = iT$. As in Sec. II C, the expression $\mathfrak{D}[i\partial_T, \partial_{\mu_a}]$ is defined by expanding the operator \mathfrak{D} in a power series in \hat{H} and the \hat{Q}_a , and then replacing each instance of \hat{H} with $i\partial_T$ and each instance of \hat{Q}_a with ∂_{μ_a} .

We may also think of this flow equation in terms of a generalization of the prescription of [35,36] for quantizing theories that are deformed by functions of the Hamiltonian. To do this, let us first recall how to derive the kernel representation of the propagator. Suppose that we can identify a complete basis of simultaneous eigenstates of the Hamiltonian \hat{H}_{λ} and each of the charge operators \hat{Q}_a . We write these simultaneous eigenstates as $|\phi_n\rangle$, which satisfy

$$\hat{H}_{\lambda}|\phi_{n}\rangle = E_{n}(\lambda)|\phi_{n}\rangle, \qquad \hat{Q}_{a}|\phi_{n}\rangle = q_{a,n}|\phi_{n}\rangle.$$
 (3.33)

In this case, we are thinking of the label n as a multi-index, which collects the quantum numbers for all of the charges. Using the completeness relation

$$\sum_{n} |\phi_n\rangle \langle \phi_n| = 1, \qquad (3.34)$$

we can evaluate the propagator as

$$K = \left\langle \vec{x}_{B} | e^{-i\hat{H}_{\lambda}T + \mu_{a}\hat{Q}_{a}} | \vec{x}_{A} \right\rangle$$

$$= \sum_{n,m} \left\langle \vec{x}_{B} | \phi_{n} \right\rangle \left\langle \phi_{n} | e^{-i\hat{H}_{\lambda}T + \mu_{a}\hat{Q}_{a}} | \phi_{m} \right\rangle \left\langle \phi_{m} | \vec{x}_{A} \right\rangle$$

$$= \sum_{n,m} e^{-iE_{n}(\lambda)T + \mu_{a}q_{n,a}} \left\langle \vec{x}_{B} | \phi_{n} \right\rangle \delta_{m,n} \left\langle \phi_{m} | \vec{x}_{A} \right\rangle$$

$$= \sum_{n} \phi_{n}^{*}(\vec{x}_{A}) \phi_{n}(\vec{x}_{B}) e^{-iE_{n}(\lambda)T + \mu_{a}q_{n,a}}.$$
(3.35)

Here we have used the orthogonality relation $\langle \phi_n | \phi_m \rangle = \delta_{m,n}$, along with the definition of the position space wave function

$$\langle \vec{x} | \phi_n \rangle = \phi_n(\vec{x}). \tag{3.36}$$

Equation (3.35) suggests a straightforward interpretation of a deformation by conserved charges. In the deformed theory, every eigenstate $|\phi_n\rangle$ of the Hamiltonian and charges remains an eigenstate of the Hamiltonian and charges. The eigenvalues of $|\phi_n\rangle$ under each of the charge operators remains unchanged, since we deform the Hamiltonian \hat{H} but not the operators \hat{Q}_a . It is easy to see that this is consistent with the fact that all of the operators \hat{Q}_a remain conserved in the deformed theory, since

$$[\hat{\mathfrak{D}}, \hat{Q}_a] = 0, \qquad (3.37)$$

by virtue of the fact that we assume all of the charges \hat{Q}_a are commuting (which is the quantum version of our assumption (2.41) that the charges are classically Poisson commuting). All that has changed is that each ket $|\phi_n\rangle$ now has a deformed energy eigenvalue $E_n(\lambda)$ that obeys the differential equation

$$\partial_{\lambda} E_n(\lambda) = \mathfrak{O}(E_n, q_{n,a}), \qquad (3.38)$$

where on the right side of Eq. (3.38), \mathfrak{D} is now a classical variable that is evaluated on the energy eigenvalue E_n and charge eigenvalues $q_{n,a}$ of the state $|\phi_n\rangle$.

For instance, in the case of the deformation by $\mathfrak{D} = \frac{H_{\lambda}^2}{\frac{1}{2}-2\lambda H_{\lambda}}$ whose classical solution was discussed around (2.46), each energy eigenstate $|\phi_n\rangle$ with undeformed energy $E_n(0)$ remains an energy eigenstate in the deformed theory, but with a new energy

$$E_n(\lambda) = \frac{1}{4\lambda} \left(1 - \sqrt{1 - 8\lambda E_n(0)} \right).$$
(3.39)

There is a sharp difference in the behavior of the deformed spectrum depending on the sign of λ . If $\lambda < 0$, then the argument of the square root remains positive for arbitrary large positive undeformed energies. This choice is called the "good sign" of the deformation parameter. However, if $\lambda > 0$, then for sufficiently large undeformed energies $E_n(0)$, the deformed energy levels become complex. This is the "bad sign" of the deformation. The same qualitative behavior occurs for the $T\bar{T}$ deformation of a 2*d* CFT.¹⁰

2. Path integral analysis

We will now see how the above flow equations can be derived using the path integral. For the same reasons as we mentioned above, if the deforming operator \mathfrak{D} depends on charges Q_a besides the Hamiltonian, we cannot obtain a flow equation for the propagator using the unflavored path integral (3.4). Instead, we must introduce a flavored version that includes sources for the various charge operators,

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A; \lambda; \mu_a) = \int_{\vec{x}(t_A) = \vec{x}_A}^{\vec{x}(t_B) = \vec{x}_B} \mathcal{D}\vec{x} \int_{\vec{p}(t_A) = \vec{p}_A}^{\vec{p}(t_B) = \vec{p}_A} \mathcal{D}\vec{p}$$
$$\times \exp\left(i \int_{t_A}^{t_B} dt \left(p^i(t) \dot{x}_i(t) - H_\lambda(\vec{p}, \vec{x})\right) + \sum_a \mu_a Q_a\right). \quad (3.40)$$

To ease notation, we will again suppress the arguments of the propagator, omit the upper and lower bounds of path integration [which are always understood to take the values in (3.40)], and write $\mu_a Q_a$ for $\sum_a \mu_a Q_a$.

Let us emphasize two points about the phase space path integral expression for the propagator. First, as is typical of path integrals, all dynamical quantities appearing inside the integrand are simply classical variables $x^i(t)$ and $p^i(t)$ rather than quantum operators \hat{x}^i and \hat{p}^i , so there are no ordering ambiguities. Second, and more importantly, it is critical that the momentum variables $p^i(t)$ inside the path integral are *not* the conjugate momenta to $x^i(t)$. The path integral runs over all phase space trajectories with the specified end points, and there is no constraint between the functions x(t) and p(t) along these trajectories. This is important because it implies that

$$\frac{\partial p^{i}(t)}{\partial \lambda} = \frac{\partial \dot{x}^{i}(t)}{\partial \lambda} = 0.$$
(3.41)

Let us contrast this situation with that of the proof of Theorem 1. In that context, the momenta and velocities PHYS. REV. D 108, 126021 (2023)

were related by Eq. (2.18), the Hamilton equation of motion:

$$\frac{\partial H_{\lambda}}{\partial p^{i}} = \dot{x}^{i}. \tag{3.42}$$

Therefore, if we choose to treat p^i as an independent variable which does not depend on λ , it follows that

$$\frac{\partial \dot{x}^{i}}{\partial \lambda} = \frac{\partial (\partial_{\lambda} H_{\lambda})}{\partial p^{i}}, \qquad (3.43)$$

which is nonzero in general. This additional term appeared in (2.22), where it was needed to demonstrate the equivalence of deformations of the Lagrangian and of the Hamiltonian.

However, since the p^i and x^i are unrelated integration variables in (3.40), no term of the form (3.43) is generated when we differentiate the propagator with respect to λ . The only λ dependence appears in the Hamiltonian itself, so one finds

$$\partial_{\lambda}K = \int \mathcal{D}\vec{x} \int \mathcal{D}\vec{p} \left(-i \int_{t_{A}}^{t_{B}} dt\mathfrak{D}\right) \\ \times \exp\left(i \int_{t_{A}}^{t_{B}} dt(p^{i}(t)\dot{x}_{i}(t) - H_{\lambda}(\vec{p},\vec{x})) + \mu_{a}Q_{a}\right).$$
(3.44)

We now use the fact that \mathfrak{D} is only a function of conserved charges, which are independent of time. This means that the path integral expectation value of \mathfrak{D} is itself also independent of time, so we may interchange the time integral with the path integral to conclude

$$\partial_{\lambda}K = -i(t_B - t_A)\langle \mathfrak{O} \rangle, \qquad (3.45)$$

where we have defined the path integral expectation value

$$\langle f(x^{j}, p^{j}) \rangle = \int \mathcal{D}\vec{x} \int \mathcal{D}\vec{p}f(x^{j}, p^{j}) \\ \times \exp\left(i \int_{t_{A}}^{t_{B}} dt(p^{i}(t)\dot{x}_{i}(t) - H_{\lambda}(\vec{p}, \vec{x})) + \mu_{a}Q_{a}\right).$$
(3.46)

Likewise, defining $T = t_B - t_A$, one has

$$\partial_T K = \langle i(p^i(t)\dot{x}_i(t) - H_\lambda(p, x)) \rangle$$

= $-i\langle H_\lambda \rangle.$ (3.47)

Here we have used that the quantity $p(t)\dot{x}(t)$ is odd in both $p^{i}(t)$ and $x^{i}(t)$, and a path integral of an odd quantity over all paths vanishes by symmetry. Therefore, the first term in the path integral expectation value of the first line of (3.47)

¹⁰In this setting, the complex spectrum can sometimes be returned to a purely real spectrum by performing sequential $T\bar{T}$ deformations with a sufficiently large positive flow parameter [51].

vanishes; the second term, H_{λ} , is constant in time because the Hamiltonian is conserved. Alternatively, one can justify the conclusion (3.47) using the Schrödinger equation, which relates the time derivative of the propagator to the Hamiltonian. This Schrödinger relation for the phase space path integral is identical to that of the familiar Feynman path integral.

