Aspect of Floquet physics in closed quantum systems

K. Sengupta

IACS, Kolkata

Outline

Driven systems: Thermalization

Prethermal Physics: emergent symmetry

Outline
Driven systems: Thermalization
Prethermal Physics: emergent symmetry
Pretehrmal Hilbert space fragmentation
Arresting heating: a brief introduction

Arresting heating: a brief introduction

Two-rate periodic protocol

Future directions.

Consider a generic state of quantum non-integrable many-body system $\frac{1}{m}$

$$
|\psi(t)\rangle = \sum_{m} C_m e^{-iE_m t} |m\rangle,
$$

D'Alessio et. al
Adv. Phys. 65, 239 (2016)
. Adv. Phys. 65, 239 (2016)

The time evolution of a generic operator for this state is given by

there is a of quantum
\n*blue many-body system*
\nThe time evolution of a generic operator for this state is given by
\n
$$
O(t) \equiv \langle \psi(t)|\hat{O}|\psi(t)\rangle = \sum_{m,n} C_m^* C_n e^{i(E_m - E_n)t} O_{mn}
$$
\n
$$
O_{mn} = \langle m|\hat{O}|n\rangle.
$$
\n
$$
= \sum_m |C_m|^2 O_{mm} + \sum_{m,n \neq m} C_m^* C_n e^{i(E_m - E_n)t} O_{mn}
$$
\n*l*\n*l*\n*ssues with long-time behavior:*
\n**a)** The steady state value of *O(t)* depends on the overlap coefficients: no thermalization
\n*(in the sense that the value does not agree with standard ME prediction)*
\n**a)** It takes an incredibly long time to reach the steady state (predicts a very large relaxation time).
\nInvolking random matrix theory remedies these problems since within RMT
\n
$$
O = O' \text{ and } O = 0. However it provides an energy independent answer
$$

Issues with long-time behavior:

- (in the sense that the value does not agree with standard ME prediction)
- relaxation time).

Invoking random matrix theory remedies these problems since within RMT O_{mm} = O' and O_{mn} =0. However it provides an energy independent answer which does not agree with standard numerical results.

Generalization of the RMT result for the matrix elements of a "typical" operator

Eigenstate Thermalization Hypothesis
Generalization of the RMT result for the matrix elements of a "typical" operator

$$
O_{mn} = O(\bar{E})\delta_{mn} + e^{-S(\bar{E})/2} f_O(\bar{E}, \omega) R_{mn}, \qquad \bar{E} \equiv (\underline{E}_m + \underline{E}_n)/2,
$$

Both O and $f_{\rm O}$ are smooth functions of their arguments, S is the entropy, and R is a gaussian random number.

It states that for a large-enough system, the answer is nearly identical to that

Eigenstate Thermalization Hypothesis
\nGeneralization of the RMT result for the matrix elements of a "typical" operator
\n
$$
O_{mn} = O(\bar{E})\delta_{mn} + e^{-S(\bar{E})/2}f_O(\bar{E}, \omega)R_{mn}, \qquad \bar{E} \equiv (E_m + E_n)/2,
$$
\nBoth 0 and f_O are smooth functions of their arguments, S is the entropy, and R is a gaussian random number.
\nIt states that for a large-enough system, the answer is nearly identical to that obtained using a microcanonical ensemble at the average energy.
\n
$$
\bar{O} \equiv \lim_{t_0 \to \infty} \frac{1}{t_0} \int_0^{t_0} dt O(t) = \sum_m |C_m|^2 O_{mm} = \text{Tr}[\hat{\rho}_{DE}\hat{O}], \qquad O_{ME} = \text{Tr}[\hat{\rho}_{ME}\hat{O}]
$$
\n
$$
\bar{O} \simeq O(\langle E \rangle) \simeq O_{ME}.
$$

This relies on the fact that energy fluctuations in a many-body system are subextensive.

$$
O_{mm} \approx O(\langle E \rangle) + (E_m - \langle E \rangle) \frac{dO}{dE} \Big|_{\langle E \rangle} + \frac{1}{2} (E_m - \langle E \rangle)^2 \frac{d^2O}{dE^2} \Big|_{\langle E \rangle},
$$

$$
\overline{O} \approx O(\langle E \rangle) + \frac{1}{2} (\delta E)^2 O''(\langle E \rangle) \approx O_{\text{ME}} + \frac{1}{2} [(\delta E)^2 - (\delta E_{\text{ME}})^2] O''(\langle E \rangle),
$$

Floquet version: ETH in a periodically driven system
namics is completely determined by the Floquet Hamiltonian: U(nT,0)= exp[- i H_F nT] Floquet version: ETH in a periodically driven system
Periodic drive: Stroboscopic dynamics is completely determined by the Floquet Hamiltonian: U(nT,0)= exp[- i H_F nT]
For low or intermediate drive frequency, H_F can no Floquet version: ETH in a periodically driven system
Periodic drive: Stroboscopic dynamics is completely determined by the Floquet Hamiltonian: $U(nT,0)$ = exp[- i H_r nT]
For low or intermediate drive frequency, H_r can

For low or intermediate drive frequency, H_F can not be typically described by a short range Hamiltonian

At long times, the system heats up to infinite temperature: rapid growth of entanglement entropy to its maximum value. At long times, the system heats up to infinite temperature: rapid growth of entanglement entropy to its maximum value
However, at high drive frequencies the thermalization timescale are exponentially large [T. Mori et al,

However, at high drive frequencies the thermalization timescale are exponentially large [T. Mori et al, PRL 2016]

This makes prethermal regime experimentally relevant.

