

Entanglement dynamics from universal low-lying modes

Shreya Vardhan

Stanford University

Workshop on Quantum Information, Quantum Field Theory and
Gravity, ICTS

August 27th, 2024

2407.16763 with Sanjay Moudgalya

Thermalization and entanglement entropy

- In this talk, we will discuss the emergence of **universal properties** of the **time-evolution** of entanglement entropy in **chaotic** quantum many-body systems.

Thermalization and entanglement entropy

- In this talk, we will discuss the emergence of **universal properties** of the **time-evolution** of entanglement entropy in **chaotic** quantum many-body systems.
- Consider the time-evolution of some initial state ρ_0 :

$$\rho(t) = e^{-iHt} \rho_0 e^{iHt}$$

Thermalization and entanglement entropy

- In this talk, we will discuss the emergence of **universal properties** of the **time-evolution** of entanglement entropy in **chaotic** quantum many-body systems.
- Consider the time-evolution of some initial state ρ_0 :

$$\rho(t) = e^{-iHt} \rho_0 e^{iHt}$$

If H is chaotic, then at late times, most details of ρ_0 are forgotten and $\rho(t)$ resembles a thermal density matrix $\rho^{(\text{eq})}$. This process is known as **thermalization**.

Thermalization and entanglement entropy

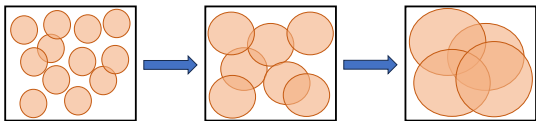
- In this talk, we will discuss the emergence of **universal properties** of the **time-evolution** of entanglement entropy in **chaotic** quantum many-body systems.
- Consider the time-evolution of some initial state ρ_0 :

$$\rho(t) = e^{-iHt} \rho_0 e^{iHt}$$

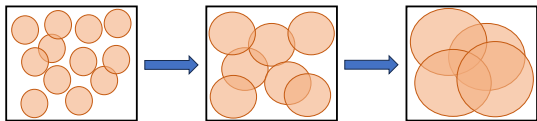
If H is chaotic, then at late times, most details of ρ_0 are forgotten and $\rho(t)$ resembles a thermal density matrix $\rho^{(\text{eq})}$. This process is known as **thermalization**.

- Holographic CFTs are an example of highly chaotic systems. Thermalization in the CFT corresponds to **black hole formation** in the bulk.

- With local interactions, increasingly large regions are thermalized as time evolves:

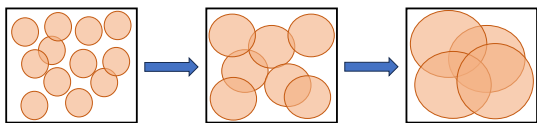


- With local interactions, increasingly large regions are thermalized as time evolves:



- For any fixed region A (smaller than half of the full system), the reduced density matrix $\rho_A(t)$ approaches thermal density matrix.

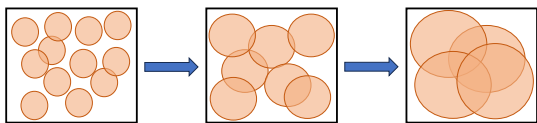
- With local interactions, increasingly large regions are thermalized as time evolves:



- For any fixed region A (smaller than half of the full system), the reduced density matrix $\rho_A(t)$ approaches thermal density matrix.
- One way to probe this process is by considering time-evolution of the Renyi entropies,

$$S_{n,A}(t) = -\frac{1}{n-1} \log \text{Tr}[\rho_A(t)^n], \quad n = 1, 2, \dots$$

- With local interactions, increasingly large regions are thermalized as time evolves:

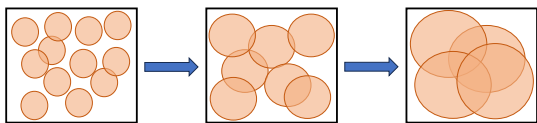


- For any fixed region A (smaller than half of the full system), the reduced density matrix $\rho_A(t)$ approaches thermal density matrix.
- One way to probe this process is by considering time-evolution of the Renyi entropies,

$$S_{n,A}(t) = -\frac{1}{n-1} \log \text{Tr}[\rho_A(t)^n], \quad n = 1, 2, \dots$$

$n \rightarrow 1$ limit is von Neumann entropy.

- With local interactions, increasingly large regions are thermalized as time evolves:



- For any fixed region A (smaller than half of the full system), the reduced density matrix $\rho_A(t)$ approaches thermal density matrix.
- One way to probe this process is by considering time-evolution of the Renyi entropies,

$$S_{n,A}(t) = -\frac{1}{n-1} \log \text{Tr}[\rho_A(t)^n], \quad n = 1, 2, \dots$$

$n \rightarrow 1$ limit is von Neumann entropy.

- Due to thermalization, these quantities will approach extensive thermal values at late times:

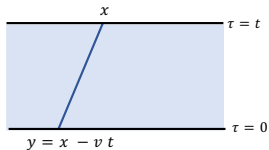
$$\lim_{t \rightarrow \infty} S_{n,A}(t) = s_{\text{eq}} V_A.$$

Universality in approach to equilibrium

- We expect there to be universality not only in the late-time saturation values of various quantities during thermalization, but also in the way in which they are approached.
- Evolution of entanglement entropy in generic chaotic time-evolutions is very difficult to study analytically.
- But the few analytically tractable examples we can study suggest a remarkable universality.
- In both **random circuits** and **holographic CFTs**, the evolution of entanglement entropy at late times can be expressed in terms of a **membrane formula**.
- Conjectured to hold universally in **Jonay, Huse, Nahum**.

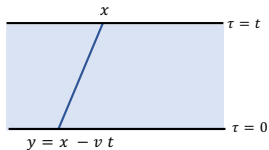
Membrane picture for entanglement growth

- In one spatial dimension, suppose we want to find S_n of the region to the left of some x at time t .
- Extend the system in time direction from $\tau = 0$ to $\tau = t$, and consider lines with different velocities v :



Membrane picture for entanglement growth

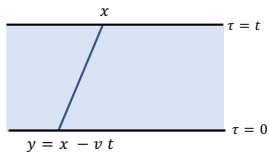
- In one spatial dimension, suppose we want to find S_n of the region to the left of some x at time t .
- Extend the system in time direction from $\tau = 0$ to $\tau = t$, and consider lines with different velocities v :



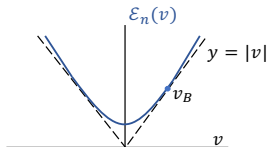
$$S_n(x, t) = \min_v [s_{\text{eq}} \mathcal{E}_n(v)t + S_n(y, t = 0)]$$

Membrane picture for entanglement growth

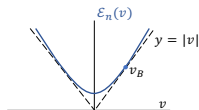
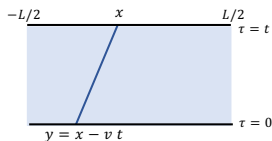
- In one spatial dimension, suppose we want to find the entanglement entropy of the left half-line at time t .
- Extend the system in time direction from $\tau = 0$ to $\tau = t$, and consider all possible curves:



$$S_n(x, t) = \min_v [s_{\text{eq}} \mathcal{E}_n(v)t + S_n(y, t = 0)]$$



Physical consequences



$$S_n(x, t) = \min_v [s_{\text{eq}} \mathcal{E}_n(v) t + S_n(y, t = 0)]$$

- Consider an initial state with volume law entanglement entropy, with coefficient s :

$$S_n(y, t = 0) = s \times (y + L/2), \quad 0 < s < s_{\text{eq}}.$$

- The membrane formula gives an s -dependent growth rate of $S_n(x, t)$:

$$S_n(x, t) = S_n(x, t = 0) + s_{\text{eq}} \Gamma_n(s) t$$

where Γ is related to \mathcal{E} by Legendre transform.