Similarly, the derivatives of the propagator with respect to the chemical potentials μ_a generate expectation values of charges:

$$\partial_{\mu_a} K = \langle Q_a \rangle. \tag{3.48}$$

Note that we have chosen conventions for the flavored phase space path integral (3.40) such that the terms $\mu_a Q_a$ do not appear under the time integral. This leads to flow equations that are most similar to the classical results obtained for the grand canonical partition function (2.57). However, because all of the charges Q_a are conserved in time, we could have alternatively defined the path integral so that the terms $\mu_a Q_a$ were instead included inside of the time integral. This would have introduced an additional factor of T in (3.48), and dependence on the charges in the computation of $\partial_T K$, which would produce a different flow equation that is related to our result by a redefinition of parameters.

We conclude that the expressions for $\partial_{\lambda}K$, $\partial_{T}K$, and $\partial_{\mu_{a}}K$ computed using the path integral formulation are identical to those computed using the canonical analysis. The phase space path integral representation for the propagator therefore obeys the same differential equation,

$$\partial_{\lambda}K = -iT\mathfrak{D}[i\partial_T, \partial_{\mu_a}]K, \qquad (3.49)$$

where the right side is defined for any analytic function \mathfrak{D} of conserved charges.

3. Comments on nonanalytic deformations

Our derivation of the phase space path integral in Sec. III A assumed that the Hamiltonian operator \hat{H} admits a power series expansion in the operators \hat{x}^i and \hat{p}^i . We also derived differential equations obeyed by the propagator, in either the canonical or path integral formalism, which involve expressions $\mathfrak{D}[i\partial_T, \partial_{\mu_a}]$ that are defined by series expanding the deforming operator \mathfrak{D} and replacing instances of H and Q_a with various derivatives. We now consider cases in which the deformation is not an analytic function of the charges, which includes the case of the 1*d* root- $T\bar{T}$ operator (2.52).

For a totally general nonanalytic Hamiltonian \hat{H} , it is not clear how to perform an analog of the time-slicing prescription of Sec. III A and obtain a path integral definition. However, for a first-order deformation of an analytic Hamiltonian by a nonanalytic function of conserved charges, the arguments of the preceding subsections apply in almost exactly the same way. At the risk of repeating ourselves, let us quickly check that this is true. We will consider a Hamiltonian with a form that is slightly more general than a first-order deformation, namely

$$\hat{H} = f_1(\lambda)\hat{H}_0 + f_2(\lambda)\hat{\mathfrak{O}}, \qquad (3.50)$$

where \hat{H}_0 is analytic and $\hat{\mathfrak{O}}$ is a (possibly nonanalytic) function of charges. When $f_1 = 1$ and $f_2 = \lambda$, this is the leading-order correction to \hat{H}_0 generated by a flow driven by \mathfrak{O} .

As before, using the kernel representation of the propagator

$$K = \sum_{n} \phi_{n}^{*}(x_{A})\phi_{n}(x_{B})e^{-iE_{n}T + \mu_{a}q_{n,a}}, \qquad (3.51)$$

one would similarly argue that a deformation of \hat{H} by a nonanalytic function \mathfrak{D} of conserved charges has the effect of leaving all of the eigenfunctions ϕ_n unchanged and only shifts the energy eigenvalues as $E_n(\lambda) = f_1(\lambda)E_n(0) + f_2(\lambda)\mathfrak{D}$. One can still differentiate

$$\partial_{\lambda}K = \sum_{n} \phi_{n}^{*}(x_{A})\phi_{n}(x_{B})(-iT(f_{1}'(\lambda)E_{n}(0) + f_{2}'(\lambda)\mathfrak{D}))e^{-iE_{n}(\lambda)T + \mu_{a}q_{n,a}},$$
(3.52)

although it may no longer be possible to express (3.52) in terms of derivatives of *K* with respect to *T* and the μ_a when \mathfrak{D} is nonanalytic.

A similar analysis is possible using the path integral. By the Baker-Campbell-Hausdorff (BCH) formula, one has

$$\exp\left(f_1(\lambda)\hat{H}_0 + f_2(\lambda)\hat{\mathfrak{O}}\right) = \exp\left(f_1(\lambda)\hat{H}_0\right)\exp\left(f_2(\lambda)\hat{\mathfrak{O}}\right),$$
(3.53)

since all commutator terms in the BCH expansion vanish by virtue of the fact that \mathfrak{D} is a function only of conserved quantities. Suppose that we repeat the time-slicing prescription of Sec. III A for this Hamiltonian. When evaluating (3.14), one has

$$\begin{aligned} \langle \vec{x}_{j+1} | \hat{U}(t_{j+1}, t_j) | \vec{x}_j \rangle &= \langle \vec{x}_{j+1} | e^{-if_1(\lambda)H_0\epsilon} e^{-if_2(\lambda)\mathfrak{D}\epsilon} | \vec{x}_j \rangle \\ &= \sum_{n,m} \langle \vec{x}_{j+1} | e^{-if_1(\lambda)\hat{H}_0\epsilon} | \phi_n \rangle \\ &\times \langle \phi_n | e^{-if_2(\lambda)\hat{\mathfrak{D}}\epsilon} | \phi_m \rangle \langle \phi_m | \vec{x}_j \rangle, \quad (3.54) \end{aligned}$$

where we have inserted two complete sets of eigenstates $|\phi_n\rangle$ of both the undeformed Hamiltonian \hat{H}_0 and the

charge operators \hat{Q}_a . We can then evaluate the middle factor by replacing $\hat{\mathfrak{D}}$ with its classical value \mathfrak{D} , giving

$$\langle \vec{x}_{j+1} | \hat{U}(t_{j+1}, t_j) | \vec{x}_j \rangle = \sum_{n,m} e^{-if_2(\lambda)\mathfrak{D}e} \langle \vec{x}_{j+1} | e^{-if_1(\lambda)\hat{H}_0e} | \phi_n \rangle \langle \phi_n | \phi_m \rangle \langle \phi_m | \vec{x}_j \rangle$$

$$= e^{-if_2(\lambda)\mathfrak{D}e} \langle \vec{x}_{j+1} | e^{-if_1(\lambda)\hat{H}_0e} | \vec{x}_j \rangle,$$

$$(3.55)$$

and then we may evaluate the remaining matrix element for the analytic part $f_1(\lambda)\hat{H}_0$ of the Hamiltonian using the same steps as before. This would lead us to the same result,

$$\partial_{\lambda}K = \left\langle -iT(f_1'(\lambda)\hat{H}_0 + f_2'(\lambda)\hat{\mathfrak{D}}) \right\rangle.$$
(3.56)

Taking $f_1 = 1$ and $f_2 = \lambda$ allows us to conclude, using either formalism, that a deformation of an analytic Hamiltonian by a nonanalytic function of charges—to leading order in the deformation parameter—is described by the same differential equation $\partial_{\lambda}K = \langle -iT\hat{\Omega} \rangle$ for the propagator. We emphasize that the only difference in the case of a nonanalytic deforming operator is that the expectation value $\langle \hat{\Omega} \rangle$ may not be expressible in terms of derivatives of the propagator with respect to *T* and the μ_a .

One might then ask how one can extend this analysis to higher orders in the deformation parameter. For instance, suppose that we use the analysis above to define the Hamiltonian $\hat{H}^{(1)} = \hat{H}_0 + \lambda \hat{\Omega}$ which solves the flow equation $\partial_{\lambda}\hat{H}_{\lambda} = \hat{\Omega}$ to first order, and we then wish to treat $\hat{H}^{(1)}$ as a new seed Hamiltonian to deform again and generate the second-order solution $\hat{H}^{(2)}$, and continue in this way to define all higher $\hat{H}^{(n)}$. Let us make two comments on this point:

(I) The first comment is that a general, first-principles analysis of this process would require a procedure for obtaining a path integral representation for the propagator of an arbitrary nonanalytic seed Hamiltonian, which is not available. Instead, we can simply *define* what we mean by the all-orders version of the deformed quantum theory by first solving the differential equation $\partial_{\lambda}H_{\lambda} = \mathfrak{D}$, and then inserting this solution for H_{λ} into the phase space path integral (3.25).

By construction, this is equivalent to using the corresponding solution for the deformed energy levels $E_n(\lambda)$ and inserting them into the kernel representation (3.51) of the propagator. In this way, we obtain a consistent prescription for the quantization of a nonanalytic theory that agrees with the above analysis to leading order around an analytic seed Hamiltonian.

(II) The second comment is that, in some cases of interest, we can sidestep the issue of performing path integral quantization of a nonanalytic seed Hamiltonian. Specifically, for the deformation of the harmonic oscillator by the 1d root- $T\bar{T}$ operator to obtain the ModMax oscillator, it turns out that the *all-orders* solution to the flow equation takes the form (3.50) for appropriately chosen functions f_1 and f_2 . In this case, the analysis that we carried out for the leading-order deformation is sufficient to derive differential equations that hold to all orders along the flow.

