

Classical Discrete Time Crystals

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(Dated: December 10, 2019)

The spontaneous breaking of time-translation symmetry in periodically driven quantum systems leads to a new phase of matter: discrete time crystals (DTC). This phase exhibits collective subharmonic oscillations that depend upon an interplay of non-equilibrium driving, many-body interactions, and the breakdown of ergodicity. However, subharmonic responses are also a well-known feature of classical dynamical systems ranging from predator-prey models to Faraday waves and AC-driven charge density waves. This raises the question of whether these classical phenomena display the same rigidity characteristic of a quantum DTC. In this work, we explore this question in the context of periodically driven Hamiltonian dynamics coupled to a finite-temperature bath, which provides both friction and, crucially, noise. Focusing on one-dimensional chains, where in equilibrium any transition would be forbidden at finite temperature, we provide evidence that the combination of noise and interactions drives a sharp, first-order dynamical phase transition between a discrete time-translation invariant phase and an activated classical discrete time crystal (CDTC) in which time-translation symmetry is broken out to exponentially-long time scales. Power-law correlations are present along a first-order line which terminates at a critical point. We analyze the transition by mapping it to the locked-to-sliding transition of a DC-driven charge density wave. Our work points to a classical limit for quantum time crystals, and raises several intriguing questions concerning the non-equilibrium universality class of the CDTC critical point.

Subharmonic entrainment occurs when the long time dynamics of a system manifest a period that is a fixed multiple of the period of the underlying equations of motion [1–7]. Such subharmonic behavior is ubiquitous in deterministic dynamical systems; most notably for example, discrete maps, e.g. $x \rightarrow f(x)$, can exhibit stable period-doubled orbits [2, 8–10], and continuous-time systems can settle down into limit cycles [7, 11]. From the point of view of many-body physics, however, in order to consider this subharmonic response characteristic of a phase of matter, the system should satisfy certain properties which embody the notion of *rigidity*. First, the system should have many locally coupled degrees of freedom so that a notion of spatial dimension and thermodynamic limit can be defined. Second, the system’s subharmonic response should be stable to arbitrary perturbations of both the initial state and the equations of motion, so long as the lat-

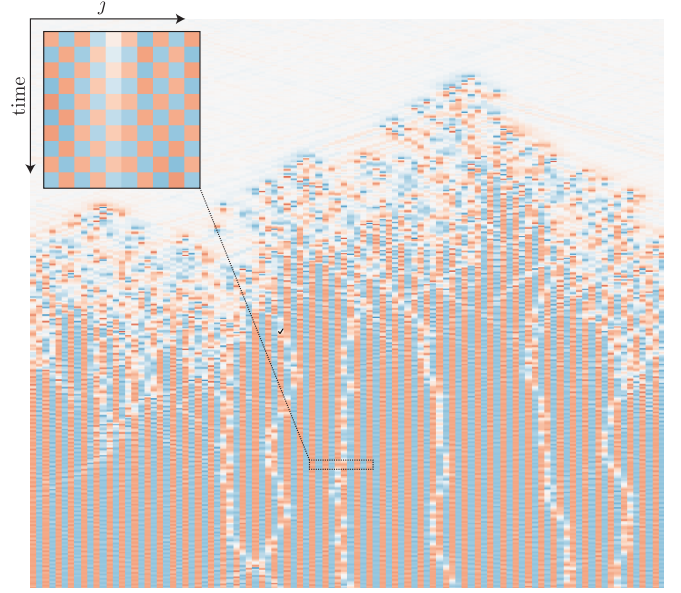


FIG. 1. **Period-doubled dynamics “boil” out of a uniform initial state.** In the main figure we present a stroboscopic view $q_j(2nt_D)$, with time n running down vertically ($0 < n < 1200$) and space j running horizontally over the $n_{\text{osc}} = 100$ oscillators. The color scale is $q_j < 0$ red, $q_j > 0$ blue, and $q_j \sim 0$ white. Note that we strobe every *two* driving periods, which is the frequency of the subharmonic response; hence the displayed phase of the oscillators varies slowly. In the inset, we show a detail of a smaller region strobbed at the driving frequency, $q_j(nt_D)$. The period-doubled oscillations $q_j(nt_D) \propto (-1)^n$ are now manifest. Strikingly, the correlations are anti-ferromagnetic both in time *and* space, even though the oscillators are coupled together ferromagnetically ($\omega_D = 1.958$, $g = 0.065$, $\delta = 0.067$, $\eta = 0.003$, $T = 0.004$). In the final state, there is a finite density of π -domain walls between the two different period-doubled solutions.

ter preserves the periodicity. Models with continuous time-translation invariance, such as the Van der Pol oscillator [11] and the Kuramoto model [12, 13] are not rigid in this sense (though the many-body synchronization transition of the latter is of interest its own right - we comment on its relation to the CDTC problem in the conclusions). Finally, the subharmonic response should have an infinite autocorrelation time, by analogy to “long-range” order.

Even within the constraints of these criteria, it turns out that general dynamical systems can still exhibit rigid subharmonic entrainment. The reason for this is that the dynamics about

fixed points can be strongly damped so that perturbations to either the state or the dynamics decay rapidly; in the presence of such contractive/dissipative dynamics, many-body subharmonic entrainment has been observed in a multitude of systems including: Faraday wave instabilities [3], driven charge density wave materials [4, 14–20] and Josephson junction arrays [21, 22].

The possibility of rigid subharmonic entrainment in the absence of contractive dynamics is significantly more subtle, but also particularly relevant [23–30]. Indeed, two broad classes of systems that fall into this category are time-periodic Hamiltonian dynamics in classical systems and unitary dynamics in quantum systems. Such systems are far more restrictive than general dynamical maps; Hamiltonian dynamics, for example, are volume preserving in phase space, thereby explicitly forbidding contractive dynamics. In the presence of an external drive (e.g. which sets the periodicity of the dynamics), energy conservation is broken and one generically expects the long-time dynamics of the many-body system to be completely ergodic. In certain fine-tuned special cases [31, 32], it may be possible for a driven, many-body system to avoid its ergodic fate [33]. An ergodic system can never exhibit true subharmonic rigidity since it is impossible for the system to remember which of the distinct subharmonic orbits (i.e. related by time translation symmetry) it began in.

To this end, a tremendous amount of recent excitement has focused on the discovery that rigid subharmonic entrainment can occur in a periodically driven (Floquet), unitary, many-body quantum system. Dubbed Floquet/discrete time crystals [34–42], this new phase of quantum matter relies crucially on many-body localization to prevent the drive-induced heating of the system to infinite temperature. While it is difficult to experimentally verify the long-time rigidity associated with a discrete time crystal, promising signatures of such behavior have been observed in spin systems [43, 44] for time-scales up to hundreds of Floquet cycles, despite the fact that localization is likely absent in such systems [41, 44–46].