- Constraints on membrane tension are equivalent to the condition that

$$\Gamma_n(s_{\text{eq}}) = 0.$$

- While this formula turns out to hold in both random circuits and holographic CFTs, the methods used in the two cases are very different.

- While this formula turns out to hold in both random circuits and holographic CFTs, the methods used in the two cases are very different.
- In random circuits, Haar averages for n -th Renyi entropy. (Explicit calculation for the second Renyi entropy, and results in certain limits for third Renyi entropy.)

- While this formula turns out to hold in both random circuits and holographic CFTs, the methods used in the two cases are very different.
- In random circuits, Haar averages for n -th Renyi entropy. (Explicit calculation for the second Renyi entropy, and results in certain limits for third Renyi entropy.)
- In holographic CFTs:

- While this formula turns out to hold in both random circuits and holographic CFTs, the methods used in the two cases are very different.
- In random circuits, Haar averages for n -th Renyi entropy. (Explicit calculation for the second Renyi entropy, and results in certain limits for third Renyi entropy.)
- In holographic CFTs:
 - Thermalization in boundary is dual to gravitational collapse in bulk.

- While this formula turns out to hold in both random circuits and holographic CFTs, the methods used in the two cases are very different.
- In random circuits, Haar averages for n -th Renyi entropy. (Explicit calculation for the second Renyi entropy, and results in certain limits for third Renyi entropy.)
- In holographic CFTs:
 - Thermalization in boundary is dual to gravitational collapse in bulk.
 - Liu and Suh, Hartman and Maldacena applied HRT formula Hubeny, Rangamani, Takayanagi to find the evolution of von Neumann entropy in this setup.

- While this formula turns out to hold in both random circuits and holographic CFTs, the methods used in the two cases are very different.
- In random circuits, Haar averages for n -th Renyi entropy. (Explicit calculation for the second Renyi entropy, and results in certain limits for third Renyi entropy.)
- In holographic CFTs:
 - Thermalization in boundary is dual to gravitational collapse in bulk.
 - Liu and Suh, Hartman and Maldacena applied HRT formula Hubeny, Rangamani, Takayanagi to find the evolution of von Neumann entropy in this setup.
 - Mezei showed that for large system size and time, we can get rid of the radial direction in the bulk, and reduce the HRT formula to a minimization problem in the boundary.

- While this formula turns out to hold in both random circuits and holographic CFTs, the methods used in the two cases are very different.
- In random circuits, Haar averages for n -th Renyi entropy. (Explicit calculation for the second Renyi entropy, and results in certain limits for third Renyi entropy.)
- In holographic CFTs:
 - Thermalization in boundary is dual to gravitational collapse in bulk.
 - Liu and Suh, Hartman and Maldacena applied HRT formula Hubeny, Rangamani, Takayanagi to find the evolution of von Neumann entropy in this setup.
 - Mezei showed that for large system size and time, we can get rid of the radial direction in the bulk, and reduce the HRT formula to a minimization problem in the boundary.
 - The resulting membrane tension satisfies non-trivial constraints from Jonay, Huse, Nahum.

Questions

- What is the physical meaning of the entanglement membrane, and the source of this universality?

Questions

- What is the physical meaning of the entanglement membrane, and the source of this universality?
- Heuristically, if there is a tensor network representation of the state, we may think of the membrane as a “minimal cut.”

Questions

- What is the physical meaning of the entanglement membrane, and the source of this universality?
- Heuristically, if there is a tensor network representation of the state, we may think of the membrane as a “minimal cut.”
- But we would like to have a more precise understanding of the following questions:

Questions

- What is the physical meaning of the entanglement membrane, and the source of this universality?
- Heuristically, if there is a tensor network representation of the state, we may think of the membrane as a “minimal cut.”
- But we would like to have a more precise understanding of the following questions:
 1. What is the source of the velocity-dependent function $\mathcal{E}(v)$?

Questions

- What is the physical meaning of the entanglement membrane, and the source of this universality?
- Heuristically, if there is a tensor network representation of the state, we may think of the membrane as a “minimal cut.”
- But we would like to have a more precise understanding of the following questions:
 1. What is the source of the velocity-dependent function $\mathcal{E}(v)$?
 2. Is there an underlying structure in terms of low-lying modes, which we could look for in a continuum theory such as a holographic CFT?

Questions

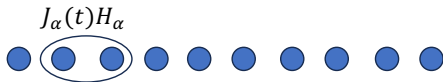
- What is the physical meaning of the entanglement membrane, and the source of this universality?
- Heuristically, if there is a tensor network representation of the state, we may think of the membrane as a “minimal cut.”
- But we would like to have a more precise understanding of the following questions:
 1. What is the source of the velocity-dependent function $\mathcal{E}(v)$?
 2. Is there an underlying structure in terms of low-lying modes, which we could look for in a continuum theory such as a holographic CFT?
 3. How does the structure of $\mathcal{E}(v)$ depend on the Renyi index n ?

- In this talk, we will propose a possible common underlying structure, involving certain universal low-lying modes.

- In this talk, we will propose a possible common underlying structure, involving certain universal low-lying modes.
- We will make use of a family of time-dependent Hamiltonians:

$$H(t) = \sum_{\alpha} J_{\alpha}(t) H_{\alpha},$$

where the H_{α} are local operators, and $J_{\alpha}(t)$ are random numbers, **uncorrelated for different times** and different α .



- In this talk, we will propose a possible common underlying structure, involving certain universal low-lying modes.
- We will make use of a family of time-dependent Hamiltonians:

$$H(t) = \sum_{\alpha} J_{\alpha}(t)H_{\alpha},$$

where the H_{α} are local operators, and $J_{\alpha}(t)$ are random numbers, **uncorrelated for different times** and different α .



- Previously, such models have allowed a derivation of **diffusion** in two-point functions [Moudgalya and Motrunich](#); [Ogunnaike, Feldmeier, Lee](#).

From Lorentzian to Euclidean time-evolution

- Observables of interest, such as the n -th Renyi entropy, can be written as a transition amplitudes under $(U \otimes U^*)^{\otimes n}$ in any system.