C. Flow of quantum partition function and comparison to 2d

In the preceding subsection, we have developed general differential equations obeyed by the propagator $K(\vec{x}_B, t_B; \vec{x}_A, t_A)$ of a quantum theory deformed by a function of conserved charges. Since the Euclidean time propagator with periodic boundary conditions is the thermal partition function, which can also be written in the trace form

$$Z(\beta) = \operatorname{Tr}(e^{-\beta\hat{H}}), \qquad (3.57)$$

one can likewise obtain differential equations obeyed by $Z(\beta)$, or more precisely for the grand canonical partition function with chemical potentials for the various charges:

$$\mathcal{Z}(\beta,\mu_1,\ldots,\mu_M) = \operatorname{Tr}(e^{-\beta\hat{H} + \sum \mu_a \hat{Q}_a}). \quad (3.58)$$

One can derive flow equations for (3.58) using manipulations that are identical to those around Eq. (2.61). In particular,

$$\partial_{\beta} \mathcal{Z} = \operatorname{Tr}(-\hat{H}e^{-\beta\hat{H} + \sum \mu_{a}\hat{Q}_{a}}) = -\langle \hat{H} \rangle,$$

$$\partial_{\mu_{a}} \mathcal{Z} = \operatorname{Tr}(\hat{Q}_{a}e^{-\beta\hat{H} + \sum \mu_{b}\hat{Q}_{b}}) = \langle \hat{Q}_{a} \rangle, \qquad (3.59)$$

and thus

$$\partial_{\lambda} \mathcal{Z} = -\beta \mathfrak{O}[-\partial_{\beta}, \partial_{\mu_1}, ..., \partial_{\mu_M}](\mathcal{Z}).$$
(3.60)

In fact, we could have derived the flow equation for the classical partition function by first arguing that the quantum partition function satisfies the differential equation (3.60) and then taking the limit $\hbar \rightarrow 0$.

We will discuss some examples of such flow equations only briefly, because they take the same form as the corresponding flow equations for the classical partition function described in Sec. II C. However, since these differential equations now hold in the quantum theory, we will comment on the relationship to the analogous flow equations for the quantum-mechanical partition functions of two-dimensional field theories which are deformed by the $T\bar{T}$ operator.

For instance, under the flow (2.33) which is the 1*d* version of the $T\bar{T}$ deformation, the propagator *K* and thermal partition function *Z* obey the differential equations

$$\begin{pmatrix} 4\lambda\partial_{\lambda}\partial_{T} + 2T\partial_{T}^{2} + i\left(1 + \frac{4i\lambda}{T}\right)\partial_{\lambda} \end{pmatrix} K_{\lambda}(T) = 0, \\ \left(4\lambda\partial_{\lambda}\partial_{\beta} + 2\beta\partial_{\beta}^{2} + \left(1 - \frac{4\lambda}{\beta}\right)\partial_{\lambda} \right) Z_{\lambda}(\beta) = 0.$$
 (3.61)

This flow equation for Z was also considered in [46]. It can be understood as follows. Suppose that we begin with a two-dimensional conformal field theory whose torus partition function is $Z_0(\tau, \bar{\tau})$, where τ is the modular parameter of the torus. One can then deform this theory by the $T\bar{T}$ operator to obtain a one-parameter family of deformed torus partition functions Z_{λ} which obey the differential equation

$$\partial_{\lambda} Z_{\lambda}(\tau, \bar{\tau}) = \left(\tau_2 \partial_{\tau} \partial_{\bar{\tau}} + \frac{1}{2} \left(\partial_{\tau_2} - \frac{1}{\tau_2}\right) \lambda \partial_{\lambda} \right) Z_{\lambda}(\tau, \bar{\tau}). \quad (3.62)$$

This differential equation can be obtained by performing a Hubbard-Stratonovich transformation which replaces the $T\bar{T}$ deformation with a random metric [11]. Modular properties of the deformed partition function were studied in [12,13]; although the $T\bar{T}$ -deformed theory no longer enjoys conformal symmetry, the partition function Z_{λ} is still invariant with respect to a modular transformation under which the parameter λ also transforms.

Let us specialize to a torus with purely imaginary modular parameter $\tau = \frac{i\beta}{8}$. Using

$$\partial_{\tau} = \frac{1}{2} (\partial_{\tau_1} - i \partial_{\tau_2}), \qquad \partial_{\bar{\tau}} = \frac{1}{2} (\partial_{\tau_1} + i \partial_{\tau_2}), \qquad (3.63)$$

with $\tau_2 = \frac{\beta}{8}$ and $\tau_1 = 0$, the flow equation (3.62) reduces to

$$\partial_{\lambda} Z_{\lambda}(\beta) = \left(2\beta \partial_{\beta}^{2} + 4\left(\partial_{\beta} - \frac{1}{\beta}\right) \lambda \partial_{\lambda} \right) Z_{\lambda}(\beta), \qquad (3.64)$$

which reproduces (3.61) after sending $\lambda \to -\lambda$, which is a choice of conventions for the deformation parameter. Thus the 1*d* $T\bar{T}$ -deformed partition function indeed descends from the 2*d* $T\bar{T}$ -deformed partition function.

It is interesting to consider a similar dimensional reduction of the two-dimensional root- $T\bar{T}$ flow. The analogous flow equation for the torus partition function of a root- $T\bar{T}$ deformed 2*d* CFT was conjectured in [28] to be

$$\partial_{\gamma}^2 Z_{\gamma}(\tau,\bar{\tau}) = (\tau_2^2 \partial_{\tau} \partial_{\bar{\tau}} + \tau_2 \partial_{\tau_2}) Z_{\gamma}(\tau,\bar{\tau}), \qquad (3.65)$$

based on a proposal for the flow of the cylinder spectrum that was justified by holographic considerations. If we take a torus with modular parameter

$$\tau = i\beta + \mu, \tag{3.66}$$

then the flow equation (3.65) reduces to

$$\left(\partial_{\gamma}^{2} - \beta \partial_{\beta} - \frac{\beta^{2}}{4}(\partial_{\beta}^{2} + \partial_{\mu}^{2})\right) \mathcal{Z}(\beta, \gamma, \mu) = 0.$$
(3.67)

The factor of $\frac{1}{4}$ multiplying the third term in (3.67) is related to the normalization of the root- $T\bar{T}$ operator and can be rescaled to 1 by an appropriate redefinition. Up to this choice of scaling, this is the same differential equation that is satisfied by the partition function of the ModMax oscillator, which we will present in Eq. (4.24).

IV. APPLICATION TO MODMAX OSCILLATOR

In this section, we will apply the general results of the preceding sections to the main example of interest in the present work, which is the ModMax oscillator. This theory was first introduced in [30] and is a particular deformation of an isotropic harmonic oscillator. Given a collection of position variables x^i , for i = 1, ..., N, we begin by defining the undeformed theory with the Lagrangian

$$L_0 = \frac{1}{2} (\dot{x}^i \dot{x}^i - x^i x^i), \qquad (4.1)$$

where we have set the mass m and frequency ω of the harmonic oscillator to 1 for convenience. This theory has a conserved energy

$$E_0 = \frac{\partial L_0}{\partial \dot{x}^i} \dot{x}^i - L_0 = \frac{1}{2} (\dot{x}^i \dot{x}^i + x^i x^i), \qquad (4.2)$$

which is the Noether current associated with time translation symmetry, along with a collection of conserved angular momenta

$$J_0^{nm} = \frac{\partial L_0}{\partial \dot{x}^n} x^m - \frac{\partial L_0}{\partial \dot{x}^m} x^n = \dot{x}^n x^m - \dot{x}^m x^n, \quad (4.3)$$

which are the conserved currents associated with rotations $x^i \rightarrow R^i_{\ i} x^j$, $R \in SO(N)$. The total angular momentum is

$$J_0^2 = J_0^{nm} J_{0,nm}.$$
 (4.4)

For any γ , we can now define the Lagrangian for the ModMax oscillator as

$$L_{\gamma} = \frac{\cosh(\gamma)}{2} (\dot{x}^{i} \dot{x}^{i} - x^{i} x^{i}) \pm \sinh(\gamma) \sqrt{E_{0}^{2} - J_{0}^{2}}.$$
 (4.5)

Note that there is a choice of the relative sign between the two terms in (4.5), which is correlated with the choice of sign for the root- $T\bar{T}$ operator that drives the flow equation (4.7) below. One can also view this sign choice as a convention for the sign of the parameter γ , since sending $\gamma \rightarrow -\gamma$ reverses the relative sign.

Even though this Lagrangian L_{γ} is written in terms of the conserved quantities E_0 and J_0^2 in the *undeformed* theory, it satisfies a differential equation that involves the conserved currents in the *deformed* theory at finite γ , namely

$$E_{\gamma} = \frac{\partial L_{\gamma}}{\partial \dot{x}^{i}} \dot{x}^{i} - L_{\gamma}, \quad J_{\gamma}^{nm} = \frac{\partial L_{\gamma}}{\partial \dot{x}^{n}} x^{m} - \frac{\partial L_{\gamma}}{\partial \dot{x}^{m}} x^{n}, \quad J_{\gamma}^{2} = J_{\gamma}^{nm} J_{\gamma,nm}.$$

$$(4.6)$$

One can show that L_{γ} obeys

$$\frac{\partial L_{\gamma}}{\partial \gamma} = \pm \sqrt{E_{\gamma}^2 - J_{\gamma}^2}, \qquad (4.7)$$

which we refer to as the 1d root- $T\bar{T}$ flow equation, and which was first written down in [30]. The corresponding Hamiltonian,

$$H_{\gamma} = \frac{\cosh(\gamma)}{2} (p^{i} p^{i} + x^{i} x^{i}) \mp \sinh(\gamma) \sqrt{E_{0}^{2} - J_{0}^{2}}, \quad (4.8)$$

satisfies a flow equation with the opposite sign,

$$\frac{\partial H_{\gamma}}{\partial \gamma} = \mp \sqrt{E_{\gamma}^2 - J_{\gamma}^2}, \qquad (4.9)$$

as required by Theorem 1. In Eqs. (4.8) and (4.9), the quantities E_0 , E_γ , J_0^2 , and J_γ^2 are defined by beginning with the appropriate Noether currents computed in the Lagrangian formulation, and then expressing these quantities in terms of conjugate momenta p^i rather than velocities \dot{x}^i .

Although the ModMax oscillator can be defined for any number N of position variables x^i , i = 1, ..., N, we will focus on the case N = 2 for simplicity. First, let us review some features of the classical dynamics of the ModMax oscillator.