A natural question thus arises: How quantum must a time crystal truly be? Is quantum mechanics important only insofar as it allows for many-body localization to prevent heating of the system? Or does it play a more fundamental role? If closed Hamiltonian dynamics cannot generically stabilize a time crystal due to heating, a natural generalization is to consider preventing such heating by coupling the system to a bath, most simply by adding friction. However, by adding *only* friction, one essentially reverts back to the damped case where the existence of rigid subharmonic entrainment is well known. On the other hand, if the bath is in equilibrium at finite temperature T , the fluctuation-dissipation theorem implies that friction must come with noise [47]. In classical systems, this noise can be captured as a Langevin force, $F_B(t) = -\eta\dot{q} + \xi(t)$, on each coordinate q , where η is the strength of the friction and $\xi(t)$ is a stochastic force with variance, $\langle \xi(t)\xi(t') \rangle = 2\eta T \delta(t - t')$. Taking $\eta > 0, T = 0$ reduces to the damped case where period-doubling is easily stabilized, while a combination of finite T and driving results

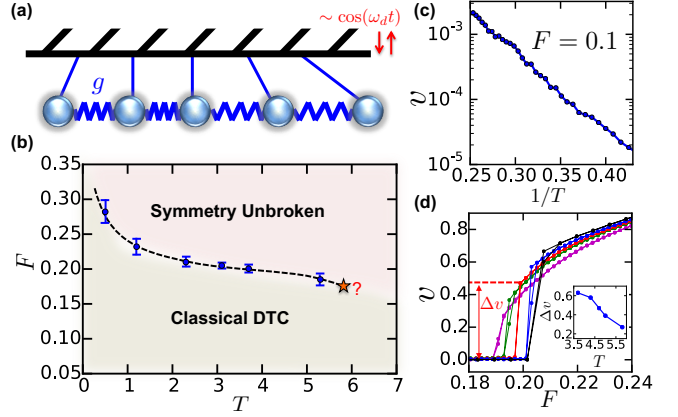


FIG. 2. Diagnostics and phase diagram of an activated classical discrete time crystal. **a)** Schematic of a one dimensional array of coupled non-linear pendula. The pendula are coupled via ferromagnetic interactions of strength g , and the system is parametrically driven at frequency ω_D . **b)** Phase diagram of the classical discrete time crystal as a function of $F = \frac{2\eta}{\delta}$ (where δ is the driving amplitude and η the damping) and the temperature T . At low- T there is a first order phase transition between the CDTC phase and the non-time-crystalline phase, while at high- T there is only a crossover. **c)** In the 1D CDTC, there is a finite rate of phase slips v between the two symmetry-breaking solutions. v fits very well to the Arrhenius form $v \sim e^{-\Delta/T}$, indicating the phase slips are activated. **d)** The phase transition is diagnosed by measuring the rate of phase slips v between the two time-translation related period doubled solutions. As we cross the first order line out of the CDTC, v jumps discontinuously.

in a truly non-equilibrium situation. The question concerning the existence of “classical discrete time crystals” (CDTC) can then be posed as follows: In what dimensions can a classical many-body system, coupled to an *equilibrium* bath, exhibit rigid subharmonic entrainment for either the closed case ($\eta = 0$), the zero-temperature case ($\eta > 0, T = 0$), or the finite temperature case ($\eta, T > 0$)?

Because the $T = 0$ (no noise) case is already known to feature rigid subharmonic entrainment, in this work we focus on $T > 0$. Furthermore, if a CDTC exists in d_* dimensions, it will presumably exist for all $d \geq d_*$, so we focus on the most delicate case: $T > 0$ in one dimension (1D). While an equilibrium phase transition is impossible in 1D, might there nevertheless be a non-equilibrium dynamical phase transition between a period-doubled CDTC and a symmetric phase?

In this work, we first argue, based on remarkable results due to Gács [48–50] and Toom [51–53], that *in principle* true CDTCs, with an infinite autocorrelation time, are possible in all dimensions $d > 0$ [54]. However, in 1D the construction is so baroque that we cannot yet explicitly prove this conjecture theoretically or numerically. To this end, we instead investigate in detail a more physical Hamiltonian, the parametrically-driven Frenkel-Kontorova (FK) model. The basic idea is simple: each non-linear oscillator in the chain undergoes a 2:1 parametric resonance, and we couple the oscillators together to try and stabilize the CDTC phase at fi-

nite temperature. We find that this model exhibits an intriguing line of first order dynamical phase transitions between an “activated” period-doubled CDTC and a symmetry unbroken phase that terminates at a critical point (Fig. 2). The activated CDTC is not a true time-crystal, rather, it has an autocorrelation time that diverges exponentially as $T \rightarrow 0$. Because of this exponentially diverging timescale, it would be extremely difficult in experiments to distinguish between an activated CDTC and a true, long-range-ordered CDTC; indeed, in order to detect this difference we must conduct careful numerical experiments over many millions of Floquet cycles.

Before diving into the details, let us emphasize the key characteristic of a true long-range-ordered CDTC; namely, the existence of period-doubling with an infinitely long autocorrelation time, τ , which is stable to small perturbations of the dynamics. More precisely, if one considers a model of periodically driven oscillators with position coordinates q_i , the auto-correlation time, τ , of period doubling can be quantified as $\langle q_i(nt_D)q_i(0) \rangle \propto (-1)^n e^{-nt_D/\tau}$, where t_D is the period of the drive and $\langle \rangle$ indicates averaging over noise realizations. For low but finite temperatures, the period doubling we observe in the driven FK model exhibits only an *activated* autocorrelation time, $\tau(T) \sim e^{\Delta_{\text{eff}}/T}$, where Δ_{eff} is an effective activation barrier. Thus, in 1D, the CDTC order of the FK model we study survives only to exponentially long, but not infinite times. Despite the activated behavior of the period doubling, we nevertheless find a first order dynamical phase transition where $\tau(T)$ drops discontinuously and period doubling is completely destroyed (Fig. 2). In 1D, such a first order phase transition would be impossible in equilibrium [55].

This leaves the question of CDTCs at a peculiar point: while we believe a true CDTC is in principle possible in 1D, the obvious candidate model is only exponentially close to one. It may be that the full complexity of Gács’ construction, which we now explain, is in fact required.

Conjecture on the existence of classical time-crystals— While our ultimate interest is to understand the possibility of time-crystals in open Hamiltonian systems, it is first worth considering a more general class of dynamical systems, “probabilistic cellular automata” (PCA). Recall that a deterministic cellular automata (CA) is a set of spins $\{\sigma_i\}$, where $\sigma_i \in \{1, 2, \dots, N\}$, with a discrete update rule $\sigma_i \rightarrow T[\sigma_{i-1}, \sigma_i, \sigma_{i+1}]$ [56]. In a PCA, the system is instead described by a probability distribution $\mathcal{P}[\{\sigma_i\}]$ over spin configurations which is updated by a local Markov process, i.e., each σ_i is updated with a probability distribution that depends only on itself and its neighbors. One way to obtain a PCA is to start with a deterministic CA and perturb it by stochastic errors, e.g., with probability $E \ll 1$, a spin can violate the transition rule and flip into a random state (the precise statement of such an error model can be found in [49]).