From Lorentzian to Euclidean time-evolution

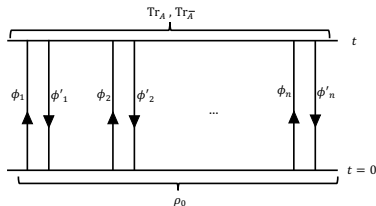
- Observables of interest, such as the n -th Renyi entropy, can be written as a transition amplitudes under $(U \otimes U^*)^{\otimes n}$ in any system.

$$e^{-(n-1)S_n^{(A)}} = \text{Tr}_A \rho_A^n = \text{Tr}_A \left(\text{Tr}_{\bar{A}} U \rho_0 U^\dagger \right)^n$$

From Lorentzian to Euclidean time-evolution

- Observables of interest, such as the n -th Renyi entropy, can be written as a transition amplitudes under $(U \otimes U^*)^{\otimes n}$ in any system.

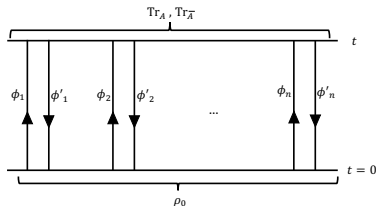
$$e^{-(n-1)S_n^{(A)}} = \text{Tr}_A \rho_A^n = \text{Tr}_A \left(\text{Tr}_{\bar{A}} U \rho_0 U^\dagger \right)^n$$



From Lorentzian to Euclidean time-evolution

- Observables of interest, such as the n -th Renyi entropy, can be written as a transition amplitudes under $(U \otimes U^*)^{\otimes n}$ in any system.

$$e^{-(n-1)S_n^{(A)}} = \text{Tr}_A \rho_A^n = \text{Tr}_A \left(\text{Tr}_{\bar{A}} U \rho_0 U^\dagger \right)^n$$



- The key simplification in Brownian models is that the Lorentzian evolution on $2n$ copies can be replaced with a **Euclidean evolution**:

$$\overline{(U(t) \otimes U(t)^*)^{\otimes n}} = e^{-P_{2n}t}$$

- Let us derive the Euclidean evolution explicitly in the two-copy case ($n = 1$).

- Let us derive the Euclidean evolution explicitly in the two-copy case ($n = 1$).
- Divide time-evolution into small steps of size ϵ .

$$H(t) = \sum_{\alpha} J_{\alpha}(t) H_{\alpha}$$

$$\overline{J_{\alpha}(t)} = 0, \quad \overline{J_{\alpha}(t) J_{\alpha'}(t')} = \frac{1}{2} \frac{\delta_{\alpha\alpha'} \delta_{tt'}}{\epsilon}$$

- Let us derive the Euclidean evolution explicitly in the two-copy case ($n = 1$).
- Divide time-evolution into small steps of size ϵ .

$$H(t) = \sum_{\alpha} J_{\alpha}(t) H_{\alpha}$$

$$\overline{J_{\alpha}(t)} = 0, \quad \overline{J_{\alpha}(t) J_{\alpha'}(t')} = \frac{1}{2} \frac{\delta_{\alpha\alpha'} \delta_{tt'}}{\epsilon}$$

- Then we have:

$$e^{iH_a(t)\epsilon} \otimes e^{-iH_b(t)^T \epsilon}$$

- Let us derive the Euclidean evolution explicitly in the two-copy case ($n = 1$).
- Divide time-evolution into small steps of size ϵ .

$$H(t) = \sum_{\alpha} J_{\alpha}(t) H_{\alpha}$$

$$\overline{J_{\alpha}(t)} = 0, \quad \overline{J_{\alpha}(t) J_{\alpha'}(t')} = \frac{1}{2} \frac{\delta_{\alpha\alpha'} \delta_{tt'}}{\epsilon}$$

- Then we have:

$$e^{iH_a(t)\epsilon} \otimes e^{-iH_b(t)^T \epsilon}$$

$$\approx (1 + iH_a(t)\epsilon - \frac{1}{2}H_a(t)^2\epsilon^2 + \dots) \otimes (1 - iH_b(t)^T \epsilon - \frac{1}{2}H_b(t)^T{}^2\epsilon^2 + \dots)$$

- Let us derive the Euclidean evolution explicitly in the two-copy case ($n = 1$).
- Divide time-evolution into small steps of size ϵ .

$$H(t) = \sum_{\alpha} J_{\alpha}(t) H_{\alpha}$$

$$\overline{J_{\alpha}(t)} = 0, \quad \overline{J_{\alpha}(t) J_{\alpha'}(t')} = \frac{1}{2} \frac{\delta_{\alpha\alpha'} \delta_{tt'}}{\epsilon}$$

- Then we have:

$$e^{iH_a(t)\epsilon} \otimes e^{-iH_b(t)^T \epsilon}$$

$$\begin{aligned} &\approx (1 + iH_a(t)\epsilon - \frac{1}{2}H_a(t)^2\epsilon^2 + \dots) \otimes (1 - iH_b(t)^T\epsilon - \frac{1}{2}H_b(t)^T{}^2\epsilon^2 + \dots) \\ &= 1 - \epsilon P_2 + O(\epsilon^2) \quad \approx \quad e^{-\epsilon P} \end{aligned}$$

where

$$P_2 = \sum_{\alpha} (H_{a,\alpha} - H_{b,\alpha}^T)^2$$

Summary

- The n -th Renyi entropy can be expressed as a transition amplitude under Euclidean evolution with a non-negative Hamiltonian P_{2n} .

Summary

- The n -th Renyi entropy can be expressed as a transition amplitude under Euclidean evolution with a non-negative Hamiltonian P_{2n} .
- The equilibrium saturation value of the n -th Renyi entropy is determined by the **zero energy states** of P_{2n} .

The result is consistent with the equilibrium approximation of [Liu and SV](#).

Summary

- The n -th Renyi entropy can be expressed as a transition amplitude under Euclidean evolution with a non-negative Hamiltonian P_{2n} .
- The equilibrium saturation value of the n -th Renyi entropy is determined by the **zero energy states** of P_{2n} .

The result is consistent with the equilibrium approximation of [Liu and SV](#).

- Approach to equilibrium is determined by **low energy eigenstates**, which have a universal structure.

- For the second Renyi entropy in models without symmetries:

- For the second Renyi entropy in models without symmetries:

P_4 has two degenerate ground states:

$$|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle, \quad |\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle$$

- The low-energy excitations include a “one-particle” band approximately given by:

$$|\psi_k\rangle = \sum_x e^{i k x} |\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle_x |\phi_{x+1,\dots,x+d}\rangle |\uparrow\rangle_{x+d+1} |\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle$$

for some $O(1)$ d .

- For the second Renyi entropy in models without symmetries:

P_4 has two degenerate ground states:

$$|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle, \quad |\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle$$

- The low-energy excitations include a “one-particle” band approximately given by:

$$|\psi_k\rangle = \sum_x e^{i k x} |\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle|\downarrow\rangle_x |\phi_{x+1,\dots,x+d}\rangle |\uparrow\rangle_{x+d+1} |\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle|\uparrow\rangle$$

for some $O(1)$ d .

- This structure leads to the membrane picture.