A. Classical aspects

We now specialize to the case of N = 2 coordinates x^i , and we will use the notation $x^1 = x$, $x^2 = y$. In terms of these variables, the general Lagrangian (4.5) for the ModMax oscillator can be written as

$$L_{\gamma} = \frac{1}{2} \left[\cosh(\gamma)(\dot{x}^2 + \dot{y}^2 - x^2 - y^2) \pm \sinh(\gamma) \sqrt{((y + \dot{x})^2 + (x - \dot{y})^2)((x + \dot{y})^2 + (y - \dot{x})^2)} \right].$$
(4.10)

The conserved angular momentum, which is the Noether charge associated with rotations in the (x, y) plane, is

$$J_{\gamma} = (x\dot{y} - y\dot{x})\cosh(\gamma) \pm \sinh(\gamma) \cdot \frac{(x\dot{y} - y\dot{x})(\dot{x}^2 + \dot{y}^2 - x^2 - y^2)}{\sqrt{((y + \dot{x})^2 + (x - \dot{y})^2)((x + \dot{y})^2 + (y - \dot{x})^2)}}.$$
(4.11)

It is interesting to express J_{γ} in terms of the conjugate momenta p_x and p_y , as appropriate for formulating flows for the Hamiltonian. The conjugate momenta computed from the Lagrangian (4.10) are

$$p_{x} = \frac{\partial L_{\gamma}}{\partial \dot{x}} = \dot{x} \cosh(\gamma) \pm \sinh(\gamma) \cdot \frac{2xy\dot{y} + x^{2}\dot{x} + \dot{x}(\dot{x}^{2} + \dot{y}^{2} - y^{2})}{\sqrt{((y + \dot{x})^{2} + (x - \dot{y})^{2})((x + \dot{y})^{2} + (y - \dot{x})^{2})}},$$

$$p_{y} = \frac{\partial L_{\gamma}}{\partial \dot{y}} = \dot{y} \cosh(\gamma) \pm \sinh(\gamma) \cdot \frac{2xy\dot{x} - x^{2}\dot{y} + \dot{y}(\dot{x}^{2} + \dot{y}^{2} + y^{2})}{\sqrt{((y + \dot{x})^{2} + (x - \dot{y})^{2})((x + \dot{y})^{2} + (y - \dot{x})^{2})}}.$$
(4.12)

After expressing the angular momentum J_{γ} in terms of p_x and p_y , one finds

$$J_{\gamma} = xp_{\gamma} - yp_{x}. \tag{4.13}$$

That is, when written in Hamiltonian variables, the deformed angular momentum J_{γ} takes the same functional form as the undeformed angular momentum J_0 . This is a special case of the observation, which we first made in the text below Eq. (3.36), that $f(H, Q_a)$ deformations modify the Hamiltonian but not the other charges such as J.

Similarly, for N = 2 one can write the Hamiltonian for the ModMax oscillator as

$$H_{\gamma} = \frac{1}{2} \left[\cosh(\gamma) (p_x^2 + p_y^2 + x^2 + y^2) \mp \sinh(\gamma) \sqrt{((p_y + x)^2 + (p_x - y)^2)((p_y - x)^2 + (p_x + y)^2)} \right].$$
(4.14)

Let us consider the symmetries of this theory in somewhat more detail. It is well-known that the undeformed theory, which is the 2*d* isotropic harmonic oscillator, enjoys an SU(2) symmetry. To see this, it is convenient to define complex variables

$$z = x + ip_x, \qquad w = y + ip_y, \tag{4.15}$$

so that the harmonic oscillator Hamiltonian is

$$H_0 = \frac{1}{2}(|z|^2 + |w|^2). \tag{4.16}$$

This Hamiltonian is invariant under any action of the form

$$\begin{bmatrix} z \\ w \end{bmatrix} \to U \begin{bmatrix} z \\ w \end{bmatrix}, \qquad U \in SU(2), \qquad (4.17)$$

since such an SU(2) transformation preserves the length of the complex vector. In these complex variables, the angular momentum is

$$J = \frac{1}{2i} (w\bar{z} - z\bar{w}) = \text{Im}(w\bar{z}).$$
(4.18)

The angular momentum J is *not* invariant under the full SU(2) symmetry group. However, it is still invariant under the restricted U(1) transformations

$$z \to e^{i\alpha}z, \qquad w \to e^{i\alpha}w, \qquad \alpha \in \mathbb{R}.$$
 (4.19)

Similarly, any deformed Hamiltonian that is a function of both the undeformed Hamiltonian and this angular momentum,

$$H = H(H_0, J), (4.20)$$

is also invariant under the U(1) transformations (4.19).

We have commented before that the ModMax oscillator is a particular dimensional reduction of the four-dimensional ModMax theory, which enjoys electric-magnetic duality invariance.¹¹ In fact, the U(1) invariance (4.19) of the ModMax oscillator descends directly from this electricmagnetic duality symmetry, which can be written as

$$z_{\mu\nu} \rightarrow e^{i\alpha} z_{\mu\nu}, \qquad z_{\mu\nu} = F_{\mu\nu} + i\tilde{F}_{\mu\nu}, \qquad (4.21)$$

where $F_{\mu\nu}$ is the field strength of the 4*d* electrodynamics theory and $\tilde{F}_{\mu\nu}$ is its Hodge dual.

It is straightforward to see that any deformation of the isotropic harmonic oscillator which is constructed from the Hamiltonian and the conserved angular momentum will preserve invariance under the U(1) duality transformation (4.19). This is the 1*d* version of the statement that any deformation of a theory of self-dual electrodynamics in four spacetime dimensions, where the deforming operator is a function of the energy-momentum tensor of the theory, will preserve electric-magnetic duality invariance. See Ref. [53] for further discussion and examples of such duality-invariance-preserving stress tensor flows.

1. Flow equation for partition functions

As we have seen in Secs. II and III, both the classical and quantum partition functions for a theory deformed by a function of conserved charges satisfy the same differential equation. We will now study this differential equation in the case of the ModMax oscillator, which obeys a flow driven by the operator \mathcal{R} or \mathfrak{R} introduced in Eq. (2.52). This falls into the class of nonanalytic deformations that we briefly considered around Eq. (2.67). In this case, the differential equation (2.73) simplifies considerably.

The reason for this simplification is the following. We have seen that the solution to the flow equation in this case is given by the Lagrangian (4.10) or Hamiltonian (4.14), which satisfies the equations

$$\frac{\partial^2 L_{\gamma}}{\partial \gamma^2} = L_{\gamma}, \qquad \frac{\partial^2 H_{\gamma}}{\partial \gamma^2} = H_{\gamma}. \tag{4.22}$$

Because $\partial_{\gamma}L_{\gamma} = \pm \mathcal{R}^{(\gamma)}$ and $\partial_{\gamma}H_{\gamma} = \mp \mathfrak{R}^{(\gamma)}$, this means

$$\frac{\partial \mathcal{R}^{(\gamma)}}{\partial \gamma} = \pm L_{\gamma}, \qquad \frac{\partial \mathfrak{R}^{(\gamma)}}{\partial \gamma} = \mp H_{\gamma}. \qquad (4.23)$$

The flow equation (2.73) then becomes

$$\partial_{\gamma}^{2} \mathcal{Z} = \langle -\beta H_{\gamma} + \beta^{2} (H_{\gamma}^{2} - J_{\gamma}^{2}) \rangle$$

= $\beta \partial_{\beta} \mathcal{Z} + \beta^{2} \partial_{\beta}^{2} \mathcal{Z} - \beta^{2} \partial_{\mu}^{2} \mathcal{Z},$ (4.24)

where in the last step we have expressed quantities in terms of derivatives of the partition function. Note that, if we had instead chosen a different normalization for the $1d \operatorname{root}-T\bar{T}$

¹¹See Ref. [52] for a manifestly duality-invariant presentation of the stress tensor for the ModMax and ModMax-Born-Infeld theories.

$$\partial_{\gamma}^2 \mathcal{Z} - \beta \partial_{\beta} \mathcal{Z} - \beta^2 c_0^2 (\partial_{\beta}^2 \mathcal{Z} - \partial_{\mu}^2 \mathcal{Z}) = 0.$$
 (4.25)

For the choice $c_0 = \frac{1}{2}$, this matches the dimensional reduction of the conjectured flow equation for the torus partition function of a root- $T\bar{T}$ deformed CFT given in (3.67).

Equation (4.24) is very nearly of a familiar form. To see this, it is convenient to Wick rotate $\mu \rightarrow i\mu$ and $\gamma \rightarrow i\gamma$, which reverses the signs on two terms. The resulting differential equation can be written as

$$0 = \frac{1}{\beta} \partial_{\beta} (\beta \partial_{\beta} \mathcal{Z}) + \frac{1}{\beta^2} \partial_{\gamma}^2 \mathcal{Z} + \partial_{\mu}^2 \mathcal{Z}.$$
(4.26)

This is identical to the Laplace equation for a function $f: \mathbb{R}^3 \to \mathbb{R}$ written in cylindrical coordinates (r, θ, z) , namely

$$0 = \frac{1}{r}\frac{\partial}{\partial r}\left(r\frac{\partial f}{\partial r}\right) + \frac{1}{r^2}\frac{\partial^2 f}{\partial \theta^2} + \frac{\partial^2 f}{\partial z^2},\qquad(4.27)$$

where the roles of the coordinates (r, θ, z) are played by (β, γ, μ) , respectively.