Since the PCA update rule is discrete-time translation invariant, one can then ask if the long-time probability distribution can ever oscillate with period two, thereby realizing a time-crystal. For a *deterministic* CA, a time-crystal is triv-

ially obtained by letting binary spins transform under the rule $0 \leftrightarrow 1$ (this is analogous to our previous discussion of fine-tuning, or introducing friction but no noise). However, any infinitesimal rate of errors will desynchronize the spins and the autocorrelations will decay over a time-scale, $\tau \sim 1/E$.

Might there be a more complicated CA which implements “error correction” enabling the system to preserve long-range (time-crystalline) order? The most obvious guess is to have each spin poll its neighbors and transition opposite to the majority, essentially implementing a type of ferromagnetism. This “anti-majority vote” rule, it turns out, is insufficient to preserve the long-range order, because if a pair of domain walls nucleates and separates beyond the range of the vote, they will diffuse randomly throughout the system and destroy the order.

For a very long time it was thought such error correction was in fact impossible in 1D as a consequence of the so-called “positive rates conjecture.” A PCA is said to be “ergodic” if it has a unique fixed-point probability distribution $\mathcal{P}_0[\{\sigma_i\}]$. The positive rates conjecture states that a PCA is generically ergodic, generalizing the physics folklore regarding the impossibility of ferromagnetism in 1D to the non-equilibrium setting [33]. On the other hand, the infinite auto-correlation time of a time-crystal requires ergodicity breaking, because there will be two probability distributions $\mathcal{P}_{\pm}[\{\sigma_i\}]$, corresponding to the two phases ($\sigma_i = \pm 1$) of the period doubling, invariant under a double Markov update. Consequently, were the positive rates conjecture true, it would forbid time-crystals as a corollary.

However, almost two decades ago, a truly remarkable counter-example to the positive rates conjecture was provided by Gács [48–50]. Gács constructed a 1D PCA with the following property: in the thermodynamic limit, there is a threshold error-rate $E < E_*$ below which a system of size L has 2^L distinct steady states in the thermodynamic limit, i.e. it “remembers” one bit per unit cell. Admittedly, the Gács PCA is very complicated: each cell has a state space whose dimension is on the order of $\sigma_i \in \{1, 2, \dots, 2^{400}\}$, which one can think of as 399 ancillary bits which implement an error-correction protocol on one protected bit. The error correction is a highly collective phenomena, requiring communication (interactions) between neighboring cells since there are occasional error events which scramble all constituent 400 bits simultaneously. The error-correction is stable against generic perturbations (for example, the errors can be adversarially biased towards a particular spin configuration, breaking anything like an Ising symmetry), so is truly a feature of a new non-equilibrium *phase*. This is somewhat analogous to MBL, insofar as the ergodicity breaking is extensive and stable to arbitrary perturbations, but is even stronger because the system is open (noisy). Furthermore, the Gács PCA can be used to simulate any deterministic CA acting on the protected subspace. Thus, by using the Gács PCA to simulate the trivial $0 \leftrightarrow 1$ CA, one obtains an example of a true classical time-crystal stable against arbitrary stochastic perturbations!

We conjecture that Gács’ result on PCAs implies the exis-

tence of true CDTCs in time-periodic Hamiltonian Langevin dynamics, because the latter can be used to approximately simulate a discrete PCA. By “simulate,” we mean that if we interpret the integer part of a classical oscillator q_i as the state of a discrete CA, $\sigma_i = \lfloor q_i \rfloor$, then by an appropriate choice of time-dependent Hamiltonian and Langevin friction, the dynamics of q_i under one Floquet-cycle can mimic the discrete update of a CA. Langevin noise then introduces stochastic errors analogous to those of the PCA. Simulating Gács’ PCA in this way would then realize a time-crystal as discussed above, and small errors in the “simulation” would be covered by Gács’ stability result.

However, at present this remains a conjecture, rather than a proof, because *a priori* it is unclear whether the errors arising from Langevin noise can be successfully mapped onto the error model covered in Gács’ result. While we would like to numerically verify this possibility, the Gács model is so complex it is daunting to implement even as a PCA, let alone simulate via continuous-time Langevin dynamics. However, given that we conjecture a CDTC is in principle possible, it is tempting to assume that by writing down a sufficiently general model, we will find this behavior somewhere in its phase diagram.

Period doubling in the Frenkel-Kontorova model—For concreteness, let us consider a parametrically driven Frenkel-Kontorova model, which describes an array of coupled, non-linear pendula (Fig. 2a) [57, 58]:

$$H = \sum_i \frac{1}{2} p_i^2 + [1 + \delta \cos(\omega_D t)] (1 - \cos(q_i)) + g \sum_{\langle i,j \rangle} \frac{(q_i - q_j)^2}{2}, \quad (1)$$

where q_i is the pendulum’s deflection from vertical and p_i its momentum, while δ and ω_D are the amplitude and frequency of the parametric driving (Fig. 2a). Such a model can be realized in many different experimental systems [59]. Note that we have normalized each pendulum’s natural frequency to one. When $\omega_D \sim 2$ and $g = 0$ (decoupled), each pendulum is susceptible to a 2:1 parametric resonance, where the dynamics are period doubled and the pendulum’s position returns only once every two driving cycles. While this subharmonic response (at $\omega_D/2$) is reminiscent of the behavior expected for a CDTC, the parametric resonance of a single oscillator does *not* exhibit the rigidity of a true time-crystal; rather, as we shall see, this sub-harmonic response of a single oscillator is destroyed in a smooth crossover for any amount of noise ($T > 0$). But, crucially, in the presence of interactions ($|g| > 0$), the sub-harmonic response of the *collective* system can undergo a sharp transition at $T > 0$ characteristic of a many-body phase.

We introduce friction η and finite temperature T through a Langevin force $F_B(t)$ which acts independently on each q_i . The pendula then evolve under a combination of this stochastic Langevin force and the ferromagnetic ($g > 0$) Frenkel-

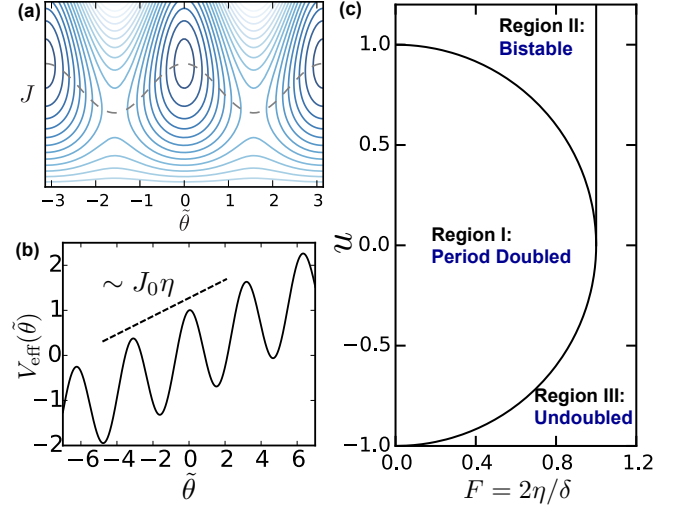


FIG. 3. Parametric resonance of a single non-linear pendulum. **a)** Equal pseudoenergy contours of the averaged Hamiltonian $\bar{H}$ in the θ, J plane. The dashed line indicates the contour $\partial_{\bar{\theta}} \bar{H} = 0$. **b)** Effective washboard potential of Eqn. (10). The slope of the potential arises from the Langevin damping η . **c)** In region I ($F > 1$ or $u < \sqrt{1-F^2}$), the particle slides and there is no period doubling. In Region III, $F < 1, u > 1-F^2$, only the locked (period doubled) phase is stable. In region II, $F < 1, u^2 < 1-F^2$, both the locked and sliding states are stable, implying bistability.