These regimes show several phenomena which do not have equilibrium analogue.

Prethermal physics and emergent symmetry

Emergent symmetry Banerjee et al arXiv 2024

Most of these prethermal phenomenon can be understood as a consequence of an approximate emergent symmetry

computed in some approximation scheme commutes with some operator O at special drive frequencies.

controls the dynamics, <O> is approximately conserved at these frequencies. This happens till a prethermal timescale \sim exp[c ω_d]

For time crystals, the argument requires a slight modification.

The first order Floquet Hamiltonian, As long as the first order Floquet Hamiltonian Example of driven Ising model. Jordan Wigner Banerjee et al arXiv 2024

$$
H_F^{(1)} = 2 \sum_{k \in BZ/2} \left[\tau_z (h_s - \cos ka + i\gamma) + \tau_x J_0(\mu) \sin ka \right]
$$

$$
H_F^{(2)} = \sum_k \left[-4\tau_z \sin^2 ka \sum_{n=0}^{\infty} \frac{J_0(\mu) J_{2n+1}(\mu)}{(2n+1)\hbar\omega_D} \right.
$$

$$
+ \tau_x \sin ka(h_s - \cos ka + i\gamma) \sum_{n=0}^{\infty} \frac{4J_{2n+1}(\mu)}{(2n+1)\hbar\omega_D}.
$$

At special frequencies $H_{\digamma}^{(1)}$ commutes with the magnetization

$$
M_z = \sum_j \sigma_j^z
$$

Dynamical freezing of magnetization A. Das, PRB 2010.

Strong Hilbert space fragmentation

Hilbert space fragmentation: Introduction

Breakdown of the Hilbert space into an exponentially large number of dynamically disconnected sectors.

The fragmentation is usually observed in the Hilbert space fragmentation: Introduction

Breakdown of the Hilbert space into an exponentially

large number of dynamically disconnected sectors.

The fragmentation is usually observed in the

computational basis; classi as number basis states $\|n_1,n_2....n_{j}...\,n_l\}$

Such a separation of Hilbert space in dynamically disconnected sectors is different from those due to global symmetries; in the latter case number of sectors scale algebraically with L.

For strong Hilbert space fragmentation (HSF), with n being the largest fragment and N being the total Hilbert space dimensions, $n/N \sim e^{-L}$

Most of the model exhibiting strong HSF are 1D models; More recently a few higher dimensional models have been put forth (see for example, Scipost Phys. 14, 146 (2023)).

Signatures of strong HSF

- 1. Memory retainment leading to finite value of the autocorrelation function at long times
- its symmetry resolved Page value.

$$
\begin{aligned}\n\leftarrow & \circ \circ \circ \circ \circ \circ \rangle \rightarrow \leftarrow & \circ \bullet \circ \circ \circ \circ \circ \circ \rangle \\
\rightarrow \leftarrow & \circ \bullet \circ \bullet \circ \circ \circ \circ \rangle \rightarrow \leftarrow & \circ \circ \bullet \circ \circ \circ \bullet \rangle \rightarrow \dots\n\end{aligned}
$$

The parity of the starting defect sites are conserved.

Counting of fragmentation

Counting of fragmento

A: Number of states in the largest symmetry sector:

We need to fill L sites with L/2 particles and L/4 bonds

To do this first fill L/2 sites with particles keeping L/2 bonds We need to fill L sites with L/2 particles and L/4 bonds

To do this first fill L/2 sites with particles keeping L/2 bonds

Next we insert L/4 empty sites so as to break L/4 of these L/2 bonds. This can be done in $L/2C_{L/4}$ ways

Next, one needs to insert L/4 more empty sites without breaking any bonds. This can be done in $L^{1/2-1}C_{L/4}$ ways.

Finally one gets a factor of 2 due to particle-hole inversion

$$
N_1 = 2^{1/2-1}C_{1/4}^{1/2}C_{1/4} = {^{1/2}C_{1/4}}^2 \sim 2^{1/2}
$$

Example for L=8

Dimension of the largest fragment

Start from an initial state for which N_b = L/4 such $($ that there are N_a= L/4 particle and hole defects. \blacksquare This also leaves $P = L/4-1$ particle-hole pairs.

For half-filling one has $P = L/2 - (N_d + 1)$

Starting from this initial state the action of H leads to diffusion of particle from R to the sequence of hole defects. Such diffusion needs an accompanying hole motion and thus leads to pair movement.

The number of possible configurations obtained by action of H is the number of ways P pairs can be distributed among 2 N_d empty space between particle and hole defects.

Total number of such configurations is $N_f = L N_0 / 2$.