One may therefore solve the flow equation for $\mathcal{Z}(\beta, \gamma, \mu)$ by separation of variables, in a manner analogous to that which is done when studying electrodynamics in cylindrical coordinates. The original differential equation (4.24), before Wick rotating the parameters μ and γ , has different signs than those that appear in the cylindrical Laplace equation, leading to the appearance of slightly different Bessel functions. Any function of the form

$$f_{a,b}(\beta,\gamma,\mu) = (c_1 e^{-a\gamma} + c_2 e^{a\gamma})(c_3 e^{-b\mu} + c_4 e^{b\mu})(c_5 I_a(b\beta) + c_6 K_a(b\beta)),$$
(4.28)

for constants *a*, *b*, and $c_1, ..., c_6$, is a solution to (4.24). Here $I_{\nu}(x)$ and $K_{\nu}(x)$ are modified Bessel functions of the first kind and second kind, respectively. A general solution to the flow equation will therefore be a sum or integral of such functions $f_{a,b}$ for various choices of *a* and *b*. Physical considerations will also restrict the choices of the parameters c_i . For instance, the modified Bessel functions of the second kind $K_{\nu}(x)$ are generically divergent as $x \to 0$, whereas we expect a partition function to be finite at $\beta = 0$, so one should normally set $c_6 = 0$ in cases of physical interest.

We note that formal analytic continuation of flow equations to imaginary values of the parameters, such as that which relates (4.24) and (4.26), has sometimes been useful in previous work. For instance, in [19] such a continuation of a $T\bar{T}$ -like parameter λ was useful in relating the flow equations that produce the 4*d* Born-Infeld and reverse-Born-Infeld theories, which are two of the four solutions to the zero-birefringence condition for 4*d* nonlinear electrodynamics [54].

2. Direct computation of classical partition function

As a warm-up for our study of the quantum partition function, and in order to illustrate an example of the Laplace equation which the deformed flavored partition functions satisfy, we will now perform a direct computation of the classical partition function for the 2d ModMax oscillator. The resulting formulas will turn out to be tidier if we evaluate this partition function with an imaginary chemical potential for the angular momentum, which merely reverses a sign in the corresponding flow equation. We will therefore compute

$$\mathcal{Z}(\beta,\gamma,\mu) = \frac{1}{(2\pi)^2} \int dx \, dy \, dp_x \, dp_y \, \exp\left(-\beta H_\gamma + i\mu J_\gamma\right),$$
(4.29)

where all integrals run from $-\infty$ to ∞ . Writing this integrand explicitly in terms of the positions and momenta, and making the sign choice for which the two terms in the Hamiltonian H_{γ} are both manifestly positive, the integral we wish to evaluate is

$$\begin{aligned} \mathcal{Z}(\beta,\gamma,\mu) &= \frac{1}{(2\pi)^2} \int dx \, dy \, dp_x \, dp_y \, \exp\left[-\beta \left(\frac{1}{2} \cosh(\gamma)(p_x^2 + p_y^2 + x^2 + y^2)\right) \right. \\ &\left. + \frac{1}{2} \sinh(\gamma) \sqrt{((p_y + x)^2 + (p_x - y)^2)((p_y - x)^2 + (p_x + y)^2)} \right) + i\mu(xp_y - yp_x) \right]. \end{aligned}$$
(4.30)

Note that we have used the expression $J_{\gamma} = xp_y - yp_x$ in Eq. (4.30) since the deformed angular momentum in Hamiltonian variables takes the same form as the undeformed angular momentum, as we pointed out around Eq. (4.13).

It is convenient to perform the change of variables

$$u_1 = p_y + x, \qquad u_2 = p_x - y, \qquad v_1 = p_y - x, \qquad v_2 = p_x + y,$$
 (4.31)

so that the integral becomes

$$\mathcal{Z}(\beta,\gamma,\mu) = \frac{1}{4(2\pi)^2} \int du_1 \, du_2 \, dv_1 \, dv_2 \, \exp\left[-\frac{\beta}{4}(\cosh(\gamma)(u_1^2 + u_2^2 + v_1^2 + v_2^2) + 2\sinh(\gamma)\sqrt{u_1^2 + u_2^2}\sqrt{v_1^2 + v_2^2}\right] + \frac{i\mu}{4}(u_1^2 + u_2^2 - v_1^2 - v_2^2)\right].$$
(4.32)

Note that this change of variables has decoupled the square-root interaction into two factors. We can now go to polar coordinates in the (u_1, u_2) and (v_1, v_2) planes as

$$u_1 = r_u \cos(\theta_u), \qquad u_2 = r_u \sin(\theta_u), \qquad v_1 = r_v \cos(\theta_v), \qquad v_2 = r_v \sin(\theta_v).$$
 (4.33)

Then our partition function is

$$\mathcal{Z}(\beta,\gamma,\mu) = \frac{1}{4(2\pi)^2} \int_0^\infty dr_u \int_0^\infty dr_v \int_0^{2\pi} d\theta_u \int_0^{2\pi} d\theta_v r_u r_v \exp\left[-\frac{\beta}{4}(\cosh(\gamma)(r_u^2 + r_v^2) + 2\sinh(\gamma)r_u r_v) + \frac{i\mu}{4}(r_u^2 - r_v^2)\right].$$
(4.34)

The angular integrals give factors of 2π , whereas the resulting radial integrals can be evaluated in closed form, and we find

$$\mathcal{Z}(\beta,\gamma,\mu) = \frac{1}{\beta^2 + \mu^2} \left(1 - \frac{\sinh(\gamma)}{\sqrt{1 + \frac{\mu^2}{\beta^2}}} \arctan\left(\operatorname{csch}(\gamma)\sqrt{1 + \frac{\mu^2}{\beta^2}}\right) \right).$$
(4.35)

This is our final expression for the classical grand canonical partition function for the 2*d* ModMax oscillator at inverse temperature β and with imaginary chemical potential $i\mu$ for the angular momentum. One can, of course, obtain the result for the opposite sign choice in the Hamiltonian by sending $\gamma \rightarrow -\gamma$. The result (4.35) clearly reduces to the corresponding partition function for the ordinary 2*d* harmonic oscillator when $\gamma = 0$. One can also check by explicit computation that it obeys the partial differential equation

$$(\partial_{\gamma}^2 - \beta^2 \partial_{\beta}^2 - \beta \partial_{\beta} - \beta^2 \partial_{\mu}^2) \mathcal{Z}(\beta, \gamma, \mu) = 0, \qquad (4.36)$$

which is equivalent to (4.24) after Wick rotating the chemical potential $\mu \rightarrow i\mu$. One can recover the corresponding solution with a real chemical potential by reversing the signs of all instances of μ^2 in Eq. (4.35).

This result gives one example of a partition function that satisfies a flow equation driven by a nonanalytic combination of charges, namely the 1*d* root- $T\bar{T}$ deformation. When the chemical potential is set to zero, the deformed partition function (4.35) is simply a γ -dependent rescaling of the undeformed partition function. However, when μ and γ are both finite, the temperature dependence is modified in a more interesting way. We will see shortly that, in the quantum theory, even for $\mu = 0$ the deformed

partition function is not simply a rescaling of the undeformed partition function.

B. Quantum aspects

We now turn to the main subject of this work, which is the quantum mechanics of the ModMax oscillator. At first glance, it is not so clear that one should be able to quantize this theory at all. Naively, one would like to begin with the classical Hamiltonian (4.14) and promote all position and momentum variables to operators. This requires one to make sense of the operator square root in the second term, which takes the form

$$\sqrt{((\hat{p}_y + \hat{x})^2 + (\hat{p}_x - \hat{y})^2)((\hat{p}_y - \hat{x})^2 + (\hat{p}_x + \hat{y})^2)}.$$
 (4.37)

It is not immediately obvious what this operator should mean. First, note that we would not expect that it is possible to define an operator square root for a generic combination of position and momentum operators, at least without additional assumptions such as positivity. For instance, the expression $\sqrt{\hat{x}}$ does not give a conventional Hermitian operator; even if one attempts to define it by diagonalizing the position operator and declaring $\sqrt{\hat{x}}|x\rangle = \sqrt{x}|x\rangle$, this operator will have imaginary eigenvalues for negative positions *x*. In our case, we are aided in interpreting the operator (4.37) by the fact that it is positive definite—which allows us to define it by diagonalization and by taking square roots and because it is a function of conserved charges in the undeformed theory, which allows us to write flow equations for quantities in the deformed theory using the results of Sec. III. We will begin by attempting to understand the operator (4.37) directly using raising and lowering operators in the theory of the undeformed harmonic oscillator.

1. Ladder operator representation

We can develop one useful perspective on the ModMax oscillator by rewriting the Hamiltonian in terms of creation and annihilation operators. As usual, when studying the undeformed theory of an isotropic 2d harmonic oscillator with Hamiltonian

$$\hat{H}_0 = \frac{1}{2}(\hat{p}_x^2 + \hat{p}_y^2 + \hat{x}^2 + \hat{y}^2), \qquad (4.38)$$

it is natural to define the annihilation operators

$$\hat{a}_x = \frac{1}{\sqrt{2}}(\hat{x} + i\hat{p}_x), \qquad \hat{a}_y = \frac{1}{\sqrt{2}}(\hat{y} + i\hat{p}_y), \quad (4.39)$$

whose Hermitian conjugates are the creation operators,

$$\hat{a}_x^{\dagger} = \frac{1}{\sqrt{2}}(\hat{x} - i\hat{p}_x), \qquad \hat{a}_y^{\dagger} = \frac{1}{\sqrt{2}}(\hat{y} - i\hat{p}_y).$$
 (4.40)

In terms of these operators the Hamiltonian takes the standard form

$$\hat{H}_0 = 1 + \hat{a}_x^{\dagger} \hat{a}_x + \hat{a}_y^{\dagger} \hat{a}_y = 1 + \hat{N}_x + \hat{N}_y, \qquad (4.41)$$

where we have defined the number operators $\hat{N}_i = \hat{a}_i^{\dagger} \hat{a}_i$. However, the angular momentum operator can be made more transparent by a change of basis. We can instead define the "circularly polarized" linear combinations

$$\hat{a}_L = \frac{1}{\sqrt{2}}(\hat{a}_x + i\hat{a}_y), \qquad \hat{a}_R = \frac{1}{\sqrt{2}}(\hat{a}_x - i\hat{a}_y).$$
 (4.42)