- The one-particle excitations have a gapped dispersion relation $E(k)$. The entanglement growth rate is given by

$$\Gamma_2(s) = E_2(k = is)/s_{\text{eq}}$$

$\mathcal{E}(v)$ can be obtained from this quantity by Legendre transform.

- The one-particle excitations have a gapped dispersion relation $E(k)$. The entanglement growth rate is given by

$$\Gamma_2(s) = E_2(k = is)/s_{\text{eq}}$$

$\mathcal{E}(v)$ can be obtained from this quantity by Legendre transform.

- Dispersion relation at $O(1)$ values of k is physically important for satisfying constraints.

- The one-particle excitations have a gapped dispersion relation $E(k)$. The entanglement growth rate is given by

$$\Gamma_2(s) = E_2(k = is)/s_{\text{eq}}$$

$\mathcal{E}(v)$ can be obtained from this quantity by Legendre transform.

- Dispersion relation at $O(1)$ values of k is physically important for satisfying constraints.
- For the third Renyi entropy, we have an analogous set of low-energy eigenstates. In addition to these, competition from another set of eigenstates leads to phase transitions in $\mathcal{E}_3(v)$ as a function of v .

Plan

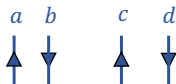
- Introduce expression for the second Renyi entropy as a transition amplitude, and the definition of $|\uparrow\rangle$ and $|\downarrow\rangle$.
- Derive the low-energy excitations in a simplifying limit.
- Discuss how the structure remains robust more generally.
- Discuss qualitatively new features of the third Renyi entropy.

Second Renyi entropy as transition amplitude

Second Renyi entropy as a transition amplitude

- Second Renyi entropy involves two forward and two backward copies of U :

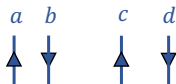
$$e^{-S_{2,A}(t)} = \text{Tr}_A \left(\text{Tr}_{\bar{A}} U \rho_0 U^\dagger \right)^2$$



Second Renyi entropy as a transition amplitude

- Second Renyi entropy involves two forward and two backward copies of U :

$$e^{-S_{2,A}(t)} = \text{Tr}_A \left(\text{Tr}_{\bar{A}} U \rho_0 U^\dagger \right)^2$$



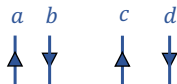
- Let us introduce the following “spins” on **four copies** of a **single site** in particular, Zhou and Nahum

$$|\uparrow\rangle = |\text{MAX}\rangle_{ab} |\text{MAX}\rangle_{cd}, \quad |\downarrow\rangle = |\text{MAX}\rangle_{ad} |\text{MAX}\rangle_{bc}$$

Second Renyi entropy as a transition amplitude

- Second Renyi entropy involves two forward and two backward copies of U :

$$e^{-S_{2,A}(t)} = \text{Tr}_A \left(\text{Tr}_{\bar{A}} U \rho_0 U^\dagger \right)^2$$



- Let us introduce the following “spins” on **four copies** of a **single site** in particular, Zhou and Nahum

$$|\uparrow\rangle = |\text{MAX}\rangle_{ab} |\text{MAX}\rangle_{cd}, \quad |\downarrow\rangle = |\text{MAX}\rangle_{ad} |\text{MAX}\rangle_{bc}$$

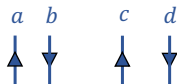
- Evolution of second Renyi entropy is given by

$$e^{-S_{2,A}(t)} = \langle D_{\Sigma_A} | (U \otimes U^*)^2 | \rho_0 \rangle | \rho_0 \rangle$$

Second Renyi entropy as a transition amplitude

- Second Renyi entropy involves two forward and two backward copies of U :

$$e^{-S_{2,A}(t)} = \text{Tr}_A \left(\text{Tr}_{\bar{A}} U \rho_0 U^\dagger \right)^2$$



- Let us introduce the following “spins” on **four copies** of a **single site** in particular, Zhou and Nahum

$$|\uparrow\rangle = |\text{MAX}\rangle_{ab} |\text{MAX}\rangle_{cd}, \quad |\downarrow\rangle = |\text{MAX}\rangle_{ad} |\text{MAX}\rangle_{bc}$$

- Evolution of second Renyi entropy is given by

$$e^{-S_{2,A}(t)} = \langle D_{\Sigma_A} | (U \otimes U^*)^2 | \rho_0 \rangle | \rho_0 \rangle$$

where

$$\langle D_{\Sigma_A} | = \otimes_{i \in A} \langle \downarrow | \otimes_{i \in \bar{A}} \langle \uparrow | \quad \overbrace{\begin{array}{c} \text{A} \\ \downarrow \downarrow \downarrow \downarrow \downarrow \downarrow \downarrow \downarrow \end{array}} \quad \uparrow \uparrow \uparrow \uparrow \uparrow \uparrow$$

Equilibrium value in models without conserved quantities

- P_4 generally has exactly two zero energy eigenstates:

$$|\uparrow \dots \uparrow\rangle, \quad |\downarrow \dots \downarrow\rangle$$

- This gives the Page value for the entropy of pure state at late times:

$$\lim_{t \rightarrow \infty} S_{2,A}(t) = \min(\log d_A, \log d_{\bar{A}})$$

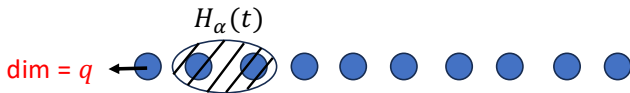
- We would now like to understand the approach to this value using the low-energy modes of P_4 .

Low energy excitations: GUE model

GUE model

- Take each $H_\alpha(t)$ to be an i.i.d. random Hermitian matrix on adjacent sites drawn from the GUE ensemble:

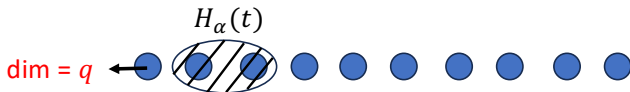
$$H_\alpha(t) = H_{i,i+1}^{(\text{GUE})}(t)$$



GUE model

- Take each $H_\alpha(t)$ to be an i.i.d. random Hermitian matrix on adjacent sites drawn from the GUE ensemble:

$$H_\alpha(t) = H_{i,i+1}^{(\text{GUE})}(t)$$



- Using the average over these random matrices, P_4 can be expressed entirely in terms of $|\uparrow\rangle, |\downarrow\rangle$.