For large L and L/4 particle and hole defects, one has $N_t/N_{1} \sim (0.8)^L$

Dimension of the largest fragment

\n

N_b = L/4 such	$(p, p, ..., h, h, ..., R, R, ...)$				
hole defects.	Na times	N _d times	P times		
+	→	→	→	→	
+	→	→	→	→	
+	→	→	→	→	
+	→	→	→	→	→

\nFrom of H leads sequence of

\n

+	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	→	
---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	---	--

$$
|\bullet\bullet\bullet\bullet\circ\circ\circ\circ\bullet\bullet\circ\bullet\rangle\rightarrow|\bullet\bullet\bullet\bullet\circ\circ\circ\bullet\bullet\circ\circ\circ\rangle.
$$

Prethermal signature of strong fragmentation

Driving a spinless fermion chain: Numerics

We start from a fermion chain and drive the interaction term

For a square pulse protocol and the state of the state of the state \sim For a continuous protocol $[V(t) = -(+)V_1$ for $t \leq (>)T/2]$

The time evolution operator

Can be expressed in terms of eigenvalues and eigenfunctions of H_+ and H_-

$$
U^s(T,0)=\sum_{m,n}c_{mn}^{+-}e^{-i(\epsilon_m^++\epsilon_n^-)T/2\hbar}|\xi_m^+\rangle\langle\xi_n^-|\ \ c_{mn}^{+-}=\langle\xi_m^+|\xi_n^-\rangle.
$$

$$
H_0(t) = V(t) \sum_{j=1..L} \hat{n}_j \hat{n}_{j+1}
$$

\n
$$
H_1 = \sum_{j=1..L} -J(c_j^{\dagger} c_{j+1} + \text{H.c.}) + \hat{n}_j (V_0 \hat{n}_{j+1} + V_2 \hat{n}_{j+2})
$$

\nFor a continuous protocol

The time evolution operator
 $= e^{-iH+T/2h}e^{-iH-T/2h}$
 $= e^{-iH+T/2h}e^{-iH-T/2h}$
 Compute the exact floquet Hamiltonian for the system
 U can be expressed in terms of eigenvalues
 U can be expressed in terms of eigenva The time evolution operator requires Suzuki-Trotter decomposition of U $U = \Pi_j \exp[-i H_j \Delta t / h]$ $\Delta t = T/N$ $\begin{aligned} &\text{H.c.}) + \hat{n}_j (V_0 \hat{n}_{j+1} + V_2 \hat{n}_{j+2}) \ &\text{inuous protocol} \ & = V_1 \cos \omega_D t \ &\text{olution operator requires} \ &\text{er decomposition of } U \ &\text{exp}[\text{-i } \textsf{H}_\text{j} \, \Delta \textsf{t}/\textsf{h}] \qquad \Delta \textsf{t} = \textsf{T}/\textsf{N} \ &\text{g} \ &\text{pressed in terms of eigenvalues} \ \text{in terms of eigenvalues} \end{aligned}$ U can be expressed in terms of eigenvalues and eigenfunctions of H_j

$$
U^{c}(T,0) = \prod_{j=1..N} \sum_{n} e^{-i\epsilon_n^j T/\hbar} |\xi_n^j\rangle\langle\xi_n^j|
$$

We consider a Hamiltonian H(t)= $H_o(t)$ + V(t) and construct the evolution operator $\bm{\mathsf{U}}_o$ corresponding to the largest term of the Hamiltonian [H $_{\mathrm{0}}$ (t)] $\hskip1cm$

Perturbative Analytics: Floquet perturbation theory

\n
$$
H_0(t) + V(t)
$$
 and

\n
$$
U_0
$$
 corresponding

\n
$$
i\hbar \frac{\partial U_0(t,0)}{\partial t} = H_0(t)U_0(t,0).
$$

Next, we construct states in the interaction picture and construct the corresponding Schrodinger equation

$$
\psi^I(t) = U_0(0, t)\psi(t). \quad i\hbar \frac{\partial \psi^I}{\partial t} = V^I(t)\psi^I(t),
$$

$$
V^I(t) = U_0(0, t)VU_0(t, 0).
$$

The evolution operator in the interaction picture reads U^I has the solution

$$
U^{I}(t,0) = \mathcal{T}e^{-(i/\hbar)\int_{0}^{t}dt'V^{I}(t')}, \quad i\hbar \frac{\partial U^{I}(t,0)}{\partial t} = V^{I}(t)U^{I}(t,0).
$$

The perturbative evolution operator is given by $U(t,0) = U_0(t,0)U^I(t,0)$.

The method reduces to the usual rotating wave approximation when the drive term is the one with largest amplitude

$$
U^{I}(t,0) = I + \left(\frac{-i}{\hbar}\right) \int_{0}^{t} dt' V^{I}(t')
$$

+ $\left(\frac{-i}{\hbar}\right)^{2} \int_{0}^{t} dt_{1} V^{I}(t_{1}) \int_{0}^{t_{1}} dt_{2} V^{I}(t_{2}) + ...$

Driven Fermi chain

Consider a chain of spinless fermions with nearest neighbor hopping and density-density interactions

$$
H = -J \sum_{j} \left(c_j^{\dagger} c_{j+1} + \text{h.c.} \right) + V_1 \sum_{j} \hat{n}_j \hat{n}_{j+1}
$$

+
$$
V_2 \sum_{j} \hat{n}_j \hat{n}_{j+2}
$$

$$
V_1(t) = V_0 + V_1 \cos \omega_0 t \qquad \text{for cosine drive}
$$

$$
V_1(t) = V_0 + (-) V_1 \text{ for } t \leq -(-1) \text{ T/2} \qquad \text{for square pulse}
$$