Kontorova Hamiltonian:

$$\frac{dq_i}{dt} = p_i \quad (2)$$

$$\frac{dp_i}{dt} = -(1 + \delta \cos(\omega_D t)) \sin(q_i) + g(q_{i+1} + q_{i-1} - 2q_i) - \eta p_i + \xi_i(t). \quad (3)$$

where $\xi_i(t)$ is a stochastic force with variance $\langle \xi_i(t) \xi_j(0) \rangle = \delta_{ij} \delta(t) 2\eta T$.

To probe the resulting dynamics, at time $t = 0$ we initialize the oscillators in $p_i(0) = q_i(0) = 0$, and integrate the equations of motion using a second-order Langevin time-stepper [59]. The resulting stroboscopic dynamics, $q_i(mt_D)$ (where $t_D = 2\pi/\omega_D$), are depicted in Fig. 1. Strikingly, the uniform initial condition gives way to a growing bubble of spatio-temporal “antiferromagnet” in which $q_i(mt_D) \propto (-1)^{i+m}$; these spatial anti-ferromagnetic correlations are particularly surprising, since the oscillators are ferromagnetically coupled. The existence of a growing bubble would seem to suggest the presence of two distinct dynamical regimes — time crystalline and not — despite the finite temperature fluctuations.

Analysis of a single non-linear pendulum—To begin, let us review the parametric resonance of a single non-linear pendulum [7]. Readers familiar with the resulting phase diagram (Fig. 3), can skip to our discussion of their coupled behavior, though the treatment in action-angle variables is not entirely standard and will prove illuminating [60]. In the action-angle coordinates of the pendulum, $q \sim \sqrt{2J} \cos(\theta), p \sim$

$-\sqrt{2J}\sin(\theta)$, the Hamiltonian [Eqn. (1)] reduces to:

$$H(t) = J - \frac{\epsilon}{2}J^2 + \delta J \cos(\omega_D t) \cos^2(\theta) + \dots \quad (4)$$

where $\epsilon = \frac{1}{8}$ and higher-order terms in J, δ are neglected [61]. In the un-driven case, $\omega(J) = \partial_J H$ sets the frequency of oscillations, so that the non-linearity is encoded in the J^2 term. Because $\epsilon > 0$, larger amplitude oscillations have lower frequency; however, we can be more general by keeping ϵ as a parameter and indeed will later explore the $\epsilon < 0$ regime.

Near a period-doubled solution, $q_i \propto \cos(\omega_D t/2 + \tilde{\theta})$, where $\tilde{\theta}$ varies slowly. To this end, we transform to the rotating frame, $\tilde{\theta} = \theta - \omega_D t/2$, wherein Eqn. (4) becomes

$$H(t) = (1 - \omega_D/2)J - \frac{\epsilon}{2}J^2 + \frac{\delta}{4}J \cos(2\tilde{\theta}) + \frac{\delta}{4}J \left[\cos(2\omega_D t + 2\tilde{\theta}) + 2\cos(\omega_D t) \right]. \quad (5)$$

At leading order in a Floquet-Magnus expansion, the average Hamiltonian over a driving period is given by:

$$\begin{aligned} \bar{H} &= \frac{1}{t_D} \int_0^{t_D} H(t) dt \\ &= \delta(1+u)J/4 - \frac{\epsilon}{2}J^2 + \frac{\delta}{4}J \left(\cos(2\tilde{\theta}) - 1 \right), \end{aligned} \quad (6)$$

where $u = 4(1 - \omega_D/2)/\delta$ is an effective detuning. The static $\bar{H}$ will govern the slow dynamics $\tilde{\theta}$. While this treatment is approximate (i.e. we neglect the off-resonant oscillatory terms), for a single pendulum one can in principle apply a convergent sequence of canonical transformations to bring the Hamiltonian to such a static form, a consequence of the KAM-theorem for small δ [60]. The equal-energy contours of $\bar{H}$ are illustrated in Fig. 3a. For $\epsilon > 0$, the landscape is that of an inverted double well potential with *maxima* at $\bar{J} = \epsilon^{-1}(1+u)\frac{\delta}{4}$ and $\tilde{\theta} = 0, \pi$, which are the two possible phase shifts of the period-doubled solutions: $q(mt_D) = \pm\sqrt{2J}(-1)^m$. Since these solutions occur at maxima, perturbations about the orbit remain bounded, oscillating at an effective frequency, $\omega_{\text{eff}}^2 = \delta^2(1+u)/4$. For weak driving δ , $\omega_{\text{eff}}/\omega_D$ is parametrically small, implying that higher order terms in our Floquet-Magnus expansion are strongly off resonant, justifying our approach. We note that period-doubled solutions exist only when $\bar{J} > 0$, yielding the period-doubling criteria $u > -1$. Finally, for $\epsilon < 0$, the above analysis remains essentially identical except that the period-doubled solutions occur at the *minima* of $\bar{H}$.

The double-well potential has a phase-shift symmetry, $\tilde{\theta} \rightarrow \tilde{\theta} + \pi$, which is unrelated to the $q \rightarrow -q$ symmetry of the original Hamiltonian, and is actually a more general consequence of period doubling: in the rotating frame, $\tilde{\theta} \rightarrow \tilde{\theta} + \pi$ *must* remain a dynamical symmetry because it corresponds to time translation symmetry, $t \rightarrow t + t_D$. In fact, the $q \rightarrow -q$ symmetry of the original Hamiltonian is inessential to the period-doubling physics we discuss, analogous to quantum case [36].

To understand the effect of the bath, consider the equations of motion when averaging the Langevin force over one period [59]:

$$\dot{\tilde{\theta}} = \partial_{\tilde{J}} \bar{H} \quad (8)$$

$$\dot{\tilde{J}} = -\partial_{\tilde{\theta}} \bar{H} - \eta \tilde{J} + \sqrt{\tilde{J}} \xi(t). \quad (9)$$

In the limit of large detuning relative to the driving ($u \gg 1$), these equations take a particularly simple form (see supplementary information [59] for the general case):

$$\bar{H} = -\frac{\epsilon \tilde{J}^2}{2} + \tilde{J} \left[\frac{\delta}{4} \cos(2\tilde{\theta}) + \eta \tilde{\theta} \right] + \dots \quad (10)$$

$$\dot{\tilde{\theta}} = \partial_{\tilde{J}} \bar{H} \quad (11)$$

$$\dot{\tilde{J}} = -\partial_{\tilde{\theta}} \bar{H} - \eta \tilde{J} + \sqrt{\tilde{J}} \xi(t), \quad (12)$$

where $\tilde{J} = J - \bar{J}$ and $\bar{J} = \epsilon^{-1}u\frac{\delta}{4}(1 + \mathcal{O}(u^{-1}))$. These equations describe a *negative* mass particle with “position” $\tilde{\theta}$ and “momentum” $\tilde{J}$ subject to Langevin damping and a washboard potential with finite slope $\tilde{J}\eta$, as shown in Fig. 3(b). We note that such equations of motion are extremely well studied in the context of RC-shunted driven Josephson-junctions [62, 63], where they can give rise to fractional Shapiro steps [21, 22], but our interest will be in their many-body behavior in the presence of noise.