The corresponding number operators, $\hat{N}_L = \hat{a}_L^{\dagger} \hat{a}_L$ and $\hat{N}_R = \hat{a}_R^{\dagger} \hat{a}_R$, count the numbers of left-moving and rightmoving circular quanta. This is a useful way to leverage the rotational symmetry of the problem, since the angular momentum operator is now simply

$$\hat{J} = \hat{N}_R - \hat{N}_L, \qquad (4.43)$$

and the Hamiltonian is

$$\hat{H}_0 = 1 + \hat{N}_L + \hat{N}_R. \tag{4.44}$$

We now turn to the 1d root- $T\bar{T}$ deformation that generates the interaction term in the ModMax oscillator Lagrangian. This operator is

$$\hat{\mathbf{\Re}} = \sqrt{\hat{H}_0^2 - \hat{J}^2} = \sqrt{(1 + 2\hat{N}_L)(1 + 2\hat{N}_R)}.$$
(4.45)

The argument of the square root factorizes into left-moving and right-moving pieces. Because the left-moving and right-moving operators commute, we are free to split the square root into separate factors:

$$\hat{\mathfrak{R}} = \sqrt{1 + 2\hat{N}_L}\sqrt{1 + 2\hat{N}_R}.$$
(4.46)

Each of the operators $1 + 2\hat{N}_L$ and $1 + 2\hat{N}_R$ have strictly positive eigenvalues, and it is therefore possible to define an operator square root. This is equivalent to defining the square-root operators by the Taylor series expansions

$$\sqrt{1+2\hat{N}_L} = \sum_{k=0}^{\infty} {\binom{\frac{1}{2}}{k}} (2\hat{N}_L)^k,$$
$$\sqrt{1+2\hat{N}_R} = \sum_{k=0}^{\infty} {\binom{\frac{1}{2}}{k}} (2\hat{N}_R)^k.$$
(4.47)

Either of these infinite sums is convergent and well-defined when acting on any state in the Hilbert space. We can therefore write the Hamiltonian of the ModMax oscillator as

$$\hat{H}_{\gamma} = \cosh(\gamma)(1 + \hat{N}_L + \hat{N}_R) + \sinh(\gamma)\sqrt{1 + 2\hat{N}_L}\sqrt{1 + 2\hat{N}_R}, \quad (4.48)$$

where we have again chosen the positive sign for concreteness, although one can obtain the other sign choice by taking γ to be negative.

One could study the quantum mechanics of this theory in essentially the same way that we have described above for the canonical quantization prescription. That is, one considers a complete basis of eigenstates $|N_L, N_R\rangle$ of the undeformed Hamiltonian and angular momentum, and then shifts each of their energy eigenvalues according to (4.48).

However, this number operator representation also allows us to see a simple way to generate the ModMax oscillator in one step from the undeformed isotropic harmonic oscillator. Let us introduce operators

$$\hat{M}_L = \sqrt{1 + 2\hat{N}_L}, \qquad \hat{M}_R = \sqrt{1 + 2\hat{N}_R}, \qquad (4.49)$$

which are defined by the convergent Taylor series expansions (4.47). Then the undeformed Hamiltonian can be written as

$$\hat{H}_0 = \frac{1}{2}(\hat{M}_L^2 + \hat{M}_R^2).$$
(4.50)

Suppose that one now performs the transformation

$$\hat{M}_L \to \hat{M}_L^{(\gamma)} = \cosh(\gamma)\hat{M}_L + \sinh(\gamma)\hat{M}_R,$$
$$\hat{M}_R \to \hat{M}_R^{(\gamma)} = \cosh(\gamma)\hat{M}_R + \sinh(\gamma)\hat{M}_L.$$
(4.51)

Then one finds

$$\hat{H}_{0} \rightarrow \hat{H}_{2\gamma} = \frac{1}{2} ((\hat{M}_{L}^{(\gamma)})^{2} + (\hat{M}_{R}^{(\gamma)})^{2}),$$

$$= \cosh(2\gamma)(1 + \hat{N}_{R} + \hat{N}_{L})$$

$$+ \sinh(2\gamma)\sqrt{1 + 2\hat{N}_{L}}\sqrt{1 + 2\hat{N}_{R}}.$$
(4.52)

This is exactly the deformed Hamiltonian of Eq. (4.48), except at parameter 2γ rather than γ . Therefore, performing a "boost" in the space of the operators \hat{M}_L and \hat{M}_R has the effect of generating the ModMax oscillator Hamiltonian in a single step, rather than via a flow equation that is defined by infinitesimally deforming by the operator $\hat{\Re}$.

This is the same structure as the classical deformation map that was introduced in Sec. 6 of [30]. In that case, a particular redefinition of positions and momenta had the effect of mapping $H_0 \rightarrow H_{2\gamma}$ in a single step, albeit at the level of the classical Hamiltonian. The transformation (4.51) on the numberlike operators \hat{M}_L and \hat{M}_R can be thought of as a quantum version of this deformation map.

2. Trace form of quantum partition function

We now consider the grand canonical partition function for the ModMax theory at the quantum level. Following the ladder operator representation of the Hamiltonian developed above, it is convenient to choose a basis of simultaneous eigenstates $|N_L, N_R\rangle$ of the number operators associated with the left- and right-circularly polarized creation and annihilation operators introduced in Eq. (4.42).

Using this basis, the flavored partition function with an imaginary chemical potential for the angular momentum $\hat{J} = \hat{N}_R - \hat{N}_L$ can be written as

$$\begin{aligned} \mathcal{Z}(\beta,\gamma,\mu) &= \operatorname{Tr}(\exp\left(-\beta\hat{H}_{\gamma} + i\mu\hat{J}_{\gamma}\right)) \\ &= \sum_{N_{L},N_{R}=0}^{\infty} \exp\left[-\beta(\cosh(\gamma)(1+\hat{N}_{R}+\hat{N}_{L})\right. \\ &+ \sinh(\gamma)\sqrt{1+2\hat{N}_{R}}\sqrt{1+2\hat{N}_{L}}) \\ &+ i\mu(\hat{N}_{R}-\hat{N}_{L})\right]. \end{aligned}$$
(4.53)

Here we have used the fact that $\hat{J}_{\gamma} = \hat{J}_0 = \hat{N}_R - \hat{N}_L$, following the comments around (4.13), and taken the positive sign choice in the Hamiltonian as usual.

Unlike in the case of the classical partition function, it does not seem to be possible to obtain a simple closed-form expression for the sum (4.53). However, it is straightforward to evaluate the trace perturbatively in the flow parameter γ . For instance, to leading order one finds

$$\begin{aligned} \mathcal{Z}(\beta,\gamma,\mu) &= \frac{1}{2\cosh(\beta) - 2\cos(\mu)} \\ &- 2\gamma\beta e^{\beta}\Phi\left(e^{-\beta - i\mu}, -\frac{1}{2}, \frac{1}{2}\right) \\ &\times \Phi\left(e^{-\beta + i\mu}, -\frac{1}{2}, \frac{1}{2}\right) + \mathcal{O}(\gamma^2), \end{aligned}$$
(4.54)

where $\Phi(z, s, a)$ is a special function known as the Lerch transcendent and defined by

$$\Phi(z, s, a) = \sum_{k=0}^{\infty} \frac{z^k}{(k+a)^s}.$$
(4.55)

Even when $\mu = 0$, the expression (4.54) for the partition function to order γ is not simply a rescaling of the undeformed partition function by a γ -dependent factor. This is unlike the classical partition function (4.35), to which the expression (4.53) reduces in the limit $\hbar \rightarrow 0$.¹² This gives one way to see that the quantum theory of the ModMax oscillator is richer than its classical counterpart, since even without a chemical potential μ , there is a nontrivial interplay between the inverse temperature β and flow parameter γ .

One can check that the infinite sum (4.53) satisfies the flow equation (4.24) for a theory deformed by the 1*d* root- $T\bar{T}$ operator, after specializing to an imaginary value of μ . As we mentioned, this differential equation can be solved by separation of variables, which gives a general solution that is a sum of factorized terms involving exponentials and Bessel functions. It is instructive to see how (4.53) can be brought into this form, since as written this sum is not obviously related to Bessel functions. This can be accomplished using a variant of the Jacobi-Anger expansion, which expresses a plane wave as a superposition of cylindrical waves. For any $z, \theta \in \mathbb{C}$, these identities take the form

$$e^{z\cos(\theta)} = I_0(z) + 2\sum_{k=1}^{\infty} I_k(z)\cos(k\theta),$$

$$e^{z\sin(\theta)} = I_0(z) + 2\sum_{k=0}^{\infty} (-1)^k I_{2k+1}(z)\sin((2k+1)\theta)$$

$$+ 2\sum_{k=1}^{\infty} (-1)^k I_{2k}(z)\cos(2k\theta).$$
(4.56)

¹²This fact is not obvious from equation (4.53) because we have set $\hbar = 1$ for simplicity. After restoring factors of \hbar and taking $\hbar \to 0$, the sum reduces to the integral (4.30), as it must.