, *J*, one can obtain the Floquet Hamiltonian using FPT
+
$$
A_j = (\hat{n}_{j+2} - \hat{n}_{j-1}), \ \gamma_0 = V_1 T/(4\hbar)
$$

We drive the chain by making V $_{1}$ ==V $_{1}$ (t) a periodic function of time characterized by an amplitude V $_1$ and frequency $\omega_{\textup{D}}$ = 2 π /T, $\hskip1cm$ $\hskip1cm$ $\hskip1cm$ Where T is the time period of the drive

 $V_1(t) = V_0 + V_1 \cos \omega_0 t$ for c $\cos\,\omega_{\rm p}$ t for cosine drive $V_1(t) = V_0 + (-) V_1$ for $t \leq (-) T/2$ for square pulse

In the high drive amplitude regime, V $_1$ >>V $_{\rm o}$, V $_{\rm 2}$, J, one can obtain the Floq

$$
H_F^{(1)} = \sum_{j=1..L} \hat{n}_j (V_0 \hat{n}_{j+1} + V_2 \hat{n}_{j+2})
$$

\n
$$
- J \sum_{i=1..L} [(1 - \hat{A}_j^2) + f(\gamma_0) \hat{A}_j^2] c_j^{\dagger} c_{j+1} + \text{H.c.},
$$

\n
$$
\hat{A}_j = (\hat{n}_{j+2} - \hat{n}_{j-1}), \ \gamma_0 = V_1 T / (4\hbar)
$$

 $f(\gamma_0) = J_0[2\gamma_0/\pi]$ Cosine protocol $f(\gamma_0) = \gamma_0^{-1} \sin \gamma_0 \exp[i\gamma_0 A_j]$ Square Pulse protocol

Realization of a Hamiltonian hosting HSF within Realization of a Hamiltonian hosting HSF within
first order Floquet at frequencies for which f(T)=0
 $H_F^{(1)} = \sum_{j=1...L} \hat{n}_j (V_0 \hat{n}_{j+1} + V_2 \hat{n}_{j+2})$

$$
H_F^{(1)} = \sum_{j=1..L} \hat{n}_j (V_0 \hat{n}_{j+1} + V_2 \hat{n}_{j+2})
$$

- $J \sum_{j=1..L} [(1 - \hat{A}_j^2) c_j^{\dagger} c_{j+1} + \text{H.c.},$

Higher order terms is expected to destroy the fragmentation.

However for large drive amplitudes such terms are small.

This leads to a large prethermal region where signatures of HSF can be seen.

The extent of the prethermal regime showing signatures of HSF can be controlled by tuning drive parameters

Entanglement entropy of the driven chain: Square-pulse protocol

$$
S_p = \ln D_{system} - 1/2
$$

 S_p^f = In D_{fragment} -1/2 Foton of

(a bit more complicated for symmetry-resolved sectors)

Entanglement entropy saturates to the Page value of the sector (S $_{\rho}^{~f})$ instead of $\hskip1cm$ Initial s that of the system (S_p) at special drive frequencies for an exponentially large prethermal timescale

.Entanglement entropy saturates to an Initial state dependent value.

Signature of prethermal HSF.

Fermion density-density autocorrelators: Square pulse protocol

Near the threshold value, the extent of the prethermal regime grows exponentially with drive amplitude

FIG. 3: (a) Plot of $\chi_1(nT)$ as a function of n for $V_1 = 19$ at ω_1^* (blue curve) and $6\omega_1^*$ (red curve) starting from a random frozen state showing lack of ETH predicted thermalization at ω_1^* . (b) Plot of $\langle \chi_1 \rangle$ as a function of V_1 at ω_1^* ; $\langle \chi_1 \rangle$ stays close to its initial value for large V_1 which is consistent with prethermal HSF. (c) Same as in (a) but for initial $|Z_2\rangle$ state showing slow oscillations at ω_1^* . (d) Schematic diagram for the Floquet quasienergies showing doubly degenerate $|Z_2\rangle$ and $|\bar{Z}_2\rangle$ with $N_d = 0$ and other states with $N_d \neq 0$. The arrows indicate transition to $|Z_2\rangle$ from $|Z_2\rangle$ using intermediate states with $N_d \neq 0$ leading to slow oscillations. For all plots $V_0 =$

Dynamics of the frozen state: Continuous protocol $\chi_j(nT) = \langle \psi_f(nT) | \hat{n}_j \hat{n}_{j+2} | \psi_f(nT) \rangle$

Oscillatory dynamics of frozen states due to residual terms in H_F beyond $H_F^{(1)}$

This requires Z $_2$ symmetry. Two states with N $_{\sf d}$ =0 which are eigenstates of H $_{\mathsf{F}}^{-1}$ with same quasienergy.

In addition, it requires fragmentation so that starting from the Z₂ state (which correspond to N_d=0), the system does not spread out in Hilbert space; the dynamics receives most significant contribution from states with $N_d = 1$.

Since χ_i =0 (1) for Z₂ and Z'₂, the oscillations occur between 0 and 1.