At zero damping, the barrier height of the washboard potential is $\bar{J}\delta/2$, but as the damping $\tilde{J}\eta$ increases the barrier height decreases. Thus, to parameterize the damping we consider a dimensionless “force”, $F = \frac{2\eta}{\delta}$, defined so that when $F \geq 1$ the stable extrema vanish and the particle slides along the washboard. In the fully sliding state, $\dot{\tilde{\theta}} = \frac{\delta}{4}u$, which in the original variables gives $\dot{\theta} = 1$, e.g. the *natural* frequency of the un-driven oscillator [59]. Thus, the “sliding regime” indicates the destruction of period doubling, while the “locked” regime ($\dot{\tilde{\theta}} = 0$) is period doubled.

This point is worth emphasizing. Naively, given that undamped Eqn. (7) is the sine-Gordon representation of the Ising model, it might seem that period doubling is analogous to the breaking of the internal Ising symmetry generated by $\tilde{\theta} \rightarrow \tilde{\theta} \pm \pi$, so that the destruction of the CDTC would occur through something like an Ising transition. But unlike an Ising model, the friction-induced slope gives π -phase slips a handedness (i.e. they in fact have a $\mathbb{Z}$ character rather than a $\mathbb{Z}_2$ Ising character): $\tilde{\theta} \rightarrow \tilde{\theta} + \pi$ is inequivalent to $\tilde{\theta} \rightarrow \tilde{\theta} - \pi$, and the phase prefers to slip one way. The actual dynamical CDTC phase transition we will observe is a collective version of the locked-to-sliding transition at which $\frac{d\tilde{\theta}}{dt}$ jumps, not an Ising transition, and hence is fundamentally distinct from the transition discussed in the context of the quantum MBL / prethermal time crystal [37, 64]. In the supplementary information [59], we demonstrate that this effect is real for the original driven pendulum.

In the absence of noise ($T = 0$), a standard stability analysis [7] reveals three distinct dynamical regimes [Fig. 3(c)]:

In Region I only the sliding regime is stable, indicating there is no period doubling. In Region III only the locked regime is stable, indicating period-doubling. Finally, Region II is a bistable regime in which *both* the locked and sliding states are stable, and the long time behavior depends on the initial state. This region will show hysteresis as F is varied.

At any finite temperature $T > 0$, activated processes cause a *single* pendulum to transition between the locked and sliding states, by analogy to Kramer’s analysis of noise-activated tunneling [65–67]. This destroys the bistable Region II, leading to a smooth (though highly non-linear) crossover between the locked and sliding regimes as either the force (F) or temperature (T) is increased [66]. This crossover in the effective static model is consistent with our numerical experiments on the original parametrically driven pendulum [Eqn. 3] [59]. Thus, for $g = 0$ (i.e. in 0D) and at finite- T there is no sharp transition between period-doubled and undoubled dynamics.

Collective behavior of coupled non-linear pendula—With the single pendulum analysis behind us, let us now turn to the collective behavior of the system at finite coupling strength g .

As before, to analyze the interactions within an effective static model we average the couplings over one Floquet period, $\bar{H}_g = -g \sum_i \sqrt{J_i J_{i+1}} \cos(\tilde{\theta}_i - \tilde{\theta}_{i+1})$ [59], leading to the effective equations of motion:

$$\bar{H} = \sum_i \left(-\frac{\epsilon \tilde{J}_i^2}{2} + \bar{J} \left[\frac{\delta}{4} \cos(2\tilde{\theta}_i) + \eta \tilde{\theta}_i \right] \right. \quad (13)$$

$$\left. - g \bar{J} \cos(\tilde{\theta}_i - \tilde{\theta}_{i+1}) \right) \quad (14)$$

$$\dot{\tilde{\theta}}_i = \partial_{\tilde{J}_i} \bar{H} \quad (15)$$

$$\dot{\tilde{J}}_i = -\partial_{\tilde{\theta}_i} \bar{H} - \eta \tilde{J}_i + \sqrt{\bar{J}} \xi_i(t), \quad (16)$$

where we have taken $J_i \approx \bar{J} + \mathcal{O}(1/u)$. Except for the negative mass and the finite slope $\eta \tilde{\theta}_i$, these equations correspond to the sine-Gordon representation of the Ising model at finite- T .

Before exploring the CDTC phase transition in this model, a few remarks are in order. First, at finite damping and zero-temperature ($\eta > 0, T = 0$), the 1D chain will rather trivially exhibit many-body period doubling which is inherited from the period doubling (Region I) of the single oscillator case. Roughly speaking, if we perturb about the period-doubled solution of a single oscillator $q(t)$, $q_j(t) = q(t) + \Delta q_j(t)$, in the limit $g, \delta \rightarrow 0$ the linearized dispersion relation of the perturbation is strongly off-resonant with the drive, $\omega_{\text{eff}}(k)/\omega_D \sim \delta$, so the finite damping η prevents heating and the system settles into a period-doubled steady state, as can be analyzed from the stability of the linearized equations of motion.

Second, the negative mass explains why the ferromagnetically coupled pendula displayed *anti*-ferromagnetic spatial synchronization in Fig. 1. Because the period doubled orbit of a single pendulum occurs at a maximum of $\bar{H}$, the volume of available phase space increases as the quasi-energy decreases. Interpreting this as a relation between entropy and energy,

the period-doubled solution is at negative temperature, even though the Langevin bath is at positive temperature. Thus, the array is entropically driven toward a *high* quasi-energy state, reversing the expected effect of the coupling g . To test this hypothesis, we can change the effective mass to be positive by changing the non-linearity ϵ of the pendulum (e.g., by replacing $\cos(q) \rightarrow \frac{1}{2}q^2 + \frac{1}{24}q^4$, giving $\epsilon = -\frac{1}{8}$). The period doubled solutions now exist at minima of $\bar{H}$, corresponding to positive temperature, and analogous simulations indeed show that the pendula now synchronize ferromagnetically [59]. While negative temperatures are familiar in models with finite phase space [68–70], the parametric resonance of a non-linear pendulum provides a novel way to dynamically generate negative temperatures in a system with an unbounded phase space. This phenomena is distinct from the dynamically stabilized inverted position of a Kapitza pendulum, which remains at a minimum of the quasi-energy; indeed, ferromagnetically coupled Kapitza pendula will synchronize ferromagnetically [71].

CDTC phase transition—While $\epsilon > 0$ provides access to an intriguing negative temperature regime, from the perspective of time translation symmetry breaking, the sign of ϵ does not appear to impact the CDTC. Thus, for the sake of simplicity and to simplify the visual presentation, we will replace the oscillator potential $\cos(q) \rightarrow \frac{1}{2}q^2 + \frac{1}{24}q^4$ (leading to $\epsilon = -\frac{1}{8}$ and hence, ferromagnetic spatial synchronization) for the remainder of the text.