See, for instance, Sec. 10.35 of [55]. To apply these identities to the partition function (4.53) for the quantum ModMax oscillator, we let $\theta = i\gamma$ in the identities (4.56). The full partition function can then be written as the expansion

$$\begin{aligned} \mathcal{Z}(\beta,\gamma,\mu) &= \sum_{N_L,N_R=0}^{\infty} \left\{ \left[\cos\left(\mu(N_R - N_L)\right) + i\sin\left(\mu(N_R - N_L)\right) \right] \right. \\ &\left. \cdot \left[I_0(-\beta(1 + N_R + N_L)) + 2\sum_{k=1}^{\infty} I_k(-\beta(1 + N_R + N_L))\cosh(k\gamma) \right] \right. \\ &\left. \cdot \left[I_0(-i\beta\sqrt{1 + 2N_R}\sqrt{1 + 2N_L}) + 2\sum_{k=1}^{\infty} (-1)^k I_{2k} \left(-i\beta\sqrt{1 + 2N_R}\sqrt{1 + 2N_L} \right) \cosh(2k\gamma) \right. \\ &\left. + 2i\sum_{k=0}^{\infty} (-1)^k I_{2k+1} \left(-i\beta\sqrt{1 + 2N_R}\sqrt{1 + 2N_L} \right) \sinh\left((2k + 1)\gamma\right) \right] \right\}. \end{aligned}$$

$$(4.57)$$

Although rather unwieldy, this expansion in Bessel functions makes the connection between the trace form (4.53) of the partition function and the Laplace-type equation, which it satisfies, more explicit.

3. Description of deformed propagator

To conclude, we will comment on the characterization of the propagator for the ModMax oscillator. Here we will be brief, since this is a direct application of the general results of Sec. III and leads to flow equations that are essentially identical to those for the classical and quantum partition function discussed above.

Because the Hamiltonian for the ModMax oscillator is of the form (3.50), there is no ambiguity in defining the propagator by a path integral representation, even at all orders in γ . Therefore, the flavored propagator is defined by the general phase space path integral given in Eq. (3.40). This propagator, with real chemical potential, satisfies

$$\partial_{\gamma}^{2}K - T\partial_{T}K - T^{2}(\partial_{T}^{2} + \partial_{\mu}^{2})K = 0, \qquad (4.58)$$

which is the same as the differential equation for the flavored partition function after making the replacement $\beta = iT$. However, we emphasize that the result is more general, since this holds for the propagator $K(\vec{x}_B, t_B; \vec{x}_A, t_A; \lambda; \mu)$ with any initial position \vec{x}_A and final position \vec{x}_B . In contrast, the thermal partition function is obtained from the Euclidean time propagator with periodic boundary conditions, which means that the initial and final positions are equal.

The differential equation (4.58) fully determines the propagator for the ModMax oscillator, given the initial condition $K(\gamma = 0)$ and the first derivative $\partial_{\gamma}K|_{\gamma=0}$, which is related to the expectation value of the 1*d* root- $T\bar{T}$ operator in the undeformed theory. The initial condition $K(\gamma = 0)$ is essentially the propagator for a 2*d* harmonic oscillator in a background magnetic field, which plays the role of the chemical potential for the angular momentum. This quantity can be computed by either canonical methods

or path integral methods; for the path integral computation, see the pedagogical review [56].

We also note that the propagator can be written using the kernel representation and the basis of states $|N_L, N_R\rangle$, which gives

$$K(\vec{x}_B, t_B; \vec{x}_A, t_A; \lambda; \mu) = \sum_{N_L, N_R=0}^{\infty} \phi_{N_L, N_R}^*(\vec{x}_A) \phi_{N_L, N_R}(\vec{x}_B)$$

$$\times \exp\left[-i\left(\cosh(\gamma)(1+\hat{N}_R+\hat{N}_L)\right.\right.$$

$$+ \sinh(\gamma)\sqrt{1+2\hat{N}_R}\sqrt{1+2\hat{N}_L}\right)$$

$$\times (t_B - t_A) + \mu(\hat{N}_R - \hat{N}_L)\right],$$

$$(4.59)$$

where the ϕ_{N_L,N_R} are harmonic oscillator wave functions, which are known in closed form.

Our characterization of the propagator K, including that it satisfies the Laplace-like differential equation (4.58), is one of the main results of this work. Because the propagator for the ModMax oscillator is completely determined by the above considerations, this essentially constitutes a full solution of the model. Any physical question involving time evolution of states can in principle be extracted from the function K. This therefore completes our study of the quantum-mechanical theory of the ModMax oscillator.

V. CONCLUSION

In this work, we have studied general deformations of 1d theories by conserved charges, at both the classical and the quantum levels. This has allowed us to obtain flow equations for quantities in the theory of the ModMax oscillator, which is the dimensional reduction of the 4d ModMax theory. In particular, we have found that the thermal partition function in the quantum theory of the ModMax oscillator—or,

relatedly, the real-time propagator—satisfies a certain partial differential equation that is related by Wick rotation to the Laplace equation in 3d cylindrical coordinates.

One way of summarizing our analysis is to say that *any* deformation of a quantum-mechanical theory by conserved charges is essentially "solvable" in the sense that one can write differential equations that relate quantities in the deformed theory, such as the propagator or partition function, to those in the undeformed theory. The results on the ModMax oscillator are a special case of this fact when the deformation is driven by the 1d root- $T\bar{T}$ operator. Furthermore, the quantization of such charge-deformed models is unambiguous, since one obtains equivalent flow equations using either canonical quantization or path integral quantization. We therefore conclude that the fairly broad class of theories obtained through deformations by conserved charges should be included among other examples of solvable deformations of quantum mechanics, such as the one in which the quadratic kinetic term is replaced by one involving a hyperbolic cosine [57].

There are several directions for future investigation, some of which we outline below.

A. One-loop calculation

In this manuscript, we have focused on finding *exact* flow equations for observables in the quantum theory of the ModMax oscillator, such as the propagator. We have also studied certain quantities in the classical theory, such as the classical partition function, for which it is possible to obtain a closed-form expression.

However, it would also be interesting to study *semi-classical* expressions for quantum observables by performing an expansion in \hbar . In the limit as $\gamma \rightarrow 0$, the ModMax oscillator reduces to the ordinary harmonic oscillator, which is one-loop exact. It seems very unlikely that the deformed theory is also one-loop exact due to the complicated nature of the interaction term. One could attempt to compute the one-loop correction around the classical solution to the equations of motion for the ModMax oscillator and examine how closely this reproduces the full quantum results.

To do this, one would need to expand the phase space path integral that defines the propagator for the ModMax oscillator and retain terms up to quadratic order in fluctuations around the classical path. Fortunately, it is straightforward to write down the general classical solution to the equations of motion for the ModMax oscillator following [30]. Given a set of initial conditions (x_0, y_0) and $(p_{x,0}, p_{y,0})$, one can evaluate the conserved energy H_0 and angular momentum J_0 corresponding to this initial condition,

$$E_0 = \frac{1}{2}(x^2 + y^2 + p_x^2 + p_y^2), \qquad J = xp_y - yp_x, \quad (5.1)$$

and then define

$$A = \cosh(\gamma) - \frac{\sinh(\gamma)H_0}{\sqrt{H_0^2 - J_0^2}}, \qquad B = \frac{\sinh(\gamma)J_0}{\sqrt{H_0^2 - J_0^2}}.$$
 (5.2)

The general solution to the deformed equations of motion is given by

$$\begin{aligned} x(t) &= \sin(At)(p_{x,0}\cos(Bt) + p_{y,0}\sin(Bt)) + \cos(At)(x_0\cos(Bt) + y_0\sin(Bt)), \\ y(t) &= \sin(At)(p_{y,0}\cos(Bt) - p_{x,0}\sin(Bt)) + \cos(At)(y_0\cos(Bt) - x_0\sin(Bt)), \\ p_x(t) &= \cos(At)(p_{x,0}\cos(Bt) + p_{y,0}\sin(Bt)) - \sin(At)(x_0\cos(Bt) + y_0\sin(Bt)), \\ p_y(t) &= \cos(At)(p_{y,0}\cos(Bt) - p_{x,0}\sin(Bt)) - \sin(At)(y_0\cos(Bt) - x_0\sin(Bt)). \end{aligned}$$
(5.3)

One could then perform a one-loop computation by defining

$$x^{i} = x^{i}_{cl} + \delta x^{i}, \qquad p^{i} = p^{i}_{cl} + \delta p^{i}, \qquad (5.4)$$

where $x_{cl}^i(t)$ and $p_{cl}^i(t)$ are solutions to the equations of motion (5.3), and then performing the phase space path integral over δp^i and δx^i . This is slightly more involved than a semiclassical computation in the ordinary Feynman path integral, which gives a one-loop determinant. In this case, one would first expand the Hamiltonian action to write

$$H = f_1(\delta x)(\delta p^2) + f_2(\delta x)\delta p + f_3(\delta x), \quad (5.5)$$

where we suppress indices on the fluctuations. It is still possible to carry out the $\mathcal{D}(\delta p)$ path integral over momentum fluctuations for such a Hamiltonian, as described in Sec. 1.4 of [50]. After performing the momentum path integral, one is left with a path integral over $\mathcal{D}(\delta x)$ with a modified Lagrangian. It should then be possible to complete the semiclassical expansion by computing the oneloop determinant associated with the fluctuations δx around the classical solution in this Feynman path integral.

B. Laplace equation for flavored partition function and Narain moduli space

One of the main results of this manuscript is that the propagator for a root- $T\bar{T}$ deformed theory satisfies a partial

differential equation that—up to signs which can be eliminated by choosing imaginary values of parameters—is identical to the Laplace equation in three-dimensional cylindrical coordinates. Similar Laplace-type equations have appeared in the description of certain torus partition functions for two-dimensional conformal field theories. For instance, in [58] the authors study flavored CFT partition functions that satisfy a Laplace equation where the Laplacian acts both on a space of Narain lattices and on a space of chemical potentials. This is of the same schematic form as the Laplace equation that we have derived for the ModMax oscillator partition function.