The oscillation time scale is determined by higher-order terms in H_F and is the energy split between bonding and antibonding states due to tunneling to $N_d = 1$ sector.

$$
|\psi_{B,A}\rangle \equiv |\mathbb{Z}_2\rangle \pm |\mathbb{Z}_2\rangle. \quad H_F|\psi_{B,A}\rangle \approx \hbar(\alpha_s \pm \alpha_d)|\psi_{B,A}\rangle
$$

$$
\chi_1(nT) \approx \sin^2(\alpha_d nT)
$$

 $10V_2 = 2, L = 14$, and all energies are scaled in units of J.
 This dynamics of frozen states has no analogue in standard HSF in equilibrium

Arresting heating and a two-rate protocol

Arresting heating during a periodic drive

The unbounded growth of entanglement in a driven system

The system reaches and infinite temperature steady state: "heat death" Any information contained in an initial state gets completely scrambled due to rapid spread of the state in the Hilbert in the presence of the drive

Is it possible to arrest this growth?

Counter-diabatic driving Counter-diabatic driv

Unfortunately both of these are difficult to implement experimentally for a generic ergodic quantum systems

Our suggestion: Use of a two-rate protocol which is experimentally more viable

A class of protocols and exact Floquet flat bands

Consider a generic non-integrable driven Hamiltonian characterized non-commuting operators O_1 and O_2

$$
H(t) = \sum_{i=1,2} \lambda_i(t) \hat{O}_i.
$$
 $[\hat{O}_1, \hat{O}_2] \neq 0.$

We choose a class of two-rate drive protocols with time periods T_1 and T_2

 $T_i = 2\pi/\Omega_i$ with $i = 1, 2, \Omega_2 = \nu \Omega_1$

Familiar examples of such class of protocols with $\pmb{\nu}$ = 3 w,λ

$$
\lambda_1(t) = \lambda_0 \cos \Omega_1 t, \quad \lambda_2(t) = w_0 + w_1 \cos \nu \Omega_1 t
$$

$$
\lambda_1(t) = +(-)\lambda_0 \quad \text{for } t \le (>)T_1/2
$$
\n
$$
\lambda_2(t) = w_0 - [+]w_1 \text{ for } \frac{(m-1)[m]T_1}{2\nu} \le t < \frac{m[(m+1)]T_1}{2\nu}
$$

A simple example: Square pulse protocol with $v = 2$

The evolution operator can be written as $U(T,0) = e^{-iH_0[-1,-1]T_1/4} e^{-iH_0[-1,1]T_1/4} e^{-iH_0[1,-1]T_1/4} e^{-iH_0[1,1]T_1/4} = I$

For the square-pulse protocol, this work for any integer v ; for the cosine protocol, this requires an odd integer $v=2p+1$ The method works for any non-commuting operators O_1 and O_2 but requires two rates. It constitutes exact dynamical localization/freezing in an otherwise ergodic many-body system.

Properties of the chosen protocol for $w_0=0$

They exhibit turning points at $t_i = \beta_i T_1$

Between any two turning points β_j and β_{j+1} there exist a point $\alpha_j = (\beta_j + \beta_{j+1})/2$ such $H(\alpha_j T_1) = 0$. that $\lambda_i(\alpha_iT_1)$ = 0 for i =1,2. This ensures **Properties of the chosen protocol for w₀=0**
 t turning points at t_j = $\beta_j T_1$
 y two turning points β_j and β_{j+1}
 y point $\alpha_j = (\beta_j + \beta_{j+1})/2$ such
 j = 0 *for* i =1,2. This ensures
 $= (\beta_{j+1} - \beta_{j$

For any $t_0 \leq (\beta_{j+1} - \beta_{j+1})T_1/2$, H(t) satisfies $H(\alpha_j T_1 + t_0) =$

The evolution operators can therefore be written as, using Suzuki-Trotter decomposition

$$
U(T_1,0) = \prod_{k=0...N_0} e^{-iH(t_k)\Delta t/\hbar} = \prod_k U(t_{k+1},t_k) = \prod_k U_k.
$$
 $\Delta t = T_1/(N_0 +$

This product can be reorganized by grouping \bm{U}_k s between turning points

The evolution operators can therefore be
\nwritten as, using Suzuki-Trotter decomposition
\n
$$
U(T_1,0) = \prod_{k=0...N_0} e^{-iH(t_k)\Delta t/\hbar} = \prod_k U(t_{k+1},t_k) = \prod_k U_k, \qquad \Delta t = T_1/(N_0+1)
$$
\nThis product can be reorganized by grouping U_k s between turning points
\n
$$
U(T_1,0) = \prod_{j=j_{\text{max}}} U(\beta_{j+1}T_1,\beta_{j+1}T_1 - \Delta t)U(\beta_{j+1}T_1 - \Delta t,\beta_{j+1}T_1 - 2\Delta t)...U(\alpha_jT_1 + 2\Delta t,\alpha_jT_1 + \Delta t)
$$
\n
$$
\times U(\alpha_jT_1,\alpha_jT_1 - \Delta t)...U(\beta_jT_1 + 2\Delta t,\beta_jT_1 + \Delta t)U(\beta_jT_1 + \Delta t,\beta_j)
$$
\n
$$
U(T_1,0) = \text{for such protocol leading to } E_n^F(T_1) = 0 \text{ for all } n
$$
\n
$$
E \text{xact Floquet flat bands for all drive frequencies}
$$