In the absence of a slope, $\eta \tilde{\theta}_i$, in the washboard potential, the effective model is the sine-Gordon representation of the equilibrium 1D Ising model, implying that at finite temperature, there will always be a finite density $\sim e^{-g_{\text{eff}}/T}$ of π -domain walls in space, as well as a finite rate of π -phase slips $\sim e^{-\Delta_{\text{eff}}/T}$ in time, where $g_{\text{eff}}, \Delta_{\text{eff}}$ are effective quasi-energy barriers. This is the reason why in 1D one naively expects that the CDTC phase will exhibit an “activated” autocorrelation time, $\tau \sim e^{\Delta_{\text{eff}}/T}$, with only a smooth crossover in T as in 0D. However, because of the finite slope $\eta \tilde{\theta}_i$ in the washboard potential, this equilibrium intuition need not apply. Indeed, there is evidence that the 1D DC-driven Frenkel-Kontorova model (DC-driving meaning a finite slope $\eta \tilde{\theta}_i$), shows a first-order locked to sliding transition at finite temperature [58, 72, 73], which would map onto a non-equilibrium phase transition between an activated CDTC and a symmetry unbroken state.

To investigate this possibility numerically, we explore the system’s behavior as a function of both temperature, T , and the dimensionless “force”, F , associated with the washboard slope. To reduce the number of parameters we fix $\eta = 0.005$, set $\delta = 2\eta/F$, and for each F we adjust ω_D, g to keep fixed $\omega_{\text{eff}} = 0.1$ and $g = 1.25\delta$. This choice ensures that throughout the phase diagram we explore $u \gg 1, F < 1$, so that the oscillators are individually in the bistable regime (Region II).

A quantitative diagnostic of period doubling is given by the “velocity” parameter, $v = \langle \dot{\tilde{\theta}} \rangle$, where the average is taken over time, space, and different realizations of the stochastic force. Period doubling corresponds to the locked state where $v = 0$,

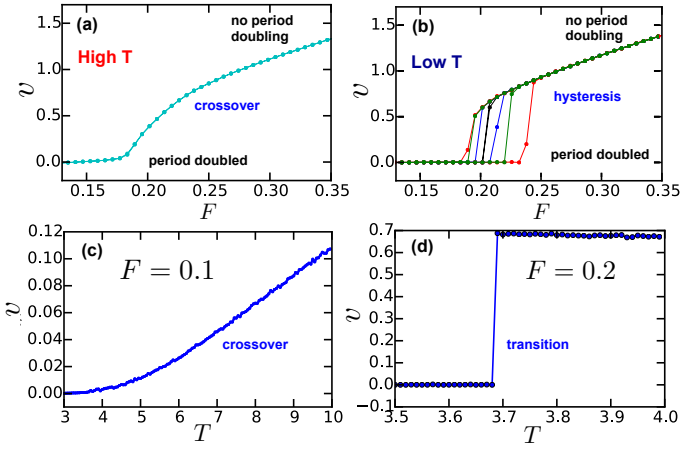


FIG. 4. **Characterizing the CDTC phase transition by measuring the rate of phase slips, $v = \langle \dot{\theta} \rangle$, as a function of damping F and temperature T .** **a)** High temperature ($T = 7$). For low damping F , $v \sim 0$ indicating period doubling, while at large F , $v > 0$ indicates that a finite rate of π -phase slips destroys true long-range order in time. There is a smooth crossover between these two regimes, with no evidence of hysteresis. **b)** At lower temperatures, ($T = 3.7$), we observe a hysteresis loop that closes into a jump discontinuity as we increase the simulation time. Curves of different color correspond to simulations of length $n_{\text{step}} \in (5k, 30k, 300k, 3m)$ for each value of F as we sweep up and down. This suggests a first-order non-equilibrium phase transition. **c)** Analogously, we can hold F fixed and vary T . At $F = 0.1$, v displays a crossover, while at $F = 0.2$, **d)**, there is a sharp transition. Repeating this analysis throughout the F, T plane allows us to trace out a first-order line which terminates, yielding the phase diagram in Fig. 2(b).

while $v \neq 0$ (sliding state) implies a finite rate of phase slips and (strictly speaking) the destruction of true period doubling.

Several representative cuts of $v(F, T)$ in the (F, T) -plane are depicted in Figure 4, which exhibit two regimes. At high temperatures (Fig. 4a), $v(F, T)$ varies smoothly with F and no transition is observed, similar to the zero dimensional case of a single pendulum. However, at low temperatures (Fig. 4b), $v(F, T)$ displays hysteretic behavior as one sweeps the force. By increasing the time scale of each sweep (e.g. increasing the number of driving periods at each F from $n_{\text{steps}} = 5 \times 10^3$ to 3×10^6), the hysteresis loop closes into a single-valued curve that exhibits an apparent jump discontinuity. This suggests that the interactions have transformed the bistable Region II of the individual pendula into a finite-temperature first order dynamical phase transition of the coupled chain! As an additional test, one can fix F to its value near the transition and slowly vary T , reproducing the same discontinuous jump (Fig. 4d).

While the jump looks sharp to the eye, it is difficult to numerically locate the transition in this manner, because the time required to close the hysteresis loop diverges at low temperature. To ameliorate this issue, we study the behavior of a “dynamical” domain wall (DDW) between the period-doubled and un-doubled states, since presumably it is the nucleation of

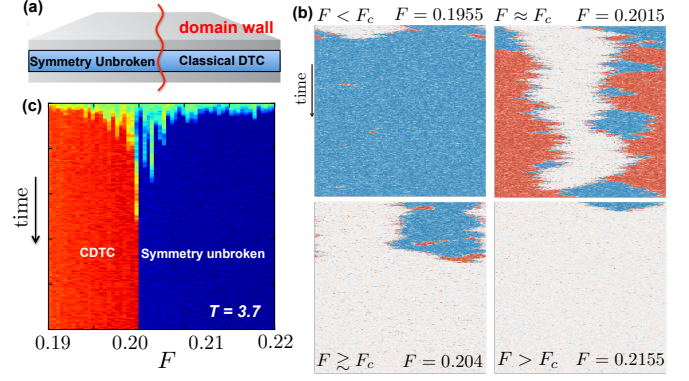


FIG. 5. **Competition between period doubled and undoubled dynamics near the putative first order transition.** By introducing a dynamical domain wall between the two regimes by hand, **a)** we circumvent the exponentially large time scale which leads to hysteresis. Similar to Fig. 1, in **b)** we present $q_j(2nt_D)$ for $n_{\text{osc}} = 10^3$ oscillators. We strobe every *two* driving periods to avoid plotting the period-doubled oscillations. Thus the red and blue regions indicate one or the other period doubled orbits, while in white regions there is no period doubling. At $t = 0$, we initialize a DDW using an initial state in which either $q_j(0) = 0$ (left half) or $q_j(0) \sim \sqrt{2J}$ (right half), and then evolve for $n_{\text{steps}} = 10^6$ periods. If a first order transition exists at critical force F_c , we expect a sensitive dependence of the DDW dynamics on F near the transition. Indeed, for $F < F_c$, the period doubled region expands and “eats” the non-time-crystalline region, while for $F > F_c$, we see the opposite. Meanwhile, close to the transition point ($F_c \sim 0.2015$ for this particular T), the competition between the two phases extends for many steps. To locate the transition precisely, in **c)** we use a color plot to display the time evolution of the average oscillation amplitude $\langle J(t) \rangle_F$ during the DDW quench. Since $\langle J \rangle$ differs in the two regimes, the long time behavior converges to one of two possible values as one or the other domain “wins”.