Another observation along similar lines is the following. We have pointed out that the flow equation we obtained for a 1*d* root- $T\bar{T}$ deformed theory descends by dimensional reduction from the flow equation (3.65) for a root- $T\bar{T}$ deformed torus partition function. Equation (3.65) is structurally similar to the equation obeyed by certain theta functions. For instance, consider a theory of *D* compact bosons that parametrize a target-space torus T^D that has some collection of moduli which we schematically indicate by *m*. The partition function for this theory is

$$Z(m,\tau) = \frac{\Theta(m,\tau)}{|\eta(\tau)|^{2D}},$$
(5.6)

where $\Theta(m, \tau)$ is a Siegel-Narain theta function, which obeys the differential equation

$$(-\tau_2^2 \partial_\tau \partial_{\bar{\tau}} - \tau_2 \partial_{\tau_2} - \Delta_{\mathcal{M}_D}) \Theta(m, \tau) = 0, \qquad (5.7)$$

where $\Delta_{\mathcal{M}_D}$ is the natural Laplacian on the Narain moduli space that parametrizes the T^D . For instance, in the case of a single compact boson, the target space is a circle of radius *R* and the Laplacian is

$$\Delta_{\mathcal{M}_1} = -\frac{1}{4} \left(R \frac{d}{dR} \right)^2. \tag{5.8}$$

The structure of Eq. (3.65) is almost identical to (5.7), except with the Laplacian on moduli space replaced with the second derivative with respect to γ . Also note that the root- $T\bar{T}$ flow equation involves the partition function itself, while (5.7) holds for the function Θ appearing in the numerator of the partition function, not for the full combination including the eta function in the denominator.

It would be interesting to understand whether there is a deeper relationship between these flow equations for root- $T\bar{T}$ deformed partition functions, and their dimensional reductions to 1*d*, and properties of Narain moduli space.

C. Coupling to worldline gravity

As we have emphasized, there may be several inequivalent quantization schemes—or "UV completions"—for a particular classical theory. In particular, this is true for theories obtained from the (0 + 1)-dimensional version of the $T\bar{T}$ deformation. One way to see the difference between two such choices of the quantization scheme is by examining the resulting thermal partition functions. As we reviewed in Sec. II C, the quantization procedure for the $1d T\bar{T}$ deformation that gives rise to the flow equation (2.65) admits a solution for the deformed partition function that can be written as an integral transform of the undeformed partition function, Eq. (2.66).

However, there is a second UV completion of this deformation that is defined by coupling the seed theory to worldline gravity [35]. In this prescription, the deformed partition function is defined by the path integral

$$Z_{\lambda}(\beta) = \int \frac{\mathcal{D}e \,\mathcal{D}X \,\mathcal{D}\sigma}{\text{Vol}} \exp\left(-S_0(e;X) - S(\lambda;e,\sigma)\right), \quad (5.9)$$

where X represents the fields of the undeformed theory, which are now minimally coupled to an einbein e, and σ is an auxiliary scalar field with an action

$$S(\lambda; e, \sigma) = \frac{1}{2\lambda} \int_0^{\beta'} d\tau \, e(e^{-1}\partial_\tau \sigma - 1)^2.$$
 (5.10)

After evaluating the path integral, one finds that this quantization scheme produces the deformed partition function

$$Z_{\lambda}(\beta) = \int_{0}^{\infty} d\beta' \frac{\beta}{\sqrt{2\pi\lambda}\beta'^{3/2}} \sum_{m} \exp\left(-\frac{(m\beta - \beta')^{2}}{2\beta'\lambda}\right) Z_{0}(\beta'),$$
(5.11)

which is a different partition function than the result (2.66), although they agree in the unit winding sector (m = 1) up to normalization for the flow parameter.¹³

It would be very interesting to find an analog of this worldline gravity prescription for the 1d root- $T\bar{T}$ -like deformation. It seems likely that one would need to couple the undeformed theory to both an einbein $e(\tau)$ and a gauge field for the SO(N) symmetry, which plays the role of a time-dependent magnetic field. This would be in accord with the fact that, in order to obtain a flow equation for the partition function using our simpler quantization prescription for the root- $T\bar{T}$ deformed quantum mechanics, we were forced to turn on a chemical potential μ for the angular momentum. Finding a path integral expression that represents this alternative, worldline gravity quantization prescription would allow us to better understand the available choices of UV completions for ModMax-like theories of quantum mechanics.

¹³In particular, one can redefine $\lambda \to -4\lambda$ in (2.66) to match the conventions of (5.11).

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APPENDIX: FIRST-ORDER ANALYSIS OF LAGRANGIAN AND HAMILTONIAN FLOWS

In this appendix, we will consider the analog of Theorem 1 where one uses the relationship between the conjugate momenta p^i and velocities \dot{x}^i in the *undeformed* theory (defined by L_0 and H_0) rather than in the deformed theory (defined by L_{λ} and H_{λ}). If one does not correct the definition of the conjugate momenta, the resulting deformations of the Lagrangian and Hamiltonian will be equivalent only to leading order in the deformation parameter. The logic of this proof was originally presented in Appendix A of [40], which focused on $T\bar{T}$ deformations of two-dimensional quantum field theories describing a field ϕ and conjugate momentum π . Here we will instead consider 1*d* theories that describe the dynamics of a collection of positions x^i and velocities \dot{x}^i , since this is the primary focus of the present work, although the reasoning is almost identical.

Let us represent the deformation of the Hamiltonian to first order as follows:

$$H(x^i, p^i) = H_0(x^i, p^i) + \epsilon \mathfrak{O}(x^i, p^i).$$
(A1)

We claim that this is equivalent to a deformation of the Lagrangian to first order as

$$L(x^{i}, \dot{x}^{i}) = L_{0}(x^{i}, \dot{x}^{i}) - \epsilon \mathfrak{O}(x^{i}, f_{0}^{i}(\dot{x}^{j})).$$
(A2)

Here we define f_0^i to be the function that relates \dot{x}^i and p^i when $\epsilon = 0$. More precisely,

$$p^{i} = \frac{\partial L_{0}}{\partial \dot{x}^{i}} = f_{0}^{i}(\dot{x}^{j}). \tag{A3}$$

Because we are defining the deformation in terms of the Hamiltonian, we view the momenta p^i as independent variables which do not depend on the deformation parameter λ ; this corresponds to the forward direction of Theorem 1, rather than the converse. Consequently the definition of the velocities \dot{x}^i changes due to the deformation.

We use the Legendre transform to write the deformed Lagrangian in terms of the Hamiltonian,

$$L(x^{i}, p^{i}) = p^{i} \dot{x}^{i} - H_{0}(x^{i}, p^{i}) - \epsilon \mathfrak{O}(x^{i}, p^{i}).$$
(A4)

Next, we solve the equations of motion to write p^i as a function of x^i and \dot{x}^i . We assume that this solution can be written as an infinite series as follows:

$$p^{i} = f^{i}(x^{j}, \dot{x}^{j}) = f^{i}_{0}(x^{j}, \dot{x}^{j}) + \epsilon f^{i}_{1}(x^{j}, \dot{x}^{j}) + \mathcal{O}(\epsilon^{2}).$$
(A5)

By Hamilton's equations of motion, $\dot{x}^i = \frac{\partial H}{\partial p^i}$. We can use the equation for p^i to first order to obtain

$$\dot{x}^{i} = \frac{\partial H_{0}}{\partial p^{i}} (x^{j}, f_{0}^{j} + \epsilon f_{1}^{j}) + \epsilon \frac{\partial \mathfrak{D}}{\partial p^{i}} (x^{j}, f_{0}^{j} + \epsilon f_{1}^{j}).$$
(A6)

We expand the first term perturbatively to obtain the following expression for \dot{x}^i :

$$\dot{x}^{i} = \frac{\partial H}{\partial p^{i}}(x^{j}, f_{0}^{j}) + \epsilon f_{1}^{j} \frac{\partial^{2} H_{0}}{\partial p^{i} \partial p^{j}}(x^{k}, f_{0}^{k}) + \epsilon \frac{\partial \mathfrak{D}}{\partial p^{i}}(x^{j}, f_{0}^{j}) + \mathcal{O}(\epsilon^{2}).$$
(A7)

Matching the terms of order ϵ^0 on either side of (A7) gives

$$\dot{x}^{i} = \frac{\partial H_{0}}{\partial p^{i}} (x^{j}, f_{0}^{j}), \tag{A8}$$

which is the Hamilton equation of motion in the undeformed theory, and equating the terms of order ϵ gives

$$f_1^j \frac{\partial^2 H_0}{\partial p^i \partial p^j} + \frac{\partial \mathfrak{D}}{\partial p^i} = 0.$$
 (A9)

This relation describes how f_1^i is implicitly determined in terms of H_0 and \mathfrak{D} . Since the Legendre transform is an involution, we can express the deformed Lagrangian in terms of the deformed Hamiltonian,

$$L = p^{i} \dot{x}^{i} - H_{0}(x^{i}, p^{i}) - e \mathfrak{O}(x^{i}, p^{i}).$$
 (A10)

Using the equation for p^i , we obtain

$$\begin{split} L &= (f_0^i + \epsilon f_1^i) \dot{x}^i - H_0(x^i, f_0^i + \epsilon f_1^i) - \epsilon \mathfrak{O}(x^i, f_0^i + \epsilon f_1^i) \\ &= f_0^i \dot{x}^i - H_0(x^i, f_0^i) + \epsilon \left(\dot{x}^i - \frac{\partial H_0}{\partial p^i} (x^j, f_0^j) \right) f_1^i \\ &- \epsilon \mathfrak{O}(x^i, f_0^i). \end{split}$$
(A11)

By equation (A8), the term proportional to ϵ must vanish, leading to

$$L = L_0 - \epsilon \mathfrak{O}(x^i, f_0^i(\dot{x}^j)).$$
(A12)

This is the claim we sought to prove since deforming the Hamiltonian will give rise to the same results as deforming the Lagrangian according to (A2) to first order in ϵ .

Since we are only looking at the first order, it does not matter that the relation between \dot{x}^i and p^i changes. However,

as we saw in Theorem 1, one can extend this argument to all orders in the deformation parameter by using the corrected relationship between velocities and conjugate momenta in the deformed theory.

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