U(T₁,0)=I for such protocol leading to E_n ^F(T₁)=0 for all n

Model for studying two-rate dynamics

Rydberg atom arrays

Effective low-energy description

d to their
\n
$$
\frac{\mathcal{H}}{\hbar} = \sum_{i} \frac{\Omega_{i}}{2} \sigma_{x}^{i} - \sum_{i} \Delta_{i} n_{i} + \sum_{i < j} V_{ij} n_{i} n_{j},
$$
\n
$$
\mathbf{which allows
$$
\n
$$
\mathbf{where \textit{g} \textit{v} the time of the time, i.e.,}
$$
\n
$$
\mathbf{v}_{0} = (1+\sigma^{2})/2
$$
\n
$$
\mathbf{v}_{ij} = \mathbf{v}_{0}/|\mathbf{r}_{ij}|^{6}
$$
\n
$$
\mathbf{v}_{0} = \mathbf{v}_{0}/|\mathbf{r}_{ij}|^{6}
$$
\n
$$
\mathbf{v}_{0} = \mathbf{v}_{0}/|\mathbf{r}_{0}|^{6}
$$
\n
$$
\mathbf{v}_{0} = \mathbf{v}_{0}/|\mathbf{
$$

 $n = (1+\sigma^2)/2$)/2 $V_{ii} = V_0 / |r_{ii}|^6$

 $V₀$ can be tuned so that Rydberg excitations in neighbouring sites are forbidden.

System of ⁸⁷Rb atoms controllably coupled to their Rydberg excited state.

atoms in their excited (Rydberg) states is denoted by V and is a tunable parameter.

One can vary the detuning parameter Δ which allows one to preferentially put the atom in a Rydberg or ground state

Similar to the transition found in tilted optical lattice

S. Sachdev et al, PRB 66, 075128 (2002), P Fendley et al PRB 69, 075106 (2004)

Mapping to a constrained model

$$
\frac{\mathcal{H}}{\hbar} = \sum_{i} \frac{\Omega_i}{2} \sigma_x^i - \sum_{i} \Delta_i n_i + \sum_{i < j} V_{ij} n_i n_j,
$$

trained model $\Delta_i n_i + \sum_{i < j} V_{ij} n_i n_j,$ Rydberg blockade on neighboring sites: V_{i,i+1} >> Δ , Ω >> V_{i,i+2}
 $P_\ell = (1-\sigma_\ell^z)/2$ Two states per site: Natural spin 1/2 representation Rydberg blockade on neighboring sites: $V_{i,i+1} \gg \Delta$, $\Omega \gg V_{i,i+2}$

A up-spin (Rydberg excitation) can be created on a site if and only if there are no up-spins (excitations) on the neighboring sites

$$
F_{\ell} = (1 - \sigma_{\ell})/2
$$
\n
$$
A \text{ up-spin (Rydberg excitation) can be created on a site}
$$
\nif and only if there are no up-spins (excitations) on the
\nneighboring sites\n
$$
\frac{\mathcal{H}}{\hbar} = \sum_{i} \frac{\Omega_{i}}{2} \sigma_{x}^{i} - \sum_{i} \Delta_{i} n_{i} + \sum_{i < j} V_{ij} n_{i} n_{j},
$$
\n
$$
= \sum_{\ell} (-w\tilde{\sigma}_{\ell}^{x} + \lambda/2\sigma_{\ell}^{x})
$$
\n
$$
= \sum_{\ell} (-w\tilde{\sigma}_{\ell}^{x} + \lambda/2\sigma_{\ell}^{x})
$$
\nDivie the system by making λ and w periodic function of time with frequencies Ω_{1} and Ω_{2} : $\Omega_{2} = 3 \Omega_{1}$

Drive the system by making λ and w periodic function of time with frequencies Ω_1 and Ω_2 : $\,\Omega_2$ =3 Ω_1

$$
\sigma_{\ell}^z = 2n_{\ell} - 1, \quad \sigma_{\ell}^{x(y)} = (i)(d_{\ell} + (-)d_{\ell}^{\dagger}).
$$

Dynamics around exact flat bands

We study the PXP Hamiltonian describing Rydberg atom array

Floquet bands and the spectral form factor

\n
$$
H_R = \lambda_1(t) \sum_j \sigma_j^z + \lambda_2(t) \sum_j \tilde{\sigma}_j^x
$$
\nUse exact di the eigenspe

Use exact diagonalization (ED) to find Use exact diagonalization (ED) to find
the eigenspectrum of U and hence H_F

Floquet bands and the sp

ve study the PXP Hamiltonian
 $H_R = \lambda_1(t) \sum_j \sigma_j^z + \lambda$

Secribing Rydberg atom array

Non-perturbative regime: w_j

Non-perturbative regime: w_j

Floquet bandwidth
 λ_F as a function of
 Ω_1 Floquet bandwidth
 A_F as a function of
 A_F as a function of Ω_1 and for $v=3$. 0.5 For single rate drive $\sqrt{\frac{6.06}{1}}$ Floquet bands and the specifies of the PXP Hamiltonian