Analytically solvable large q limit

- In the large q limit, P_4 is exactly solvable, and has a very simple action on a single domain wall

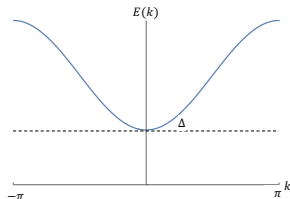
$$\langle D_x | \equiv \langle \downarrow \downarrow \dots \downarrow_x \uparrow_{x+1} \uparrow \dots \uparrow |$$

$$\langle D_x | P_4 = \langle D_x | - \frac{1}{q} (\langle D_{x-1} | + \langle D_{x+1} |)$$

- This leads to the following band of lowest excited states:

$$\langle \psi_k | = \sum_x e^{ikx} \langle D_x |$$

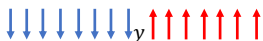
$$E(k) = 1 - \frac{2}{q} \cos k$$



Second Renyi entropy for half-line region

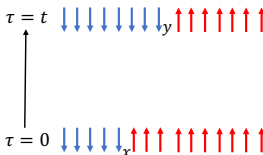
- Let us return to the second Renyi entropy of a half-line region:

$$e^{-S_2(y,t)} = \langle D_y | e^{-P_4 t} | \rho_0, e \rangle$$



- Since $\langle D_y |$ only evolves to a superposition of $\langle D_x |$ at other locations,

$$\begin{aligned} e^{-S_2(y,t)} &= \sum_x \langle D_y | e^{-P_4 t} | \bar{D}_x \rangle \langle D_x | \rho_0, e \rangle \\ &= \sum_x \langle D_y | e^{-P_4 t} | \bar{D}_x \rangle e^{-S_2(x,t=0)} \end{aligned}$$



Membrane picture from one domain wall band

- Using one-particle eigenstates in domain wall propagator:

$$\langle D_y | e^{-P_4 t} | \bar{D}_x \rangle = \sum_k e^{ik(x-y)} e^{-E(k)t}$$

Membrane picture from one domain wall band

- Using one-particle eigenstates in domain wall propagator:

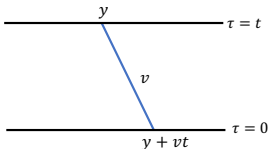
$$\langle D_y | e^{-P_4 t} | \bar{D}_x \rangle = \sum_k e^{ik(x-y)} e^{-E(k)t}$$

- At late times: using saddle-point approximation for the propagator,

$$S_2(y, t) = \min_v [s_{\text{eq}} \mathcal{E}(v) t + S_2(y + vt, t = 0)]$$

where

$$\mathcal{E}(v) = \frac{E(k_v) - ik_v v}{s_{\text{eq}}}, \quad k_v \text{ is solution to } E'(k_v) = iv.$$



Membrane picture from one domain wall band

- Using one-particle eigenstates in domain wall propagator:

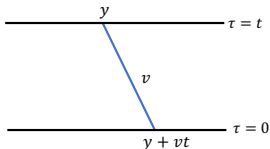
$$\langle D_y | e^{-P_4 t} | \bar{D}_x \rangle = \sum_k e^{ik(x-y)} e^{-E(k)t}$$

- At late times: using saddle-point approximation for the propagator,

$$S_2(y, t) = \min_v [s_{\text{eq}} \mathcal{E}(v) t + S_2(y + vt, t = 0)]$$

where

$$\mathcal{E}(v) = \frac{E(k_v) - ik_v v}{s_{\text{eq}}}, \quad k_v \text{ is solution to } E'(k_v) = iv.$$



- We can also check that for an initial state with volume law entropy with coefficient s ,

$$\Gamma(s) = E(k = is) / s_{\text{eq}}.$$

Finite q in Brownian GUE model

- Away from large q limit, interactions can cause domain walls to split, so the eigenstates and eigenvalues are modified.

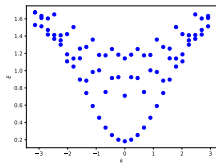
$$\langle D_x | P_4 = \langle D_x | - \frac{1}{q} (\langle D_{x-1} | + \langle D_{x+1} |) + \frac{1}{q^2} \langle D_{x-1,x,x+1} |$$

- Away from large q limit, interactions can cause domain walls to split, so the eigenstates and eigenvalues are modified.

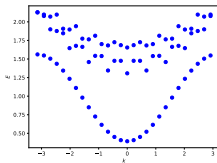
$$\langle D_x | P_4 = \langle D_x | - \frac{1}{q} (\langle D_{x-1} | + \langle D_{x+1} |) + \frac{1}{q^2} \langle D_{x-1,x,x+1} |$$

- From numerical diagonalization of P_4 :

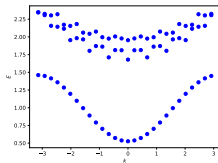
$q = 2$



$q = 3$



$q = 4$



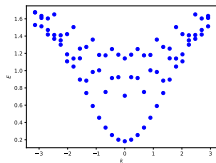
Gapped spectrum in all cases, which implies $v_E \neq 0$.

- Away from large q limit, interactions can cause domain walls to split, so the eigenstates and eigenvalues are modified.

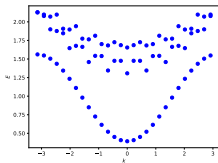
$$\langle D_x | P_4 = \langle D_x | - \frac{1}{q} (\langle D_{x-1} | + \langle D_{x+1} |) + \frac{1}{q^2} \langle D_{x-1,x,x+1} |$$

- From numerical diagonalization of P_4 :

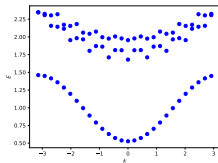
$q = 2$



$q = 3$



$q = 4$



Gapped spectrum in all cases, which implies $v_E \neq 0$.

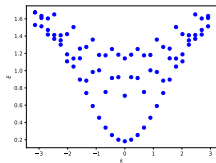
- Is the structure of the eigenstates robust?

- Away from large q limit, interactions can cause domain walls to split, so the eigenstates and eigenvalues are modified.

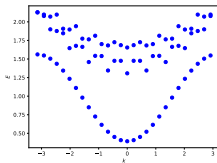
$$\langle D_x | P_4 = \langle D_x | - \frac{1}{q} (\langle D_{x-1} | + \langle D_{x+1} |) + \frac{1}{q^2} \langle D_{x-1,x,x+1} |$$

- From numerical diagonalization of P_4 :

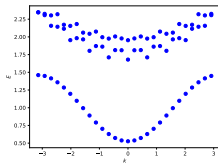
$q = 2$



$q = 3$



$q = 4$



Gapped spectrum in all cases, which implies $v_E \neq 0$.

- Is the structure of the eigenstates robust?
- Is there still a well-defined one-particle band within the continuum for $q = 2$?

Structure of eigenstates at finite q

- Let us consider a variational ansatz for the eigenstates:

$$|\psi_k\rangle = \sum_x e^{-ikx} |\downarrow \dots \downarrow_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |\uparrow_{x+\Delta+1} \dots \uparrow\rangle$$

Structure of eigenstates at finite q

- Let us consider a variational ansatz for the eigenstates:

$$|\psi_k\rangle = \sum_x e^{-ikx} |\downarrow \dots \downarrow_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |\uparrow_{x+\Delta+1} \dots \uparrow\rangle$$

- We can increase the value of Δ , and at each Δ , minimize

$$E_{\text{var}}(k) = \langle \psi_k | P_4 | \psi_k \rangle$$

over all choices of $|\phi\rangle$.

Structure of eigenstates at finite q

- Let us consider a variational ansatz for the eigenstates:

$$|\psi_k\rangle = \sum_x e^{-ikx} |\downarrow \dots \downarrow_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |\uparrow_{x+\Delta+1} \dots \uparrow\rangle$$

- We can increase the value of Δ , and at each Δ , minimize

$$E_{\text{var}}(k) = \langle \psi_k | P_4 | \psi_k \rangle$$

over all choices of $|\phi\rangle$.