the first such DDW that requires the largest time. Specifically, we initialize the left half of the system to be in the symmetry unbroken state with $q, p \sim 0$ and the right half of the system to be in the period doubled CDTC with $q, p \sim \sqrt{2J}$ (Fig. 5a). We then time evolve for $n_{\text{steps}} = 10^6$ to determine which state “wins.” In Figure 5(b,c), we fix $T = 3.7$ and repeat this experiment for a very narrow window of F around the putative transition at F_c . For $F < F_c$, we observe the CDTC region expand and “eat” the non-time-crystalline region, while for $F > F_c$, we see the opposite behavior. Meanwhile, close to the critical point, F_c , the competition between the two phases extends for many steps as the location of the DDW fluctuates, indicating coexistence [74].

To quantify this competition between the two domains, we measure the average oscillator amplitude, $\langle J(t) \rangle$, as the system evolves after the quench [Fig. 5(c)]. Since J differs between the CDTC and symmetry unbroken states, its spatial average indicates which domain is winning (although other local observables would serve just as well). Far from F_c , J converges rapidly in time to the value it takes in either the CDTC ($F < F_c$) or symmetry unbroken ($F > F_c$) state, indicating that one or the other domain has taken over. As $F \rightarrow F_c$, the

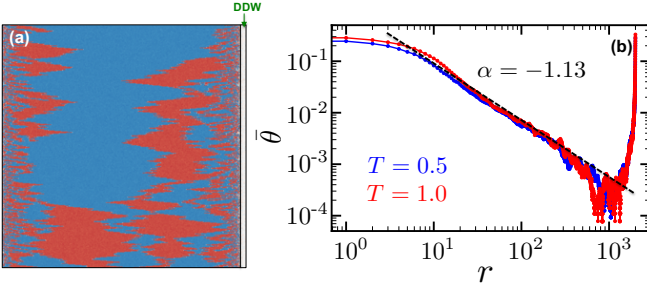


FIG. 6. **Depicts the presence of power law correlations at the first order CDTC transition.** **a)** To probe these correlations, we pin a pair of dynamical domain walls by spatially modulating F above and below F_c (note the small white strip on the right, which is the high- F region), and observe the resulting dynamics $q_j(2nt_D)$. The DDW emits a constant stream of π -domain walls into the CDTC region, visible as boundaries between red and blue domains. Their density is quantified by the average velocity $v(r) = \langle \dot{\theta}(r) \rangle$. **b)** The density of π -domain walls follows a power law $v(r) \sim r^{-\alpha}$ away from the DDW, where $\alpha \approx 1.13$. The exponent appears to be constant over several temperatures (here $T = 0.5, 1$).

time required for convergence increases, and we utilize the long time behavior to accurately determine F_c . One expects that for larger systems, this convergence time scale will diverge, owing to the diffusive dynamics of the DDW, although we have not investigated this quantitatively [74].

By repeating this analysis as a function of F and T , we obtain the CDTC phase diagram depicted in Fig. 2(b). We observe a line of first-order dynamical phase transitions terminating at a point in the (F, T) plane. As expected, in the CDTC region of the phase diagram, the rate of phase slips exhibits activated behavior: $v \sim e^{-\Delta_{\text{eff}}(F)/T}$ [Fig. 2(c)], while at the first order transition, v jumps discontinuously into a regime with *complete* destruction of period doubling.

The nature of the end-point of the first order line is an intriguing question for future study. In Fig. 2(d), we show the magnitude of the jump discontinuity, Δv , across the first order line for a range of temperatures. The magnitude of the jump decreases as we approach the end-point of the first order line, consistent with a scenario in which the phase transition becomes continuous at the critical end-point. Understanding this critical point would be a fruitful starting point for a field-theoretic understanding of the CDTC transition.

If the coarse-grained behavior of the DDWs were governed by an effective free-energy functional with short-range interactions, then entropic arguments would imply that the putative transition is in fact rounded-out to a crossover [55]. However, far from equilibrium, it is unclear that such a free energy based argument has any relevance. Moreover, a non-equilibrium system can generically develop power-law correlations which may mediate power-law interactions between the DDWs. To explore this, we pin a pair of DDWs by spatially modulating F slightly above and below its critical value. As depicted in Fig. 6(a), we find that the DDW boundary between the CDTC and symmetry-unbroken regions emits a finite density of π -

phase slips. These π -phase slips contribute to a finite velocity, $v(r)$, that depends on the distance r from the DDW. We observe $v(r) \sim r^{-\alpha}$ over almost two decades, where $\alpha \sim 1.13$ [Fig. 6(b)]; within the accuracy of our numerics, we obtain the same exponent α for cuts across the transition at two different temperatures $T = 0.5, 1$. This power law behavior is certainly distinct from the expectations for an equilibrium first order transition.

Finally, we will now consider a diagnostic of the CDTC phase which is particularly amenable to experiments [59], namely, the power spectrum of q_j ; in the supplementary information [59], we propose two possible experimental realizations of our model using either a trapped-ion nano-friction simulator or a coupled array of superconducting junctions [75–77]. We define the “stroboscopic” Fourier transform by $q(\omega, k) \equiv \sum_{n,j} (-1)^n e^{i(\omega n - k j)} q_j(nt_D)$. In order to estimate the power spectrum, $S(\omega, k) = \langle q(-\omega, -k)q(\omega, k) \rangle / n_{\text{osc}} n_{\text{steps}}$, we utilize Welch’s method and average over the stochastic noise [78]. A typical spectral function is illustrated in Figure 7(a) and reveals the dispersion relation of the effective Hamiltonian in Eqn. (13).

Perfect ferromagnetic period-doubling would manifest as a δ -function peak at $\omega, k = 0$. A shift in the peak (away from $\omega = 0$) indicates unlocking from the subharmonic response $\omega_D/2$. Since $(-1)^n q(nt_D) \sim \sqrt{2J} \cos(\tilde{\theta}(t))$, this shift is precisely the velocity $v = \langle \dot{\theta} \rangle$ we’ve discussed, while the broadening is analogous to a Debye-Waller factor. In Figure 7(b), we depict $S(\omega, 0)$ for a range of temperatures across the phase transition and observe a qualitative transformation below T_c . As the temperature decreases even further, the spectral function has to be averaged over extremely long times in order to detect the exponentially small shift and broadening of the peak away from $\omega = 0$; nevertheless, it can be used to experimentally detect the first-order jump in v shown in Fig. 4(d).

Discussion—We have demonstrated that a periodically driven, one-dimensional system at finite temperature can exhibit a first-order dynamical phase transition between an activated classical discrete time crystal and a symmetry-unbroken phase. This behavior depends crucially on an interplay between interactions and non-equilibrium driving, without which a transition would be forbidden.

Our work opens the door to a number of intriguing future directions. First, while the 1D CDTC model studied here is thermally activated, our understanding of Gács’ result suggests that true CDTCs are in principle possible. If so, is there a simple “physical” model, like the driven-FK chain studied here, which exhibits a true CDTC, or is the whole complexity of the Gács result required?

In $d > 1$, where ergodicity can certainly be broken without requiring the full complexity of Gács’ result (for example in Toom’s model [51]), the possibility of a true CDTC is still complex. This issue was considered in the context of certain 2D PCAs, where it has been argued that a higher-order subharmonic response (e.g. periodic-tripling, $k = 3$) cannot have an infinite auto-correlation time, while a period-doubled

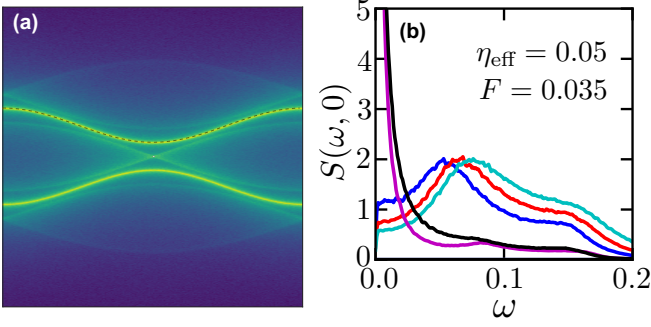


FIG. 7. Probing the CDTC using the stroboscopic spectral function. **a)** The stroboscopic spectral function $S(\omega, k) \propto \langle |\text{DFT}[(-1)^n q_j(nt_D)]|^2 \rangle$. In this convention, the period doubled component is mapped to $\omega = k = 0$. Since the autocorrelation time is unmeasurable over the $n_{\text{steps}} = 15000$ used to take the data, there is a δ -function peak at the origin we have removed by hand to preserve the scale. The residual noise spectrum reveals the mode $\omega_{\text{eff}}(k)$ of the effective Floquet-Magnus Hamiltonian Eqn. (13). **b)** $S(\omega, 0)$ as a function of frequency at several temperatures, holding the other parameters (e.g., F, η, g) fixed. The lowest temperatures (purple, black) lie below the first-order transition, $T < T_c$ and the peaks (if resolved) would exhibit a very small shift away from $\omega = 0$ due to the exponentially-rare activated phase slips. For the remaining curves $T > T_c$, and the peaks are strongly shifted away from $\omega = 0$ indicating unlocking of the subharmonic response.

response could [54]. The basic argument is that if a fluctuation nucleates a bubble of the neighboring subharmonic orbit (the generalization of our π -phase slips), the domain wall will generically experience a force which causes the bubble to expand, destroying the CDTC. They argue that this force is absent for $k = 2$ because an Ising-like domain wall does not have an orientation, leading to the distinction between $k = 2$ and $k > 2$.

Our mapping to a tilted sine-Gordon model provides a new perspective on their analysis. For a k -th order subharmonic response, as would occur for a $k : 1$ parametric resonance, one can generalize our effective Hamiltonian by simply replacing the period of the washboard potential with $\cos(k\theta_j)$. As we have emphasized, when the domain wall is smooth it still has a definite handedness even when $k = 2$, so the distinction between $k = 2$ and $k > 2$ no longer seems so important. Indeed, in the continuum limit, k drops out of the resulting sine-Gordon model. For a tilted sine-Gordon model satisfying detailed balance, the “locked” phase, with true long-range order at the subharmonic period, is unstable in any dimension: i.e. for any non-zero tilt there is only a finite energy barrier to create a phase slip domain, which, once nucleated, grows rapidly to infinite size due to the tilt. However, nucleation could in principle be prevented by non-equilibrium effects. Furthermore, the locking to a commensurate frequency and the existence of spontaneous oscillations are potentially distinct: even the “sliding” phase may exhibit long-range or quasi-long-range [79] order at a frequency *shifted* from $\omega_D/2$. Such “incommensurate” temporal order was discussed in the context of driven periodic media [20].

Furthermore, while the handedness (and hence force) on domain walls is well defined in the sine-Gordon continuum limit, the lattice allows for 2π -vortices at which the orientation of the domain wall reverses. These vortices cost finite energy when localized to the domain wall, where they may proliferate due to fluctuations. This may provide a mechanism which renormalizes the force on the domain walls to zero, stabilizing a true CDTC for $k = 2$ and $d > 1$.

Another interesting scenario is to consider the $\eta, T \rightarrow 0$ limit of our model, which reduces to closed Hamiltonian dynamics. Here, we observe that the dynamics remain period-doubled out to extremely long time scales (e.g. many millions of driving cycles). This behavior appears to be the classical analog [80, 81] of a “prethermal” time crystal and arises from a mismatch between the driving frequency ω_D and the *effective* resonance frequency ω_{eff} [31, 64, 80–82].

The existence of a first-order line terminating at a critical point is reminiscent of the equilibrium liquid-gas transition. To this end, at the critical point, one might hope to develop a non-equilibrium field theory for the transition within the Martin-Siggia-Rose path integral formalism [83, 84]. Such a field theory might also be used to determine the precise relation between classical and quantum time-crystals by analyzing the semiclassical limit of its Keldysh path integral.

Finally, we comment on the relation of CDTC order to synchronization transitions, which arise, for example, in the Kuramoto model [12]. Adding a weak external periodic drive to a synchronized model, at a frequency which is a multiple of the synchronized one, presumably converts the synchronized phase into a rigid subharmonic response [1]. However, this does not mean the Kuramoto model is an example of a CDTC. First, the Kuramoto model is usually understood as an all-to-all model, which is non-local, so one needs to begin by introducing a geometry to the couplings (chain, array, etc.). Second, the usual Kuramoto model is dissipative, for which we have emphasized there are already many examples of rigid subharmonic responses. So one should either (a) consider the Hamiltonian realization of the Kuramoto model, with no friction [85] (the closed case) or (b) couple the model to spatio-temporal noise [86] e.g. a Langevin force (the open case). The stability of synchronization, in the sense of an infinite autocorrelation time, in a *local* model, is to our knowledge unknown for both the closed and open cases [86].

Acknowledgments—We gratefully acknowledge the insights of and discussions with E. Altman, D. Huse, S. Gazit, R. Goldstein, L. Sieberer, S. Sondhi, and B. Zhu. This work was supported in part, by the DARPA DRINQS program (D18AC00033), the David and Lucile Packard Foundation and the W. M. Keck Foundation. L.B. was supported by the NSF Materials Theory program through grant DMR1506119.

Contributions—All authors contributed extensively to all aspects of this work.

Data Availability Statement—The data represented in Figs. 2b-d, 4a-d, 6b and 7b are available as Source Data 2b-d, 4a-d, 6b and 7b. All other data that support the plots within this paper and other findings of this study are available from the

corresponding author upon reasonable request.

Competing interests—The authors declare no competing financial interests.

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