Exerciting Rydberg atom array
 $H_R = \lambda_1(t) \sum_j \sigma_j^z + \lambda$

Non-perturbative regime: w_j

Non-perturbative regime: w_j

Floquet bandwidth
 A_F as a function of
 Ω_1 $\Lambda_{\mathsf{F}} = h\Omega_1$ 0.0 $\begin{bmatrix}\n\vdots & \vdots & \vdots & \vdots \\
\hline\n\vdots & \vdots & \vdots \\
\hline\n\$ paret bandwidth

and for $v=3$.

and for $v=3$.

and $v=4$.
 $\frac{10}{2}$ and $\frac{10}{2}$ and $\frac{1}{2}$ and $\frac{1}{2}$ and $\frac{1}{2}$ around the first Floquet Billouin zone
 $v=6$.
 $\frac{1}{2}$ and $\frac{1}{2}$ and $\frac{1}{2}$ around Strong deviation from prediction of Floquet ETH (realized for single rate drive protocol).

Strong deviation from prediction of Floquet ETH (realized for single rate drive protocol).

Strong deviation from prediction of F

Plot of distribution, P, of the act diagonalization (ED) to find
renspectrum of U and hence H_F
Plot of distribution, P, of the
Floquet eigenenergies within
the first Floquet Brillouin zone
for $\Omega_1 = w_1 = \lambda_1$.
For w_o=0, P is a delta function act diagonalization (ED) to find
yenspectrum of U and hence H_F
Plot of distribution, P, of the
Floquet eigenenergies within
the first Floquet Brillouin zone
for Ω_1 = w₁= λ_1 .
For w₀=0, P is a delta function
pe for $\Omega_1 = w_1 = \lambda_1$.

for \varOmega_1 = w_1 = λ_1 .
For w₀=0, P is a delta function peaked at zero while Floquet ETH predicts a flat distribution

Red dashed line $\rightarrow w_0 = 0$ where one has exact Floquet flat band Black dashed line $\,\Rightarrow$ single rate drive protocol with $w_{\text{\scriptsize{$1$}}}$ =0 and $w_{\text{\scriptsize{$0$}}}$ =1 $=1$

Non-perturbative regime: $\bm{w}_1 \!\!=\!\! \lambda_1 \!\!=\!\! \Omega_1$ and $\bm{w}_0 \!<\!\! \bm{w}_1$

Spectral form factor

SYK model for

L=16, two rate protocol

For the two-rate protocol, the dip time can be extended for small w_0

The ramp region signifying thermalization always occur after the dip 10^{-7}

This indicates a large prethermal timescale and slow thermalization at small $w₀$

Numerically, we find $1/t_{dip} \sim A_F$ and hence w_o : w_o provides a knob for 10^{-10} slowing down thermalization

Dependence of dip time on Floquet bandwidth
To obtain an estimate of the dip time we note the following: To obtain an estimate of the dip time we note the following:

1) At short times, the spectral form factor can be written as

2) The sum over eigenstates can be converted to an integral over energy gaps using a density of states ρ the of dip time on Floquet bandwidth

To obtain an estimate of the dip time we note the following

1) At short times, the spectral form factor can be written
 $K(nT_1) \approx 1 - (nT_1)^2 \sum_{p,q} (E_p^F - E_q^F)^2 / (2\hbar^2 \mathcal{D}^2)$.

2) T 1) At short times, the spectral form factor can be written
 $K(nT_1) \approx 1 - (nT_1)^2 \sum_{p,q} (E_p^F - E_q^F)^2 / (2\hbar^2 \mathcal{D}^2)$.

2) The sum over eigenstates can be converted to an integ

over energy gaps using a density of states \r 2) The sum over eigenstates can be converted to an integral
over energy gaps using a density of states ρ
 $\int_{-A_F/2}^{A_F/2} \rho d\epsilon = \mathcal{D} = \sum_{m_A}$
3) We choose $\rho = \rho_0 \Lambda_F^{-1} f(\epsilon/\Lambda_F)$ $\rho_0 \sim \mathcal{D}$
4) Using this DOS, one fi

$$
\int_{-\Lambda_F/2}^{\Lambda_F/2} \rho d\epsilon = \mathcal{D} = \sum_{m} \chi
$$

$$
t_{\text{dip}} = n_d T_1
$$
, where $n_d \sim \text{Int}[\hbar/(T_1 \Lambda_F)]$

The steps occur due to integer nature of n_d

Correlation and entanglement

Half chain entanglement Starting from the Rydberg vacuum (all spin-down state)

The entanglement and the correlation function stays at their initial value in the flat band limit: perfect freezing.

Both the entanglement and correlation shows a broad dip over a wide range of frequency for finite w_o

correlation stays nears its diagonal ensemble predicted value while the entanglement saturates to its Page value

The half-chain entanglement, for two rate-protocol at finite $w₀$, saturates to a much lower value signifying a lower spread of the initial state in the Hilbert space.

Rapid decay of fidelity and growth of entanglement for single rate drive protocol (black-dashed line) In contrast, F(t) remains close to zero for two-rate drive protocols both zero and finite w_o The growth of entanglement is also much smaller for the two-rate protocol The spread of the driven state in the Hilbert space and hence heating is drastically reduced. For w_o=0, one has an exact symmetry in micromotion $\ O(t)$ = $O(T_{\text{1}}\text{-}t)$ for all t. The dynamics involves coherent reversal of excitations and is reminiscent of spin echoes.

Analytical result at high drive amplitude

Floquet Pertubation theory (single rate protocol for Rydberg model)
mall parameter: accurate for large drive amplitude and intermediate drive frequencies.
r, the evolution operator receives contribution from the σ^{z} Uses w/λ as the small parameter: accurate for large drive amplitude and intermediate drive frequencies.

At the zeroth order, the evolution operator receives contribution from the σ^2 term.

Example 18 Follow Finding Theory (single rate protocol for Rydberg model)

Uses w/λ as the small parameter: accurate for large drive amplitude and intermediate drive frequencies.

At the zeroth order, the evolution op

Using standard perturbation theory, one obtains the first order contribution to H_F as follows

$$
U_0(t,0) = e^{i\lambda t \sum_j \sigma_j^2/2} \text{ for } t \leq T/2, \qquad \langle m|U_0(t,0)|n\rangle = \delta_{mn}e^{im\lambda t/2} \text{ for } t \leq T/2, \qquad \hat{C}^2
$$
\n
$$
= e^{i\lambda(T-t) \sum_j \sigma_j^2/2} \text{ for } T/2 \leq t \leq T, \qquad \text{Using standard perturbation theory, one obtains the first order contribution to } H_f \text{ as follows}
$$
\n
$$
U_1(T,0) = -i \int_0^T dt H_1(t) \qquad \qquad \langle m|U_1(T,0)|n\rangle = \delta_{m,n+s} \frac{2w}{\lambda s} \left(e^{i\lambda sT/2} - 1\right) \qquad \text{s = \pm 1.}
$$
\n
$$
U_1(T,0) = -i \int_0^T dt H_1(t) \qquad \qquad \langle m|U_1(T,0)|n\rangle = \delta_{m,n+s} \frac{2w}{\lambda s} \left(e^{i\lambda sT/2} - 1\right) \qquad \text{s = \pm 1.}
$$
\n
$$
U_1(T,0) = \sum_m \sum_j \sum_{s_j = \pm 1} c_{s_j}^{(1)} |m\rangle \langle m + s_j|, \qquad \text{Magnus in rotated frame}
$$
\n
$$
c_s^{(1)} = \frac{4iw}{\lambda} \sin(\lambda T/4) e^{i\lambda Ts/4}, \qquad \text{Magnus in rotated frame}
$$
\n
$$
\text{Floguet Hamiltonian}
$$
\n
$$
H_F^{(1)} = -w \frac{\sin(\gamma)}{\gamma} \sum_j [\cos(\gamma)\tilde{\sigma}_j^x + \sin(\gamma)\tilde{\sigma}_j^y], \qquad \text{Resummation of the standard Magnus expansion}
$$

Application to two-rate protocol

Application to two-rate pr

\n
$$
H_F^{(1)} = \frac{w_0 \sin(\lambda_0 T_1/2)}{\lambda_0 T_1/2} \sum_{j,s=\pm} \tilde{\sigma}_j^s e^{-i\lambda_0 T_1 s/2}
$$
\nThe drive amplitude $2\lambda_0 >> w_o w_1$ for validity

\nof the perturbation theory.

\nFloquet Hamiltonian vanishes at w_0 =0; consistent with exact flat band.

\nThe presence of second order term is different from analogous expansion in single-rate protocol where

The drive amplitude $2\lambda_0 >> w_ow_1$ for validity w_1 of the perturbation theory.

Floquet Hamiltonian vanishes at $w_0=0$; consistent with exact flat band.

The presence of second order term is different from analogous expansion in single-rate protocol where only odd order terms are present.

This is due presence of chirality operator $\boldsymbol{\mathsf{C}}_o$ for single rate protocol satisfying

$$
C_0 = \prod_j \sigma_j^z \qquad \qquad {\{H_F, C_0\}} = 0.
$$

 w_1 and w_0

$$
C = \frac{2w_0w_1C}{\lambda_0} \sum_j (\sigma_j^z + (\tilde{\sigma}_j^+ \tilde{\sigma}_{j+1}^- + \text{h.c.}))
$$

\n
$$
C = 6[2\sin(x/6) - 2\sin(x/3) + \sin(x/2)]/x - 1,
$$
\n
$$
0.15 - P
$$
\n
$$
P
$$
\n
$$
0.10 - P
$$
\n
$$
0.05 - P
$$
\n
$$
0.00 - P
$$
\n $$

FIG. 6. Plot of the eigenvalue distribution P within the first FBZ for $\hbar\Omega_1/w_0 = 15$, $w_1/w_0 = 0.1$, and $\lambda_0/w_0 = 20$. The red This property is absent for the two-rate protocol for finite
squares indicate those from exact numerics obtained using ED. For both plots $L = 24$, $K_0 = 0$ and we have used square-pulse protocol with $\nu = 3$.

Generic properties

At the initial time, C=0 and F=1.

As the operator spreads, F -> 1 and C->0

This spread of this correlation at initial times is linear and is bounded by the Lieb-Robinson velocity.

For ergodic systems, F shows a rapid convergence to its steady state value ($F \sim 1$).

For fragmented systems, the late time values of F depends on the initial condition.