- This minimization maps to the problem of diagonalizing an effective Hamiltonian on a 2^Δ -dimensional Hilbert space for each k .

Structure of eigenstates at finite q

- Let us consider a variational ansatz for the eigenstates:

$$|\psi_k\rangle = \sum_x e^{-ikx} |\downarrow \dots \downarrow_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |\uparrow_{x+\Delta+1} \dots \uparrow\rangle$$

- We can increase the value of Δ , and at each Δ , minimize

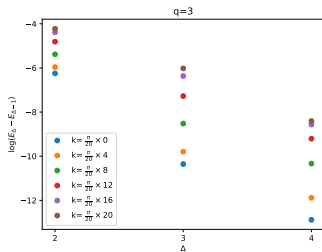
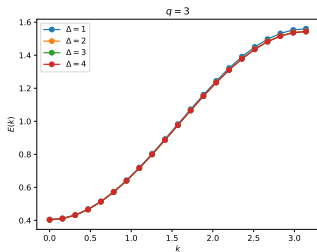
$$E_{\text{var}}(k) = \langle \psi_k | P_4 | \psi_k \rangle$$

over all choices of $|\phi\rangle$.

- This minimization maps to the problem of diagonalizing an effective Hamiltonian on a 2^Δ -dimensional Hilbert space for each k .
- Rapid convergence of $E_{\text{var}}(k)$ with Δ would tell us that the eigenstates are well-approximated by $|\psi_k\rangle$ for $O(1)$ Δ . Haegeman, Spyridon,

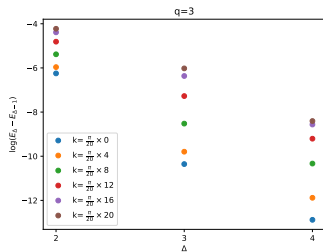
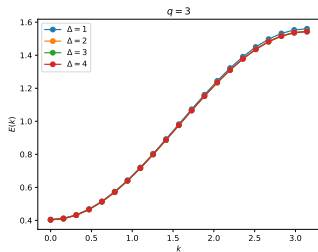
Variational results for $q = 3$

From minimizing $\langle \psi_k | A | \psi_k \rangle$ over all choices of $|\phi\rangle$ for various Δ :

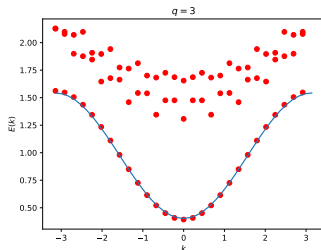


Variational results for $q = 3$

From minimizing $\langle \psi_k | A | \psi_k \rangle$ over all choices of $|\phi\rangle$ for various Δ :

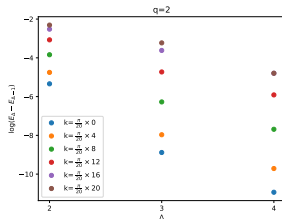
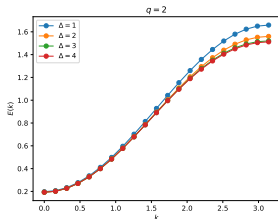


Good agreement with exact diagonalization results:



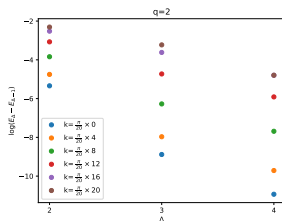
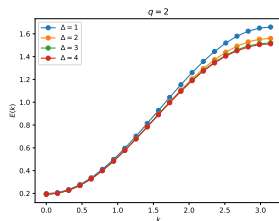
Variational results for $q = 2$

Still very good convergence with Δ :

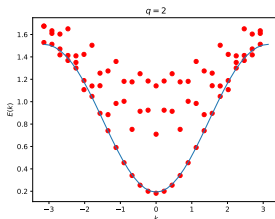


Variational results for $q = 2$

Still very good convergence with Δ :

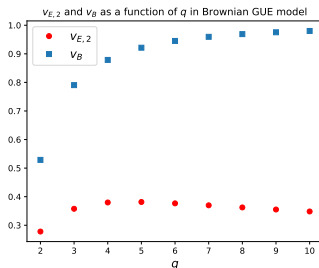
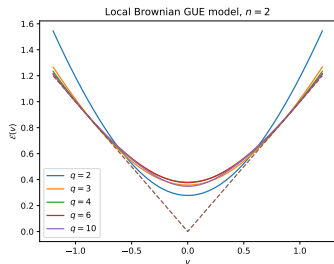


- Confirms that there is still a well-defined domain wall band within the continuum. ($\langle D_x |$ will only have significant overlap with this band.)



Membrane picture at finite q

- In the limit of large system sizes and late times, the local dressing with $O(1)$ Δ can be neglected.
- We still get the membrane formula at finite q , with $\mathcal{E}(v)$ given by Legendre transform of **exact dispersion relation**.
- $\mathcal{E}(v)$ can be found numerically from $E(k)$, and in particular we can check that the constraints on the membrane tension are satisfied:



Brownian mixed-field Ising model

- Let us now consider a spin-1/2 system, with the following time-dependent Hamiltonian.

$$H(t) = \sum_i J_Z(t) Z_i + J_X(t) X_i + J_{ZZ}(t) Z_i Z_{i+1}$$

where the variances are proportional to g_Z , g_X , g_{ZZ} .

- Let us now consider a spin-1/2 system, with the following time-dependent Hamiltonian.

$$H(t) = \sum_i J_Z(t) Z_i + J_X(t) X_i + J_{ZZ}(t) Z_i Z_{i+1}$$

where the variances are proportional to g_Z , g_X , g_{ZZ} .

- Let us fix $g_X = g_{ZZ} = 1$, and consider various values of g_Z .

- Let us now consider a spin-1/2 system, with the following time-dependent Hamiltonian.

$$H(t) = \sum_i J_Z(t) Z_i + J_X(t) X_i + J_{ZZ}(t) Z_i Z_{i+1}$$

where the variances are proportional to g_Z , g_X , g_{ZZ} .

- Let us fix $g_X = g_{ZZ} = 1$, and consider various values of g_Z .
- In this case,

- Let us now consider a spin-1/2 system, with the following time-dependent Hamiltonian.

$$H(t) = \sum_i J_Z(t) Z_i + J_X(t) X_i + J_{ZZ}(t) Z_i Z_{i+1}$$

where the variances are proportional to g_Z , g_X , g_{ZZ} .

- Let us fix $g_X = g_{ZZ} = 1$, and consider various values of g_Z .
- In this case,
 - Ground state subspace of P_4 is still spanned by $|\uparrow \dots \uparrow\rangle$, $|\downarrow \dots \downarrow\rangle$ (except at $g_Z = 0$).

- Let us now consider a spin-1/2 system, with the following time-dependent Hamiltonian.

$$H(t) = \sum_i J_Z(t) Z_i + J_X(t) X_i + J_{ZZ}(t) Z_i Z_{i+1}$$

where the variances are proportional to g_Z , g_X , g_{ZZ} .

- Let us fix $g_X = g_{ZZ} = 1$, and consider various values of g_Z .
- In this case,
 - Ground state subspace of P_4 is still spanned by $|\uparrow \dots \uparrow\rangle$, $|\downarrow \dots \downarrow\rangle$ (except at $g_Z = 0$).
 - Subspace spanned by arbitrary strings of \uparrow, \downarrow is no longer closed.

- Let us now consider a spin-1/2 system, with the following time-dependent Hamiltonian.

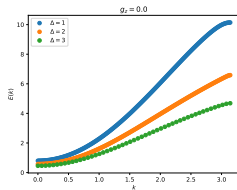
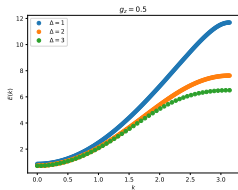
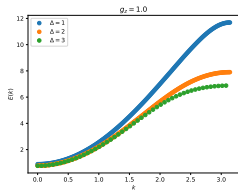
$$H(t) = \sum_i J_Z(t) Z_i + J_X(t) X_i + J_{ZZ}(t) Z_i Z_{i+1}$$

where the variances are proportional to g_Z , g_X , g_{ZZ} .

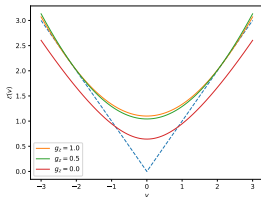
- Let us fix $g_X = g_{ZZ} = 1$, and consider various values of g_Z .
- In this case,
 - Ground state subspace of P_4 is still spanned by $|\uparrow \dots \uparrow\rangle$, $|\downarrow \dots \downarrow\rangle$ (except at $g_Z = 0$).
 - Subspace spanned by arbitrary strings of \uparrow, \downarrow is no longer closed.
- Again, let us consider a variational ansatz

$$|\phi_k\rangle = \sum_x e^{-ikx} |\downarrow \dots \downarrow_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |\uparrow_{x+\Delta+1} \dots \uparrow\rangle$$

where now, $|\phi_{x+1, \dots, x+n}\rangle$ is an arbitrary state in a 16^Δ -dimensional Hilbert space.

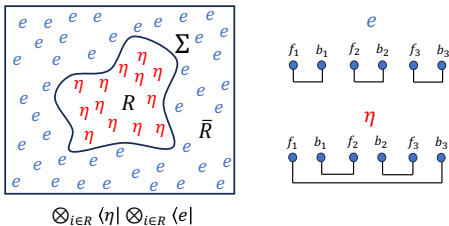


- Generic values of $g_z \neq 0$ should correspond to chaotic systems.
- For $g_z = 0$, $H(t)$ can be rewritten in terms of free fermions. The gap of P_4 vanishes and the ground state subspace is larger. [Swann, Bernard, Nahum](#)
- We see better convergence of $E(k)$ with Δ for $g_z \neq 0$.
- Membrane tensions from $\Delta = 3$ dispersion relations:



Higher Renyi entropies in Brownian GUE model

The n -th Renyi entropy can again be written as a transition amplitude, now with the final state:



where e is associated with identity permutation, and η with the cyclic permutation $(n \ n - 1 \dots 1)$.

- For the superhamiltonian in the Brownian GUE model, we now have an $n!$ -dimensional Hilbert space at each site.

- For the superhamiltonian in the Brownian GUE model, we now have an $n!$ -dimensional Hilbert space at each site.
- The state $\langle \eta \dots \eta_x e_{x+1} \dots e |$ evolves to a state consisting of other permutations $\sigma \in \mathcal{S}_n$.

- For the superhamiltonian in the Brownian GUE model, we now have an $n!$ -dimensional Hilbert space at each site.
- The state $\langle \eta \dots \eta_x e_{x+1} \dots e |$ evolves to a state consisting of other permutations $\sigma \in \mathcal{S}_n$.
- Let us make the variational ansatz that there are eigenstates of the form

$$|\psi_k\rangle = \sum_x e^{-ikx} |\eta \dots \eta_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |e_{x+\Delta+1} \dots e\rangle$$

where now $|\phi\rangle$ is an arbitrary state in an $(n!)^\Delta$ -dimensional Hilbert space.

- For the superhamiltonian in the Brownian GUE model, we now have an $n!$ -dimensional Hilbert space at each site.
- The state $\langle \eta \dots \eta_x e_{x+1} \dots e |$ evolves to a state consisting of other permutations $\sigma \in \mathcal{S}_n$.

- Let us make the variational ansatz that there are eigenstates of the form

$$|\psi_k\rangle = \sum_x e^{-ikx} |\eta \dots \eta_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |e_{x+\Delta+1} \dots e\rangle$$

where now $|\phi\rangle$ is an arbitrary state in an $(n!)^\Delta$ -dimensional Hilbert space.

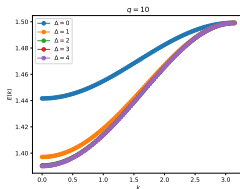
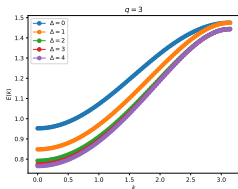
- For $n = 3$, we can use the variational method up to $\Delta = 4$.

- For the superhamiltonian in the Brownian GUE model, we now have an $n!$ -dimensional Hilbert space at each site.
- The state $\langle \eta \dots \eta_x e_{x+1} \dots e |$ evolves to a state consisting of other permutations $\sigma \in \mathcal{S}_n$.
- Let us make the variational ansatz that there are eigenstates of the form

$$|\psi_k\rangle = \sum_x e^{-ikx} |\eta \dots \eta_x\rangle |\phi_{x+1, \dots, x+\Delta}\rangle |e_{x+\Delta+1} \dots e\rangle$$

where now $|\phi\rangle$ is an arbitrary state in an $(n!)^\Delta$ -dimensional Hilbert space.

- For $n = 3$, we can use the variational method up to $\Delta = 4$.
- We find good convergence of $E_3(k)$ with Δ , indicating that we do have low-energy eigenstates of this form.



- By plugging this form of the eigenstates into the time-evolution of S_n for initial state with entropy density s , we now get

$$e^{-(n-1)S_n(x,t)} = \int_{-\infty}^{\infty} dv \int_{-\pi}^{\pi} \frac{dk}{2\pi} e^{-[E_n(k)+ikv-(n-1)sv]t}$$

- By plugging this form of the eigenstates into the time-evolution of S_n for initial state with entropy density s , we now get

$$e^{-(n-1)S_n(x,t)} = \int_{-\infty}^{\infty} dv \int_{-\pi}^{\pi} \frac{dk}{2\pi} e^{-[E_n(k)+ikv-(n-1)sv]t}$$

- Using the saddle-point equations, we get the entanglement growth rate

$$\Gamma_n(s) = \frac{E_n(-(n-1)is)}{(n-1)s_{\text{eq}}}$$

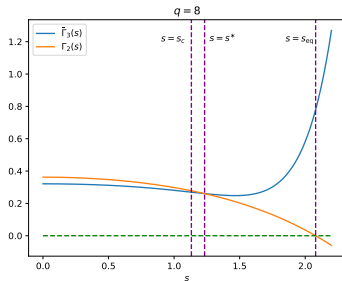
- By plugging this form of the eigenstates into the time-evolution of S_n for initial state with entropy density s , we now get

$$e^{-(n-1)S_n(x,t)} = \int_{-\infty}^{\infty} dv \int_{-\pi}^{\pi} \frac{dk}{2\pi} e^{-[E_n(k)+ikv-(n-1)sv]t}$$

- Using the saddle-point equations, we get the entanglement growth rate

$$\Gamma_n(s) = \frac{E_n(-(n-1)is)}{(n-1)s_{\text{eq}}}$$

- Using $E_3(k)$ from the variational calculation, we find:



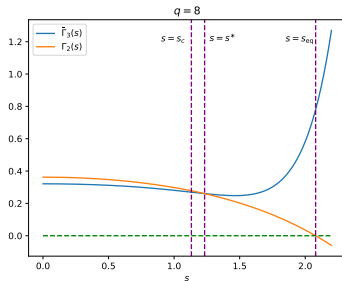
- By plugging this form of the eigenstates into the time-evolution of S_n for initial state with entropy density s , we now get

$$e^{-(n-1)S_n(x,t)} = \int_{-\infty}^{\infty} dv \int_{-\pi}^{\pi} \frac{dk}{2\pi} e^{-[E_n(k)+ikv-(n-1)sv]t}$$

- Using the saddle-point equations, we get the entanglement growth rate

$$\Gamma_n(s) = \frac{E_n(-(n-1)is)}{(n-1)s_{\text{eq}}}$$

- Using $E_3(k)$ from the variational calculation, we find:

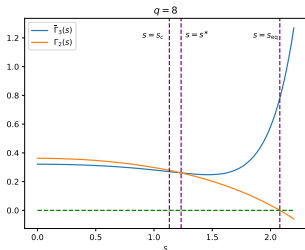


- This seems to give the unphysical prediction that the $\Gamma_3(s_{\text{eq}}) \neq 0!$

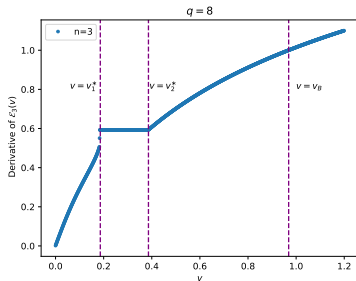
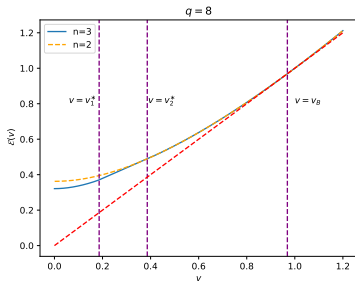
- Unphysical prediction must be corrected by contributions to e^{-2S_3} from some other set of eigenstates of P_6 .
- We can argue that there is another natural set of eigenstates of P_6 , such that

$$e^{-2S_3(x,t)} \propto e^{-2s_{\text{eq}}\bar{\Gamma}_3(s)t} + e^{-2s_{\text{eq}}\Gamma_2(s)t}$$

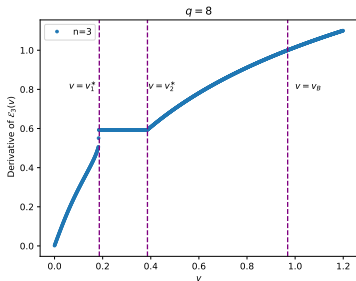
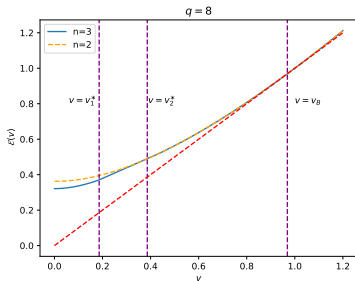
$$\Rightarrow \Gamma_3(s) = \min(\bar{\Gamma}_3(s), \Gamma_2(s))$$



- $\mathcal{E}(v)$ is related to $\Gamma(s)$ by Legendre transformation. We find that $\mathcal{E}_3(v)$ has two phase transitions, of first and second order respectively.



- $\mathcal{E}(v)$ is related to $\Gamma(s)$ by Legendre transformation. We find that $\mathcal{E}_3(v)$ has two phase transitions, of first and second order respectively.



- For $v > v_2^*$, $\mathcal{E}_3(v) = \mathcal{E}_2(v)$.

In particular v_B is the same for $n = 2$ and $n = 3$. (Independent hints from holography that it should be the same for all n .)

Is there a physical reason for this?

Summary and further questions

- In Brownian models without conserved quantities, the membrane picture is a result of gapped low-energy modes that resemble plane waves of domain walls between permutations.

Questions:

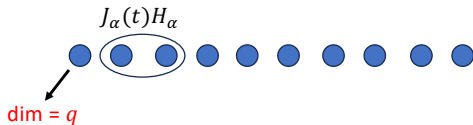
- How does this picture generalize to finite temperature?
 - How does the picture change in Brownian circuits with conserved quantities? *work in progress with Sanjay Moudgalya*
 - How can a similar set of modes emerge in systems without random averaging? Can they be seen in holographic CFTs? *work in progress with Mark Mezei and Zhencheng Wang*
- What is the physical interpretation of the phase transitions in the higher Renyi membrane tensions?
- Can we quantitatively analyse the higher-dimensional case?
- Can these modes be used to formulate an effective field theory of hydrodynamics for entanglement?

Thank you!

$$\mathcal{E}(v) = \max_s \left(\frac{vs}{s_{\text{eq}}} + \Gamma(s) \right)$$

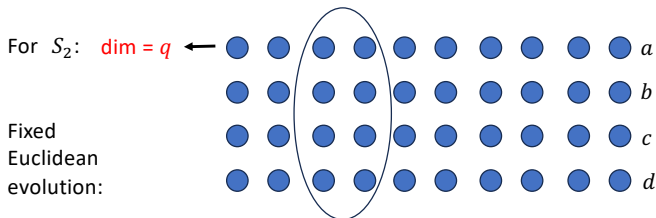
$$\mathcal{E}_3(v) = \begin{cases} \bar{\mathcal{E}}_3(v) & v \leq v_1^* \\ \bar{\Gamma}_3(s_*) + \frac{s^*}{s_{\text{eq}}} v & v_1^* \leq v \leq v_2^* \\ \mathcal{E}_2(v) & v \geq v_2^* \end{cases}$$

Random Lorentzian evolution:



average over $J_\alpha(t)$

For S_2 : $\dim = q$



Fixed Euclidean evolution:

Random GUE $H_\alpha(t)$

In terms of

$$|\uparrow\rangle = |\text{MAX}\rangle_{ab} |\text{MAX}\rangle_{cd}, \quad |\downarrow\rangle = |\text{MAX}\rangle_{bc} |\text{MAX}\rangle_{ad}$$

Fixed Euclidean evolution:

