

emergence of hydrodynamics / classical simulations of quantum transport

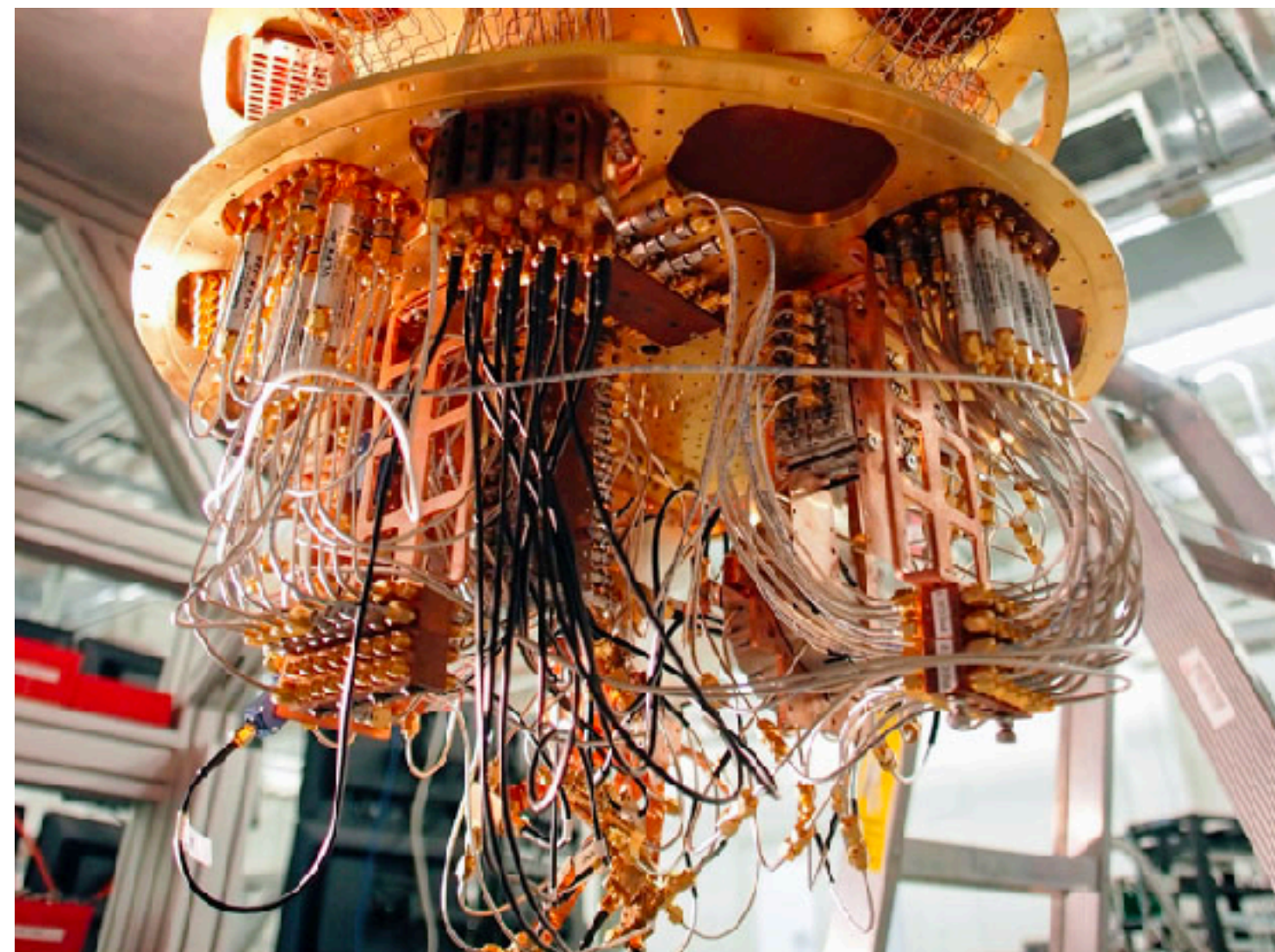
sarang gopalakrishnan (princeton)

limitations of phenomenological hydro

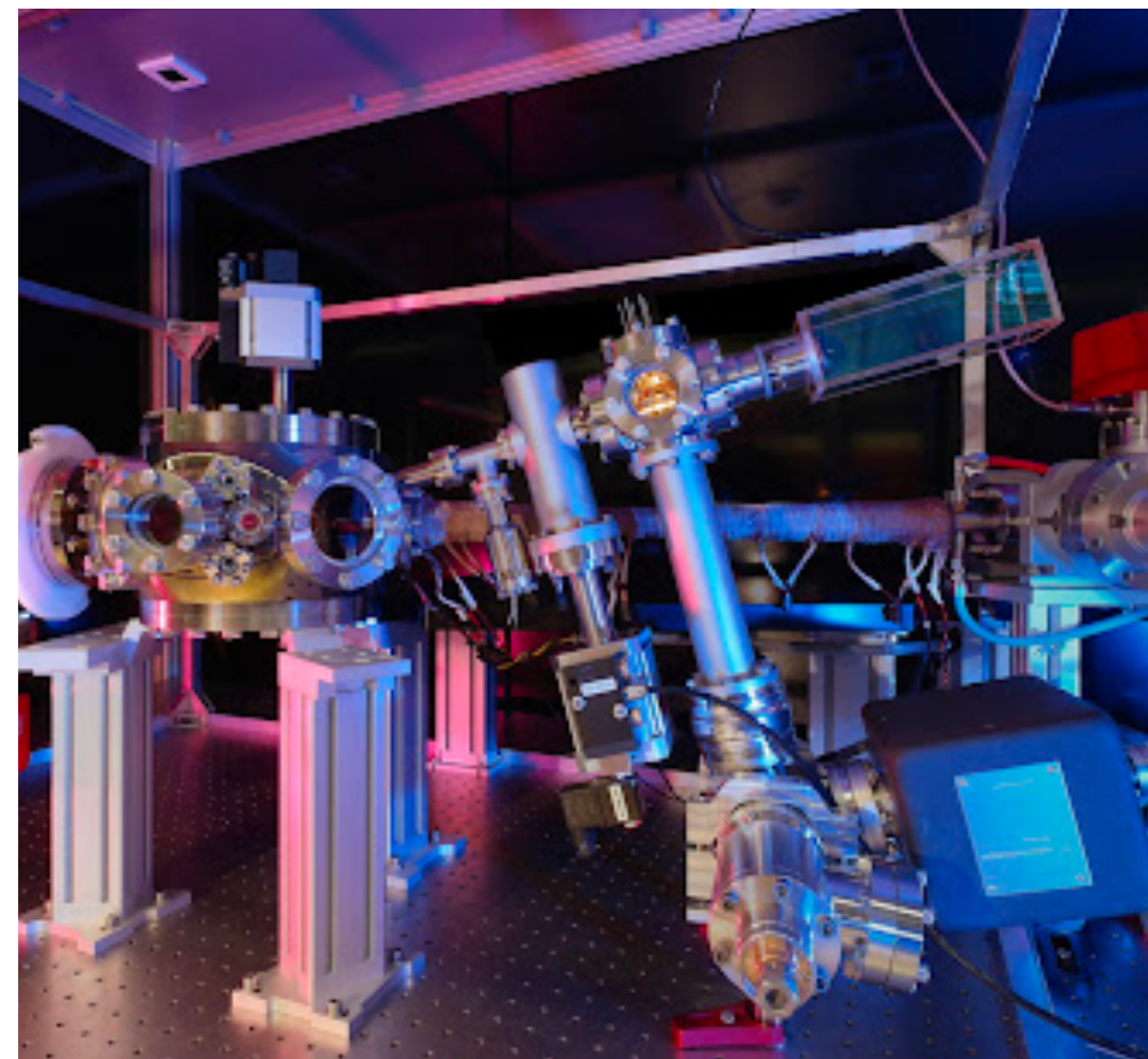
- “Standard” hydro philosophy:
 - Start with symmetries, identify conserved charges and Goldstone modes
 - Write down all terms in gradient expansion, with general coefficients
 - Gives dynamical implications of a particular symmetry structure, could allow one to infer symmetries from experiment
- Basic limitation: disconnected from microscopics
 - Attempts to derive hydro from microscopics: Boltzmann kinetic theory, etc., controlled in specific cases (e.g., weakly interacting systems)
 - But we care about generic strongly interacting systems
- Closely related question: how hard is it to classically predict quantum dynamics?

worst-case scenario for numerics?

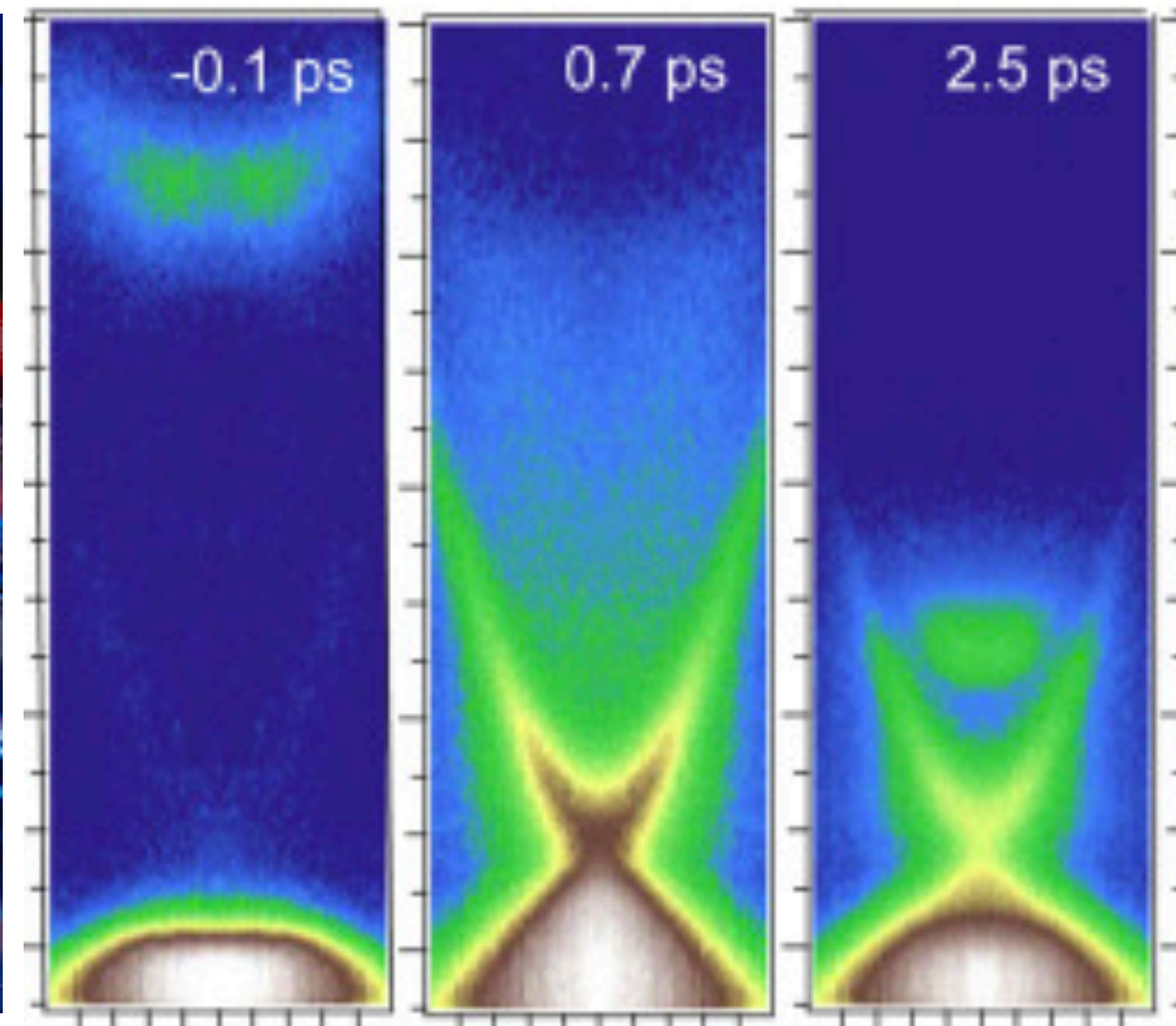
- Tensor-network methods work well when entanglement is low
- Dynamics starting from states that are high-energy-density (far from ground state) generate a lot of entanglement (exponentially hard in time to simulate the state exactly)
- Are we stuck?



superconducting qubit arrays



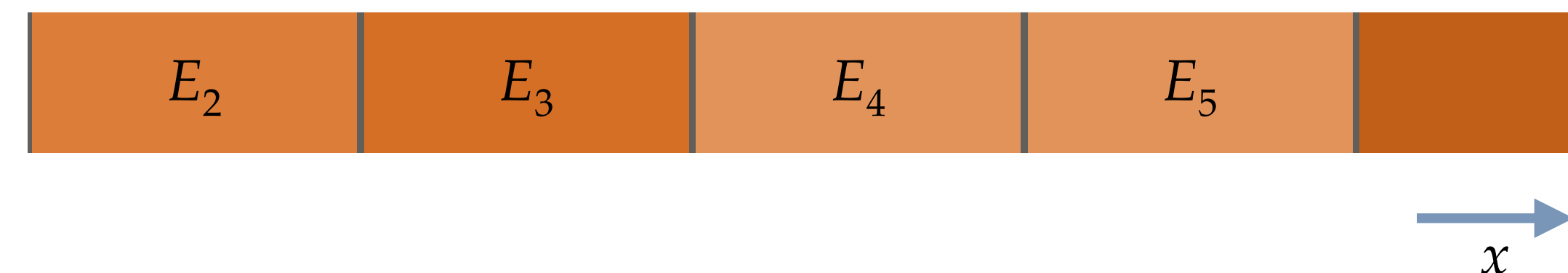
ultracold atoms



ultrafast probes of materials

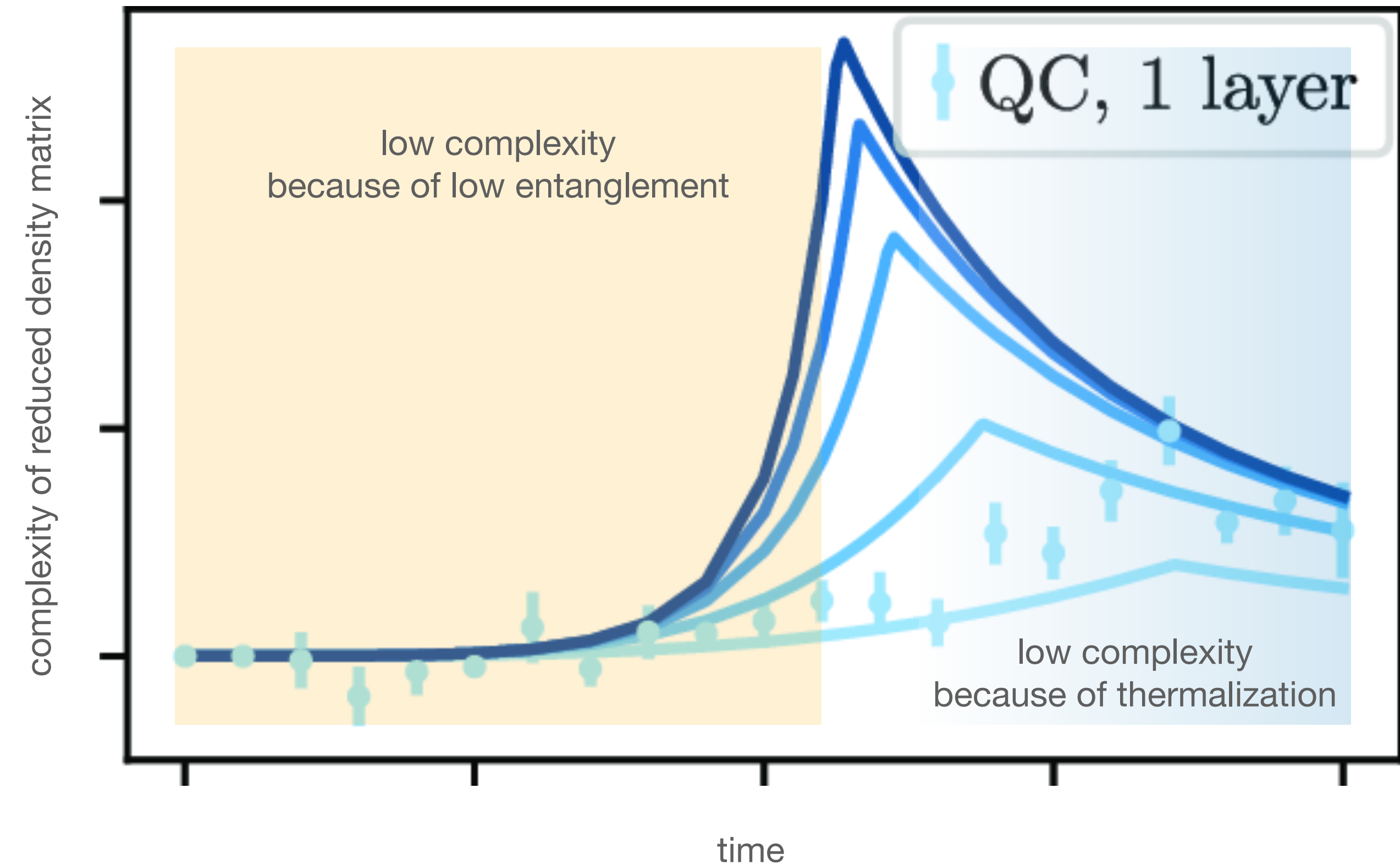
thermalization and hydrodynamics

- High temperature dynamics is complex, chaotic
- Chaos leads to effective randomization of state
 - Local subsystems have thermal density matrices, full system does not but we only care about local properties
- System goes to maximum entropy state subject to conservation laws (“thermalization”)
- Hydrodynamics:
 - Assume system is *locally* in some thermal state (described by local values of conserved variables)
 - Write down equations of motion for conserved variables by gradient expansion (assume that variations are smooth)



entanglement barrier

- Entanglement grows linearly in time, complexity is exponential in entanglement
- High-temperature thermal states like $\exp(-\beta H)$ have low complexity
- Thermalization: reduced density matrix for subsystem can be replaced by thermal state
- Problem: how to bridge between the early and late time physics? (“Crossing the entanglement barrier”)
- Equivalent to asking: how do we derive hydrodynamics?






why do we want to do this?

- Intellectual satisfaction
- Might want precise estimates of diffusion constant, etc. (Trivedi et al., 2024)
- We are interested in specific models, we want to be open to discovering new structures beyond what hydro puts in (e.g., quantum scars, strong zero modes, many-body localization)

Article | Published: 30 November 2017

Probing many-body dynamics on a 51-atom quantum simulator

[Hannes Bernien](#), [Sylvain Schwartz](#), [Alexander Keesling](#), [Harry Levine](#), [Ahmed Omran](#), [Hannes Pichler](#), [Soonwon Choi](#), [Alexander S. Zibrov](#), [Manuel Endres](#), [Markus Greiner](#) , [Vladan Vuletić](#)  & [Mikhail D. Lukin](#) 

[Nature](#) **551**, 579–584 (2017) | [Cite this article](#)

finding slow operators

- Picture behind hydro:
 - Most operators relax fast and can be treated as white noise
 - Remaining operators form a “slow operator Hilbert space”
 - Want to perform “adiabatic elimination”/Schrieffer-Wolff to eliminate the fast operators

- Naive idea: work with Liouville superoperator:

$$\mathcal{L}(O) = i[H, O]$$

- Annoyingly, the eigenmodes of \mathcal{L} are just $|E_i\rangle\langle E_j|$, eigenstates of the true evolution; too nonlocal to be helpful

the rest of this talk

- Noise as a way to cut off the entanglement barrier
- Case with no conservation laws: Liouvillian gap and Ruelle resonances
- Case with charge conservation; hydrodynamic projections
- Note: for simplicity we will be working with discrete-time/Floquet dynamics
- Key refs:

Prosen, J Phys A 35, L737 (2002)

Von Keyserlingk, Pollmann, Rakovszky, arXiv:2111.09904

Nahum, Roy, Vijay, Zhou, arXiv:2205.11544

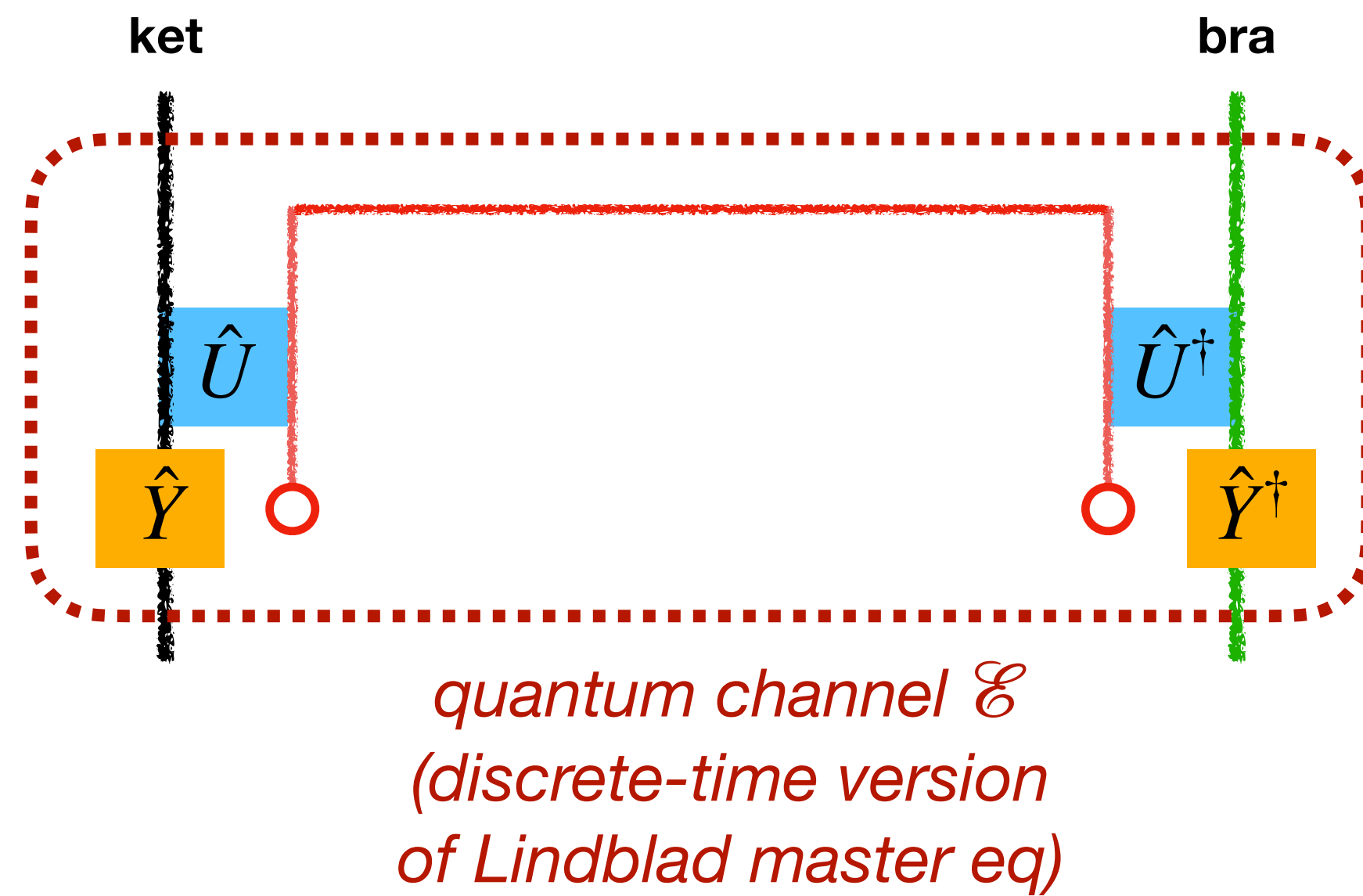
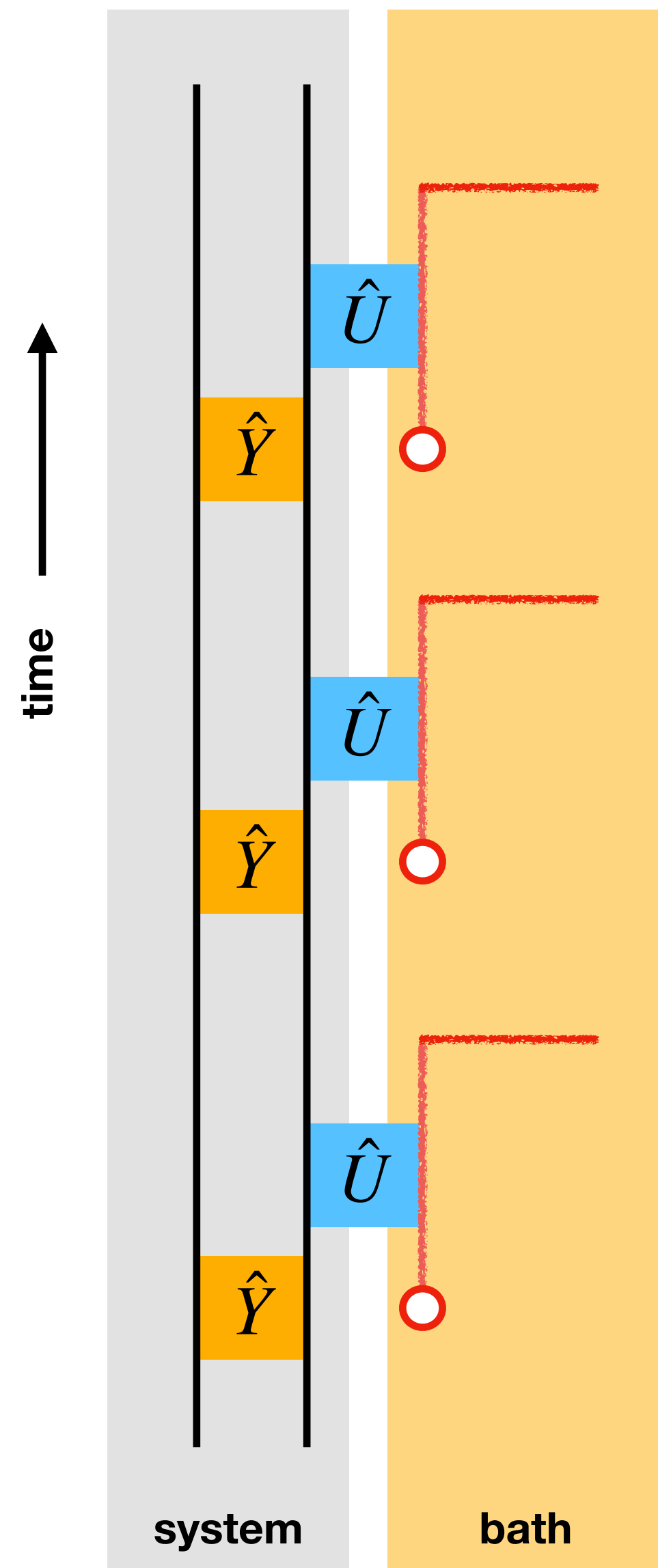
Mori, arXiv:2311.10304

Jacoby, Huse, SG, arXiv:2409.17238

Zhang, Nie, Von Keyserlingk, arXiv:2409.17251

weak dissipation + no conserved quantities

markovian open quantum systems



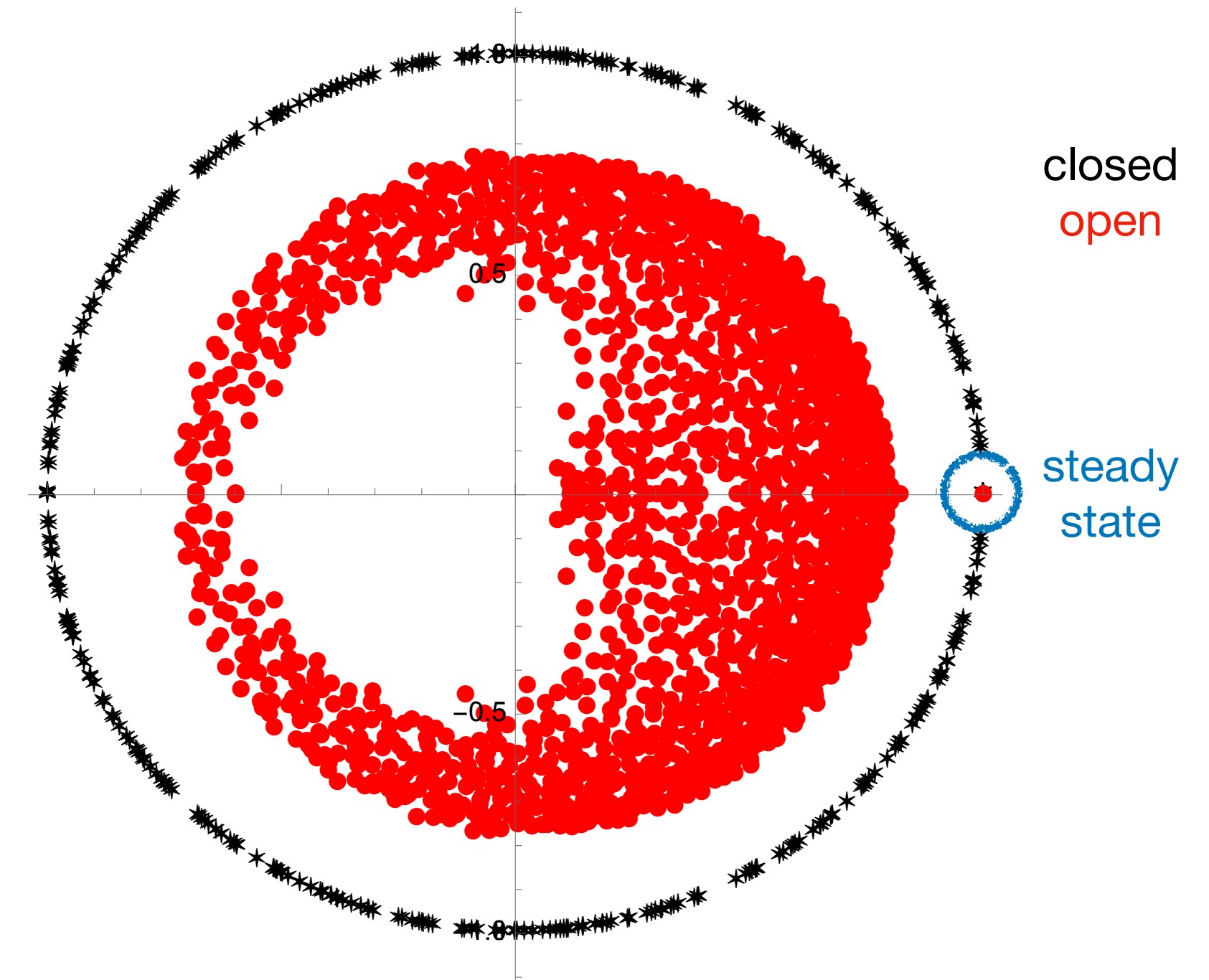
- Q: how does the spectrum of \mathcal{E} evolve in the weak-dissipation limit?

mori's result

- Specifically, consider the following Lindblad master equation with $H(t + \tau) = H(t)$:

$$\partial_t \rho = -i[H(t), \rho] + \gamma \sum_i (\sigma_i^z \rho \sigma_i^z - \rho)$$

- Integrating this equation for a time τ gives one step of evolution under \mathcal{E}

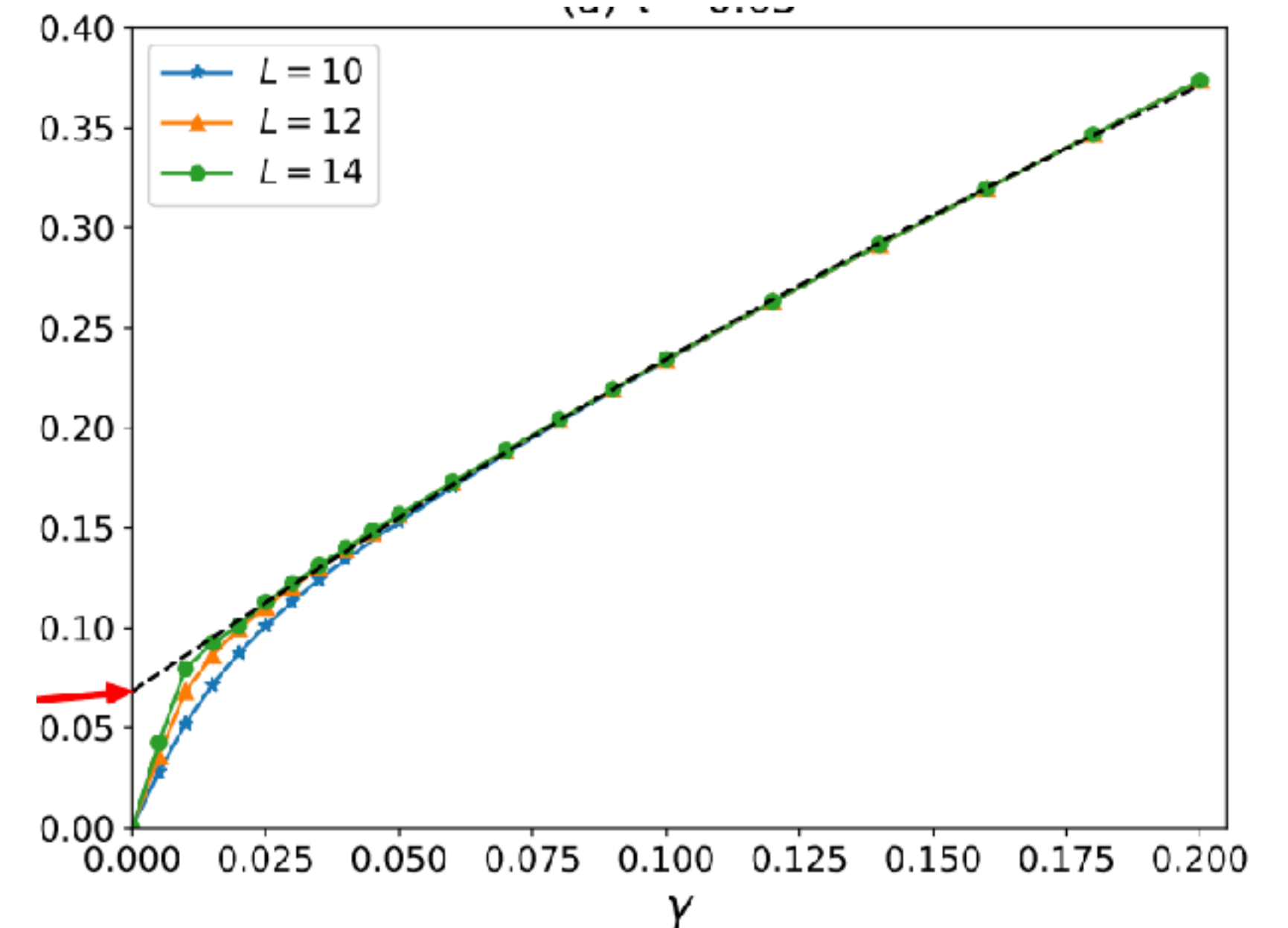
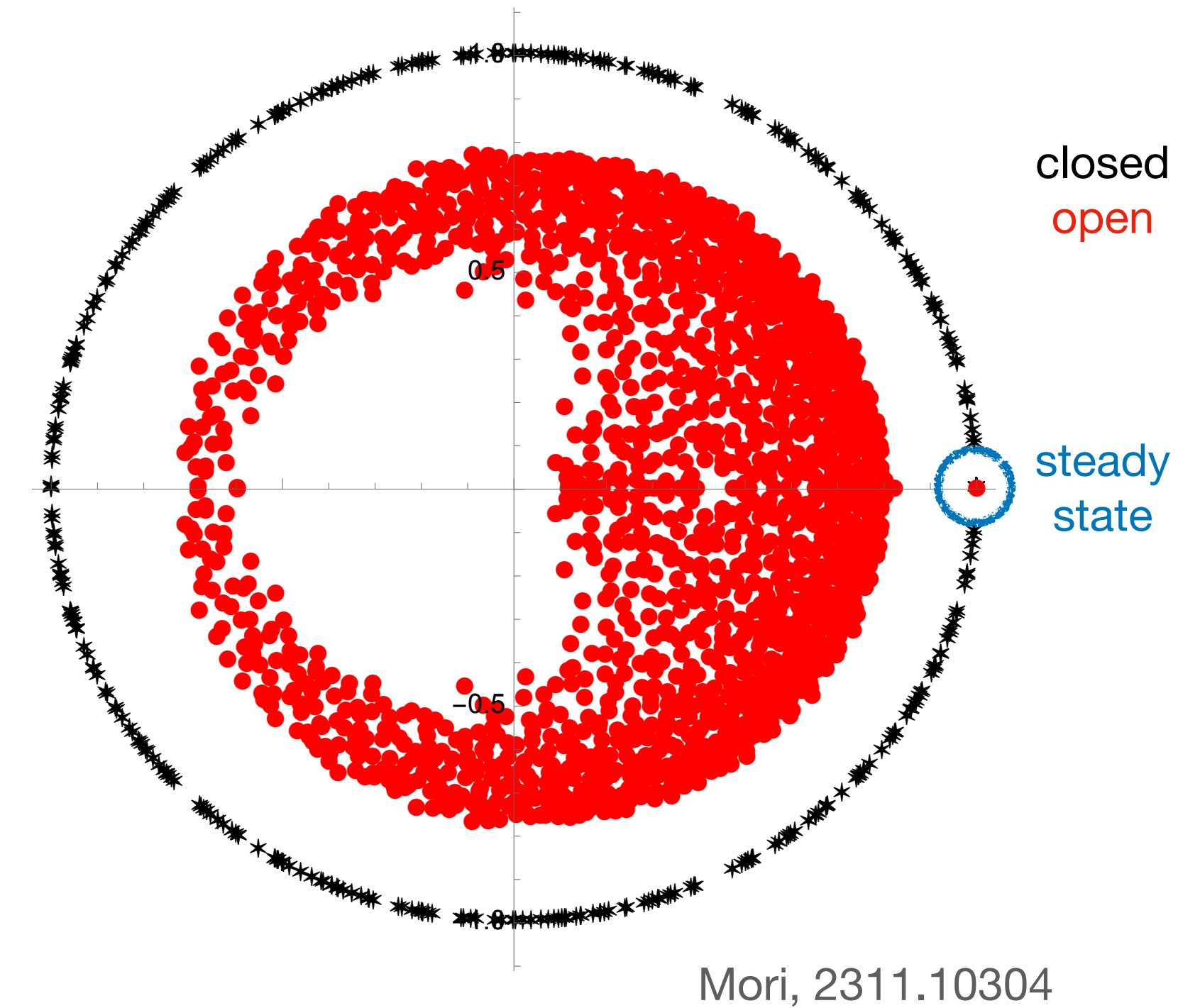


mori's result

- Specifically, consider the following Lindblad master equation with $H(t + \tau) = H(t)$:

$$\partial_t \rho = -i[H(t), \rho] + \gamma \sum_i (\sigma_i^z \rho \sigma_i^z - \rho)$$

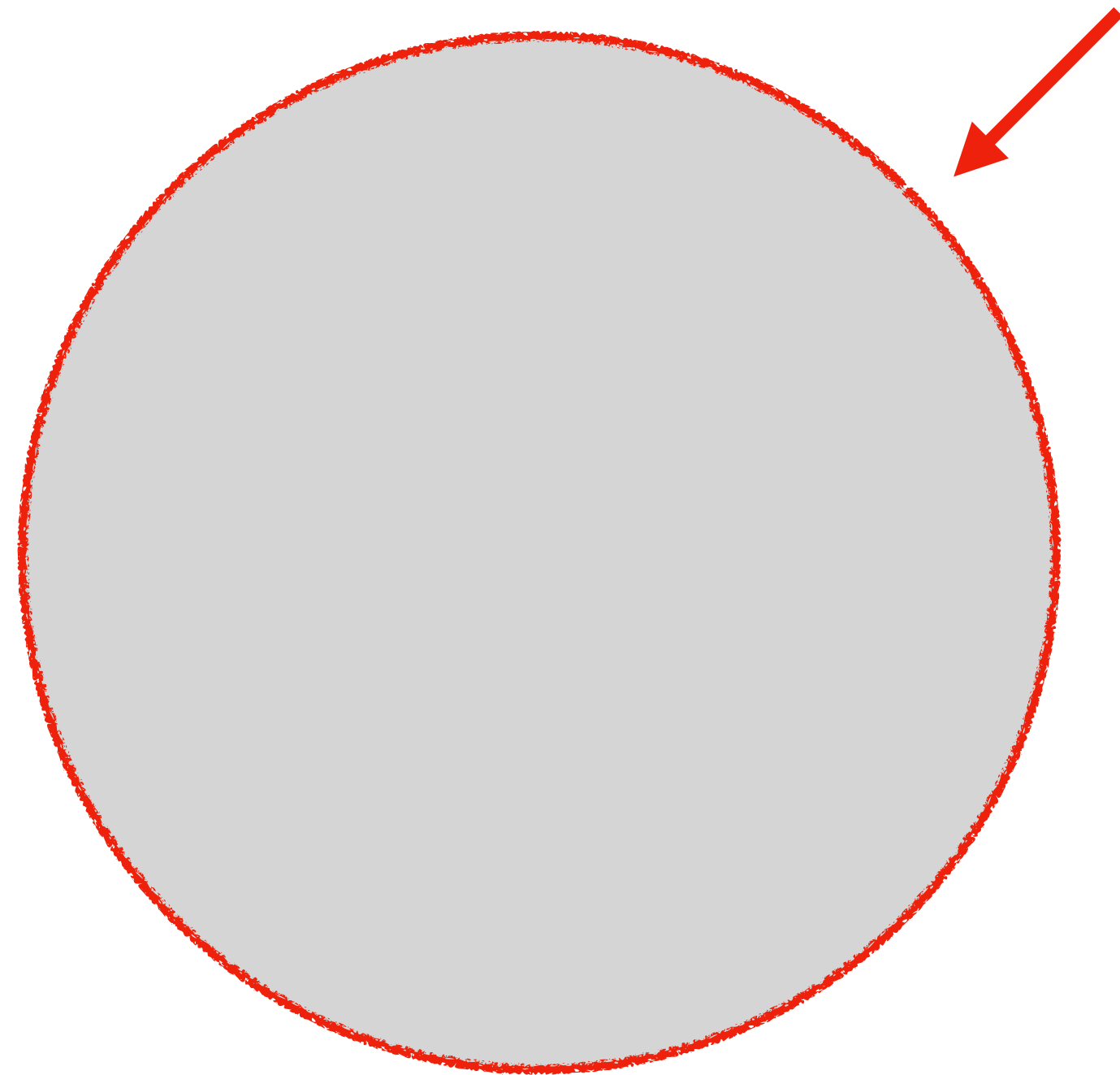
- Integrating this equation for a time τ gives one step of evolution under \mathcal{E}
- Why time-periodic? Otherwise, for $\gamma \rightarrow 0$, H itself becomes a zero mode
- Mori's finding:
 - Gap stays open and $O(1)$ as $\gamma \rightarrow 0$
 - Gap of \mathcal{E} related to decay of local correlation functions



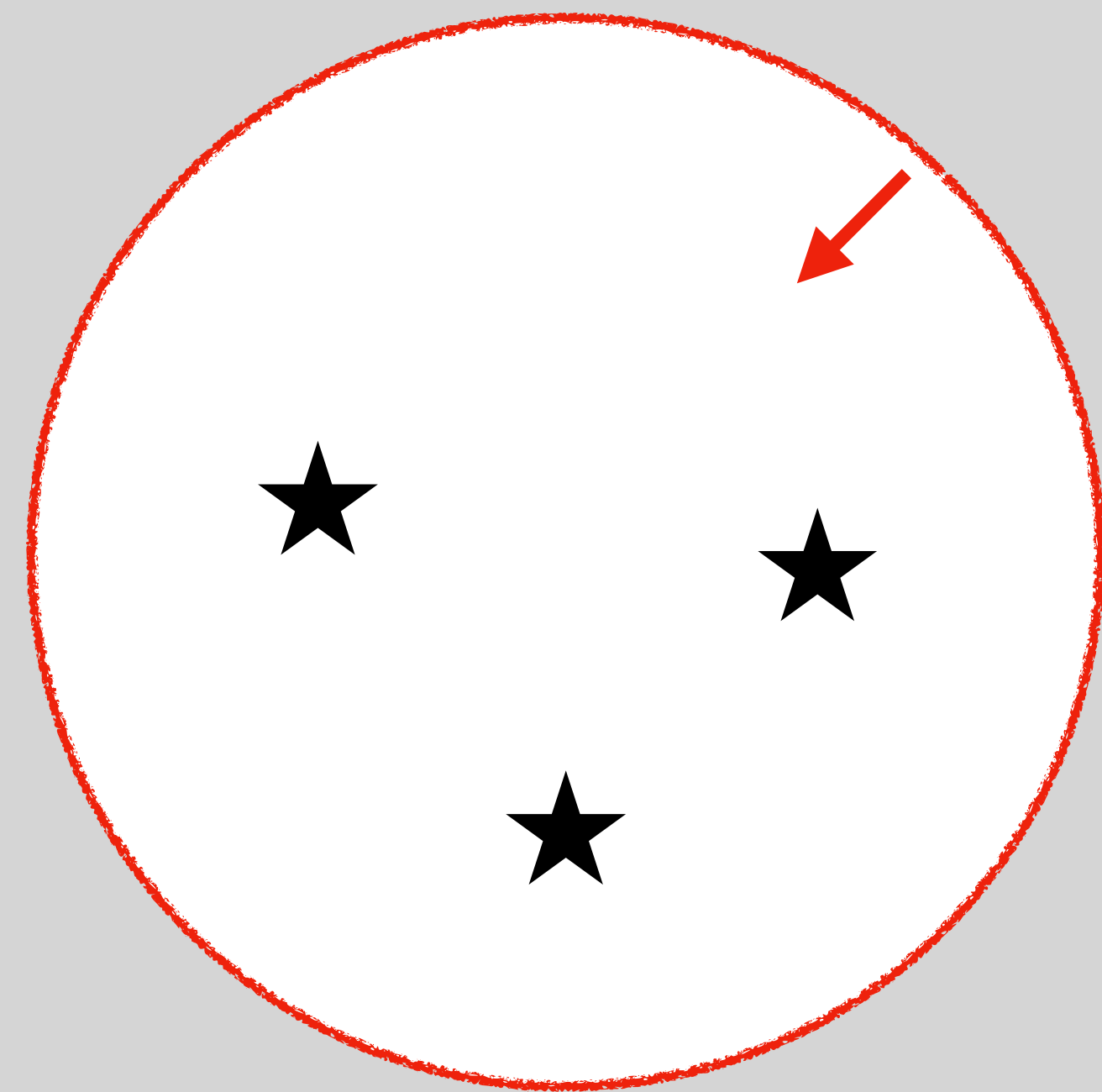
Mori, 2311.10304

ruelle resonances: analytic structure of $(z - \mathcal{E})^{-1}$

first Riemann sheet

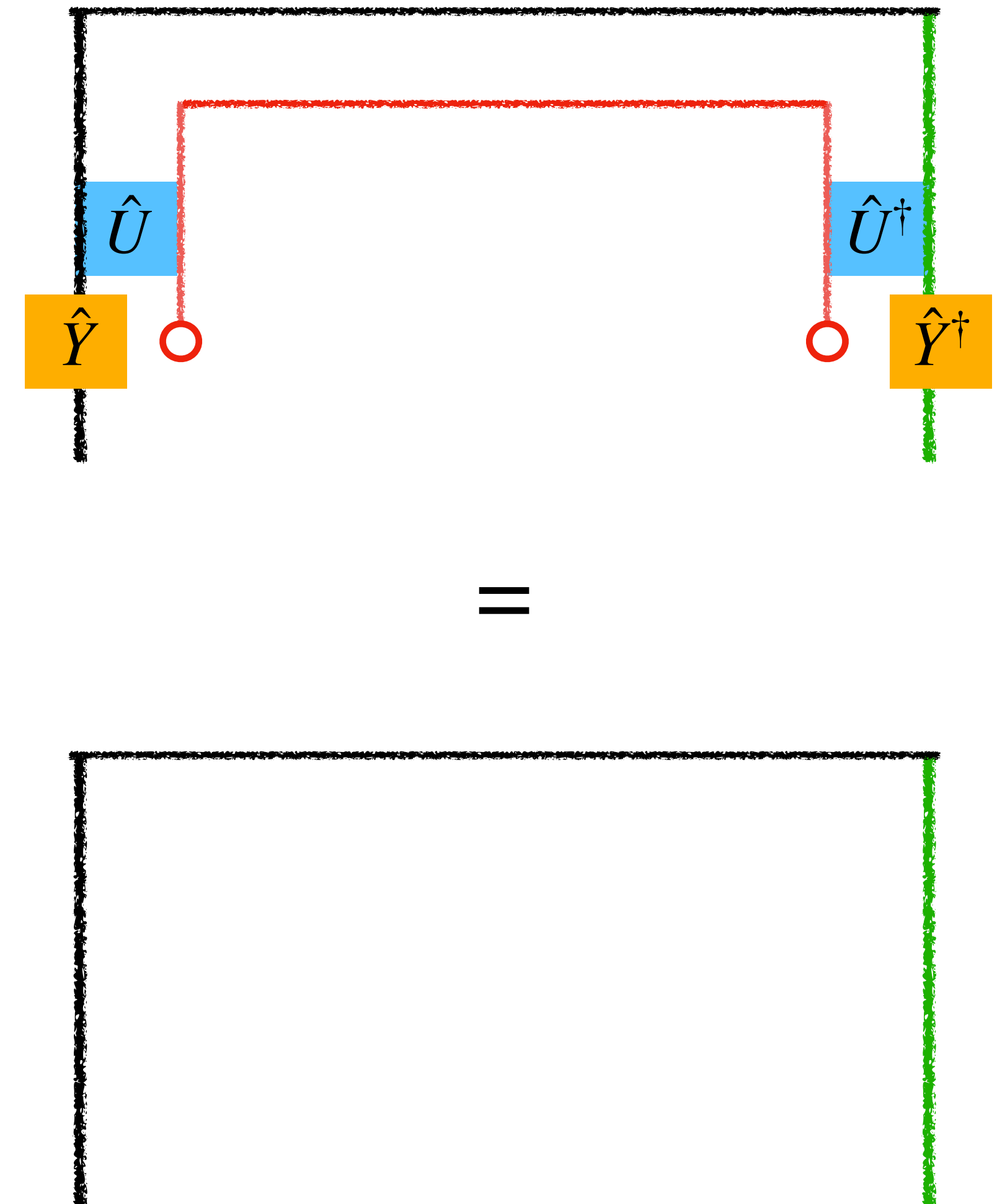


second Riemann sheet



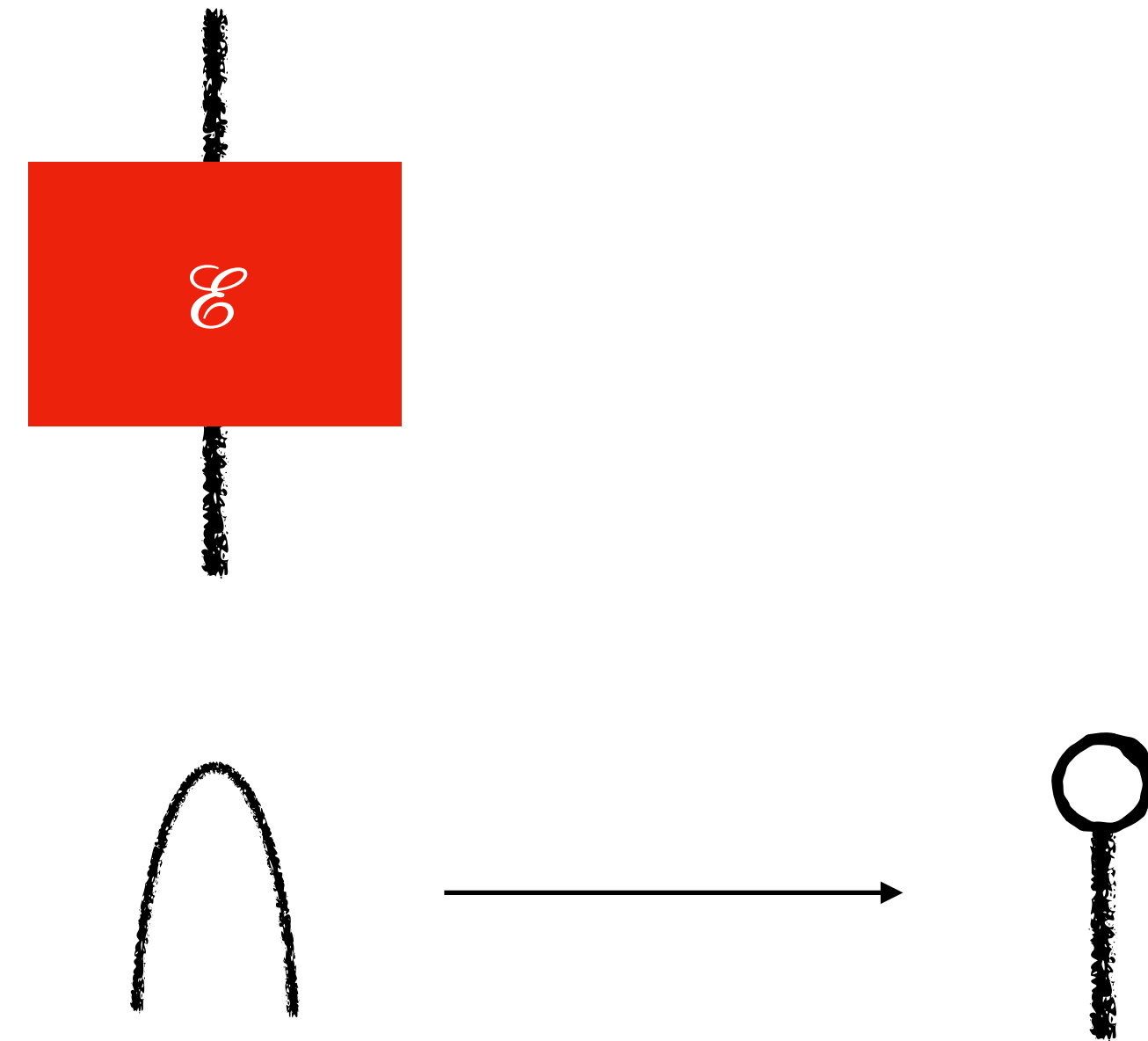
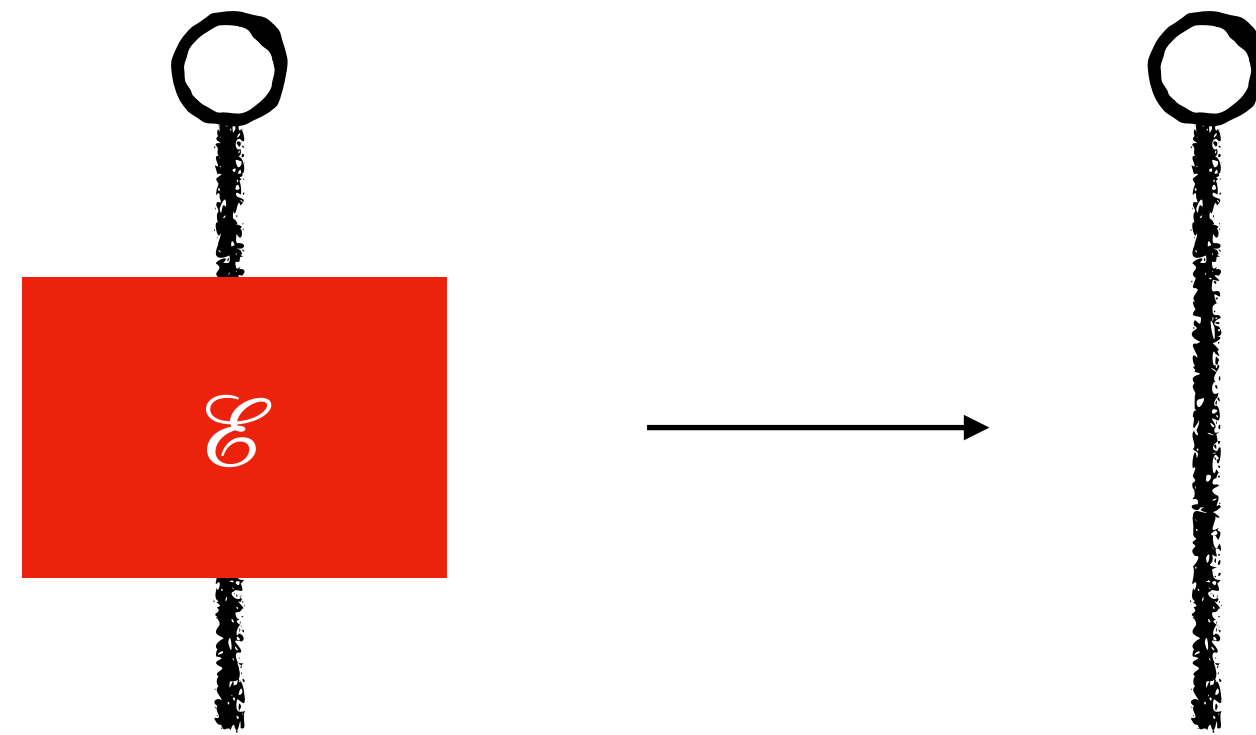
quantum channels in the heisenberg picture

- Density matrix evolves under CPTP map, $\rho \mapsto \mathcal{E}(\rho)$
- We want to compute exp val of some observable, $\langle O(t) \rangle = \text{Tr}(O \mathcal{E}_t(\rho))$
- Can define a Heisenberg picture for CPTP maps, $\text{Tr}(\mathcal{E}^A(O)\rho) \equiv \text{Tr}(O \mathcal{E}(\rho))$
- Helpful picture for superoperators:
 - States are column vectors (“superkets”) $|\rho\rangle$
 - Operators are “superbras” $\langle O|$
 - Bracket $\langle O|\rho\rangle \equiv \text{Tr}(O^\dagger \rho)$
 - Expectation value looks like $\langle O|\hat{\mathcal{E}}|\rho\rangle$
 - Left and right action have the same eigenvalues (but different eigenvectors)
- Trace preservation implies $\langle \mathbb{1}|\hat{\mathcal{E}}|\rho\rangle = \langle \mathbb{1}|\rho\rangle \equiv \text{Tr}(\rho)$, so $\langle \mathbb{1}|\hat{\mathcal{E}} = \langle \mathbb{1}|$, identity is always a left eigenvector of the adjoint map (“unital property”)



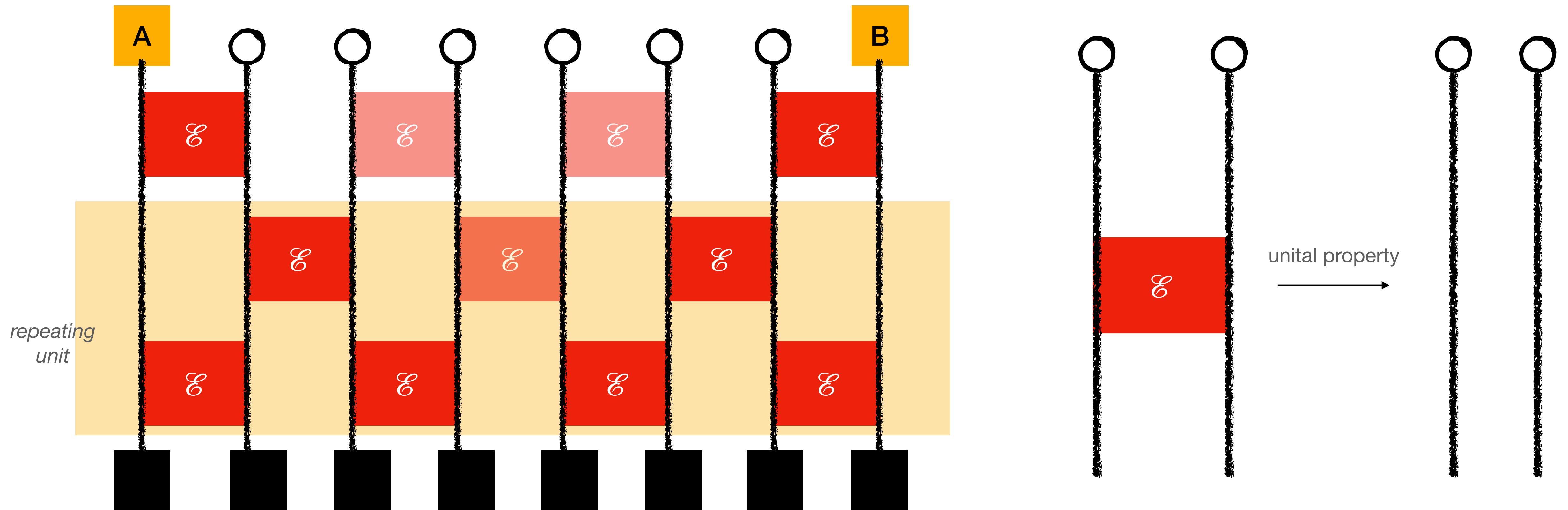
fat-line notation

- Fat lines carry density matrices, a channel is a superoperator on density matrices
- Observables are vectors acting from the top, identity
- Unital property:



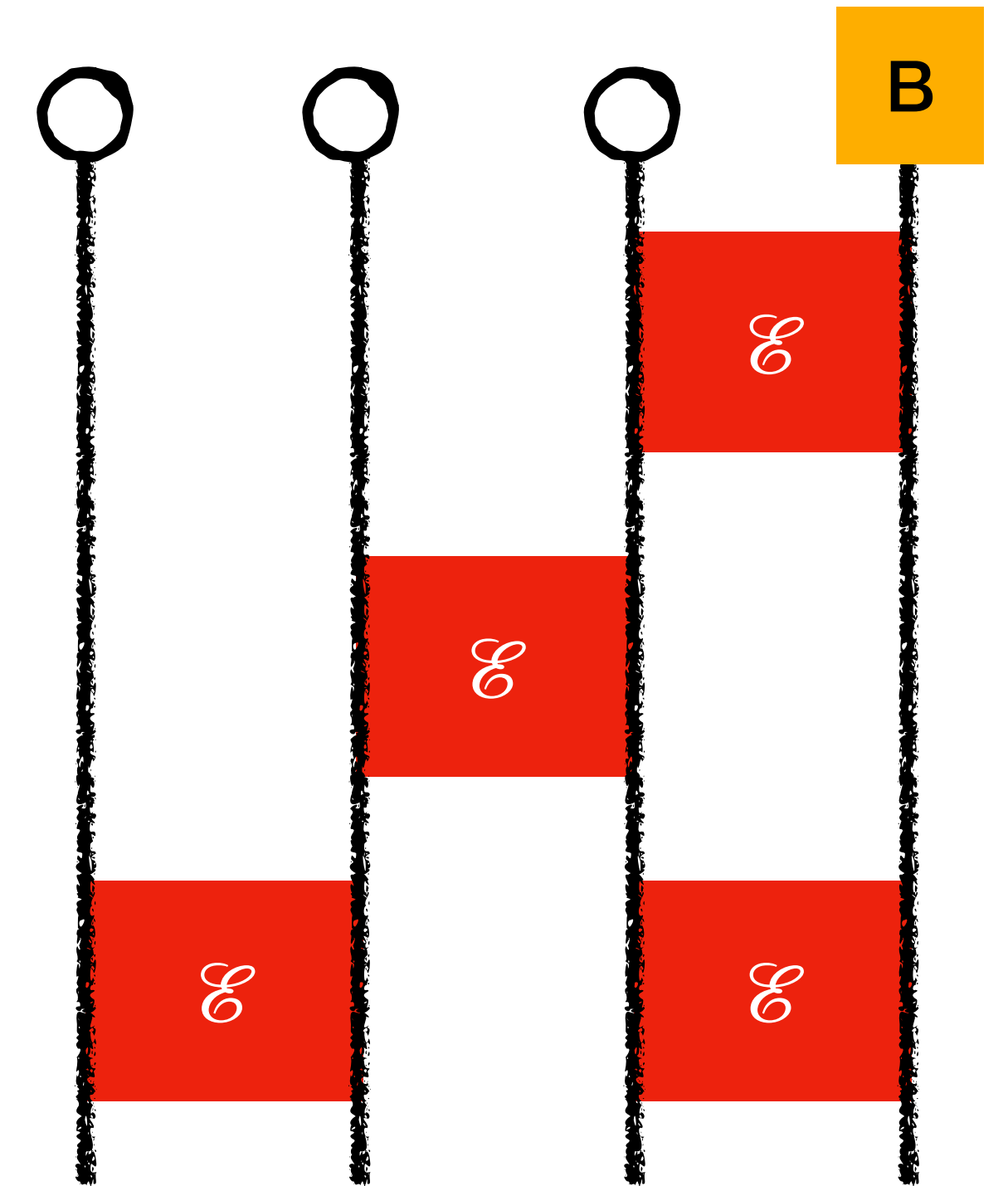
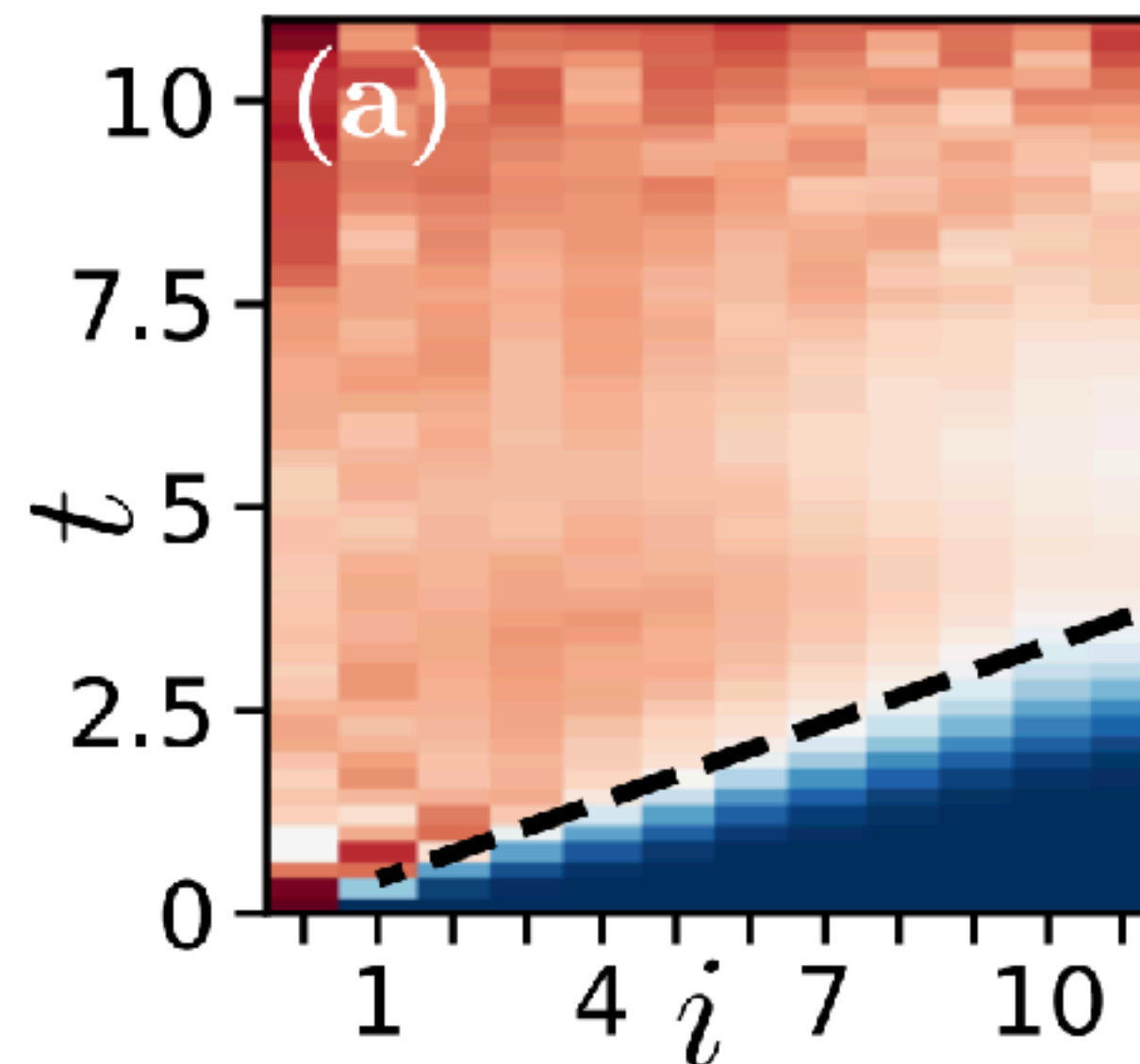
circuits composed of quantum channels

- Lieb-Robinson arguments carry over more or less directly from unitary systems

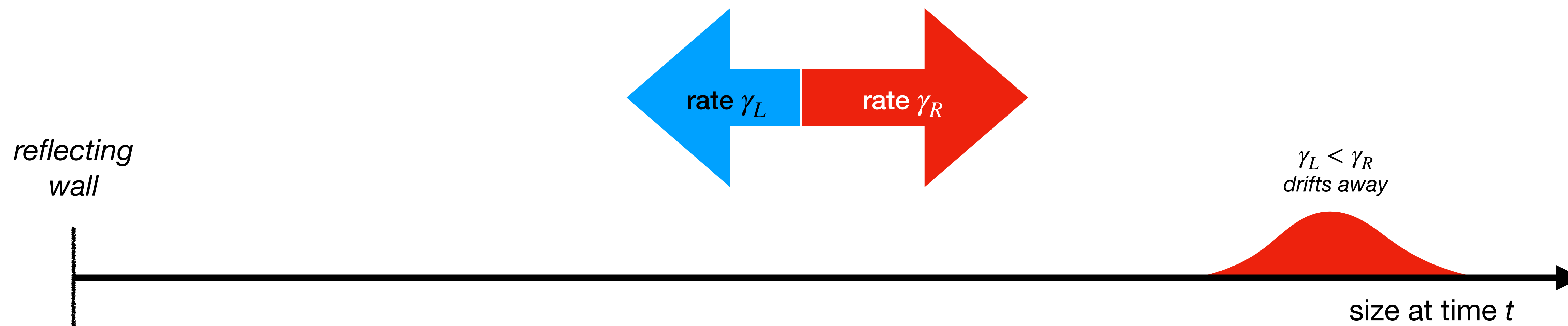


operator growth

- Over time an initially local operator “grows” to fill in its light cone
- Basis for operator space: Pauli strings
- Support of an operator: set of sites with Paulis that are not the identity
- There are four Paulis and only one is the identity, so operator growth is *entropic* in origin
- Can solve explicitly for random quantum circuits



markov process for operator growth



- Operator growth in chaotic systems is a biased random walk (toward larger size)
- Effect of dissipation increases linearly as one goes to larger size, effective process:

$$\dot{\rho}_x = w_+ \rho_{x-1} + w_- \rho_{x+1} - (w_- + w_+ + \gamma x) \rho_x$$

- What are the spectrum and eigenstates of the non-Hermitian operator $M_{x,x'}$?

“hermitian” frame

- $\dot{\rho}_x = w_+ \rho_{x-1} + w_- \rho_{x+1} - (w_- + w_+ + \gamma x) \rho_x$
- Perform a similarity transform $\tilde{M} = T^{-1} M T$, where $T_{x,x'} = e^{ax} \delta_{x,x'}$ (leaves spectrum unchanged)

- For appropriate a , \tilde{M} is a Hermitian matrix

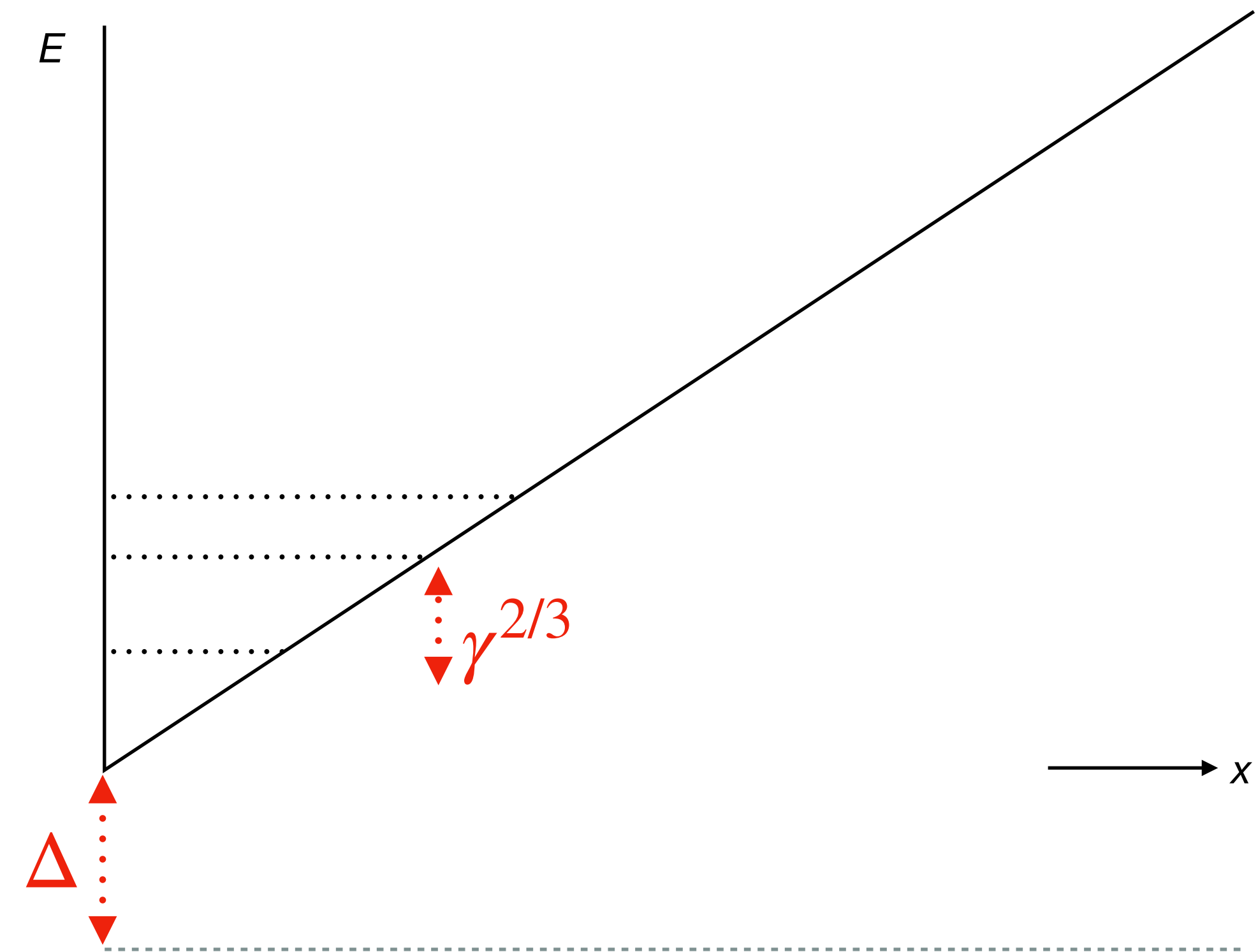
- Continuum limit of eigenvalue problem:

$$(\lambda - \Delta) \psi(x) = -w \psi''(x) + \gamma x \psi(x).$$

$$w = \sqrt{w_+ w_-}, \quad \Delta = \left(\sqrt{w_+} - \sqrt{w_-} \right)^2$$

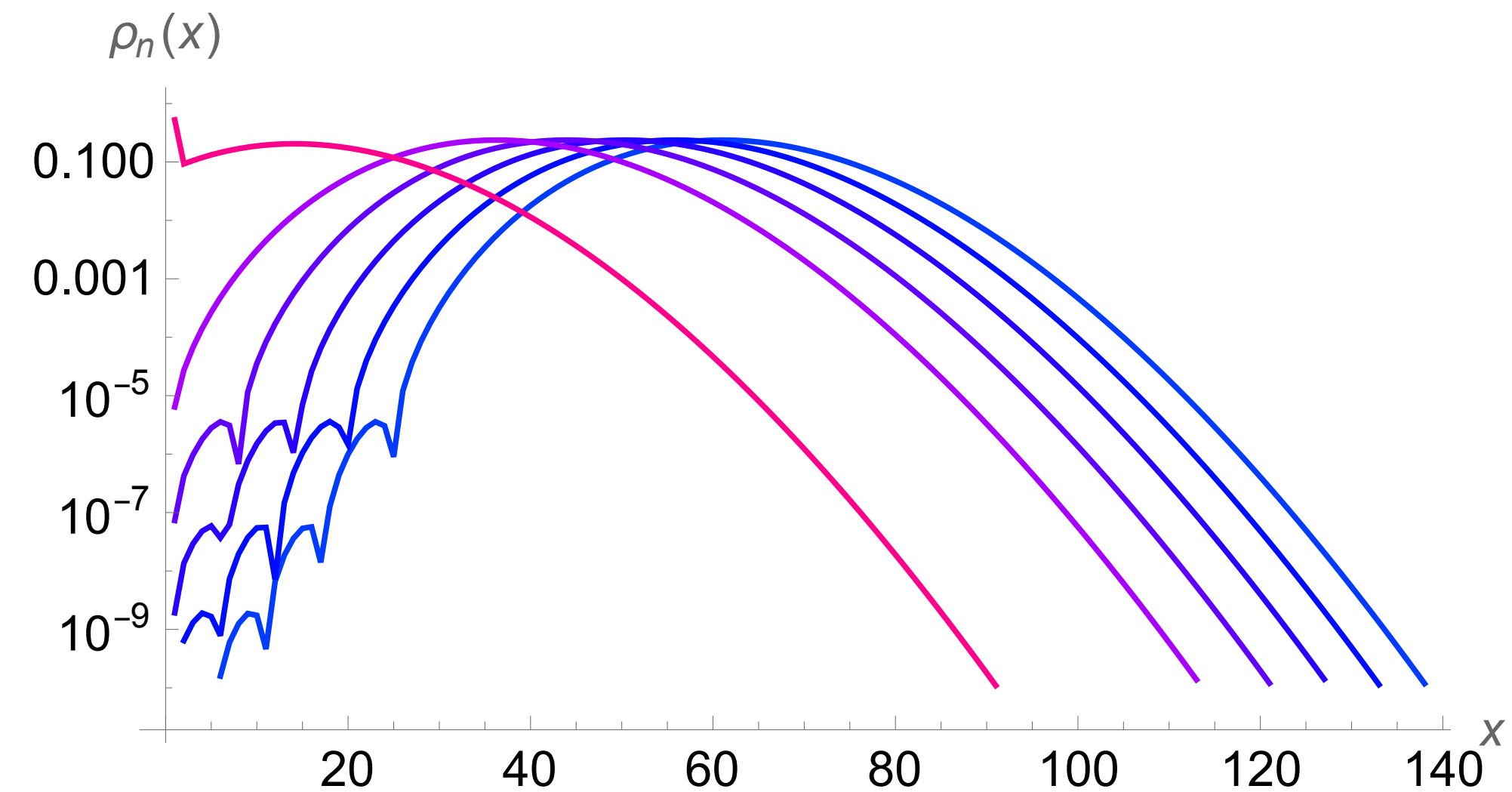
- Gap arises from coordinate transformation

- Eigenstates delocalized (in this frame) on a length-scale $\gamma^{-1/3}$



low-lying eigenstates

- Eigenstates are orthogonal in Hermitian frame, but not after transformation $T|\psi_n^H\rangle = |\phi_n^{lab}\rangle$
- Location of maximum is set by balancing operator growth e^{ax} vs. Airy function decay
- Peak at $1/\gamma$
- Why is the decay rate $O(1)$? Intuitively: eigenmodes have characteristic operator size $1/\gamma$, their decay rate is $\gamma \times 1/\gamma = O(1)$
- Eigenstates are increasingly non-orthogonal as $\gamma \rightarrow 0$
- Bottom line: in the absence of conservation laws, the evolution operator is **gapped!**



correlation functions

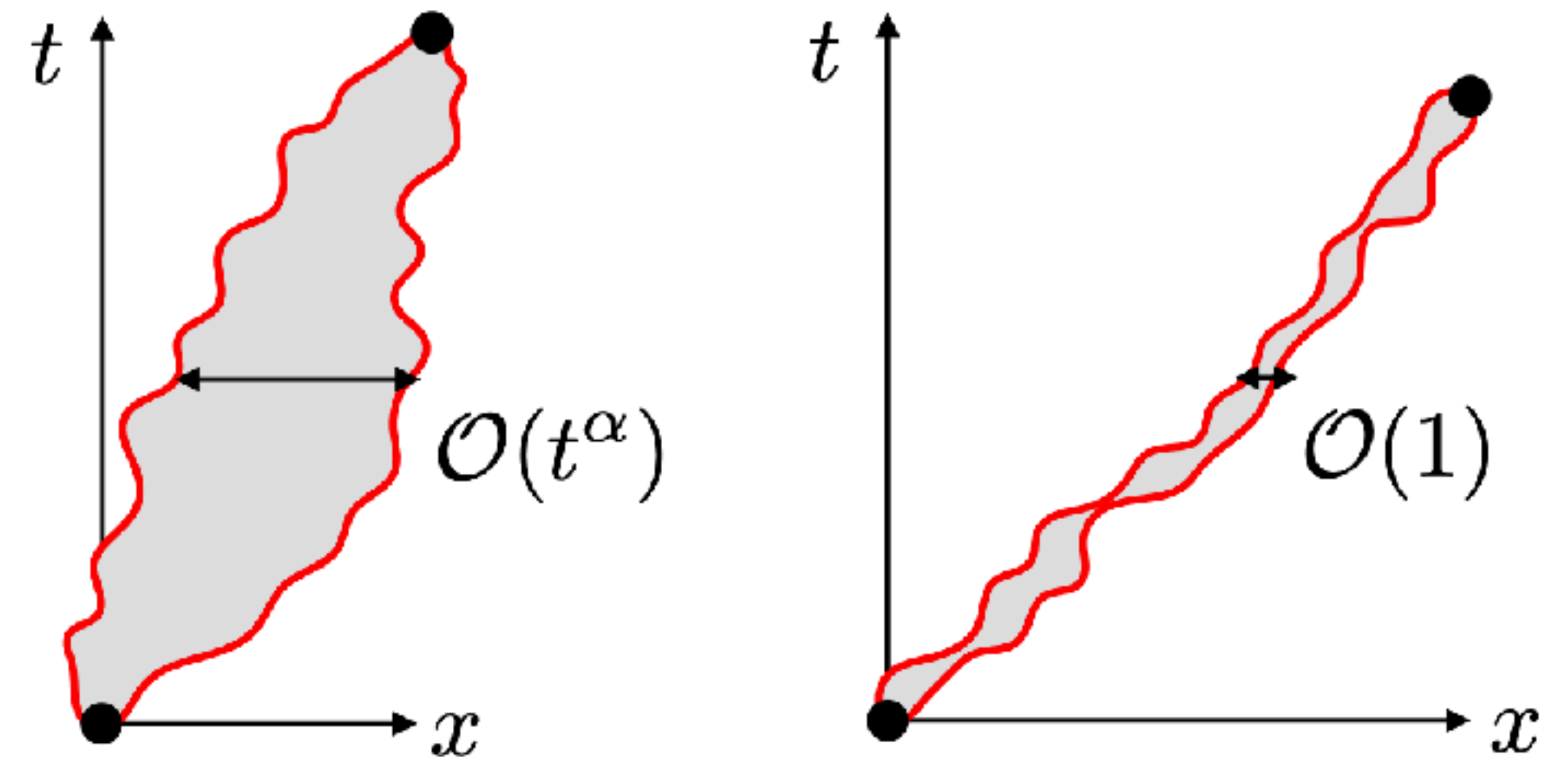
- Return probability for an operator to start and end at size 1

- Recall, generator $M = T\tilde{M}T^{-1}$

where $\tilde{M} = \Delta\mathbb{1} + \text{Schroedinger eq}$

$$\text{and } T = \sum_{xx'} e^{ax} \delta_{xx'}$$

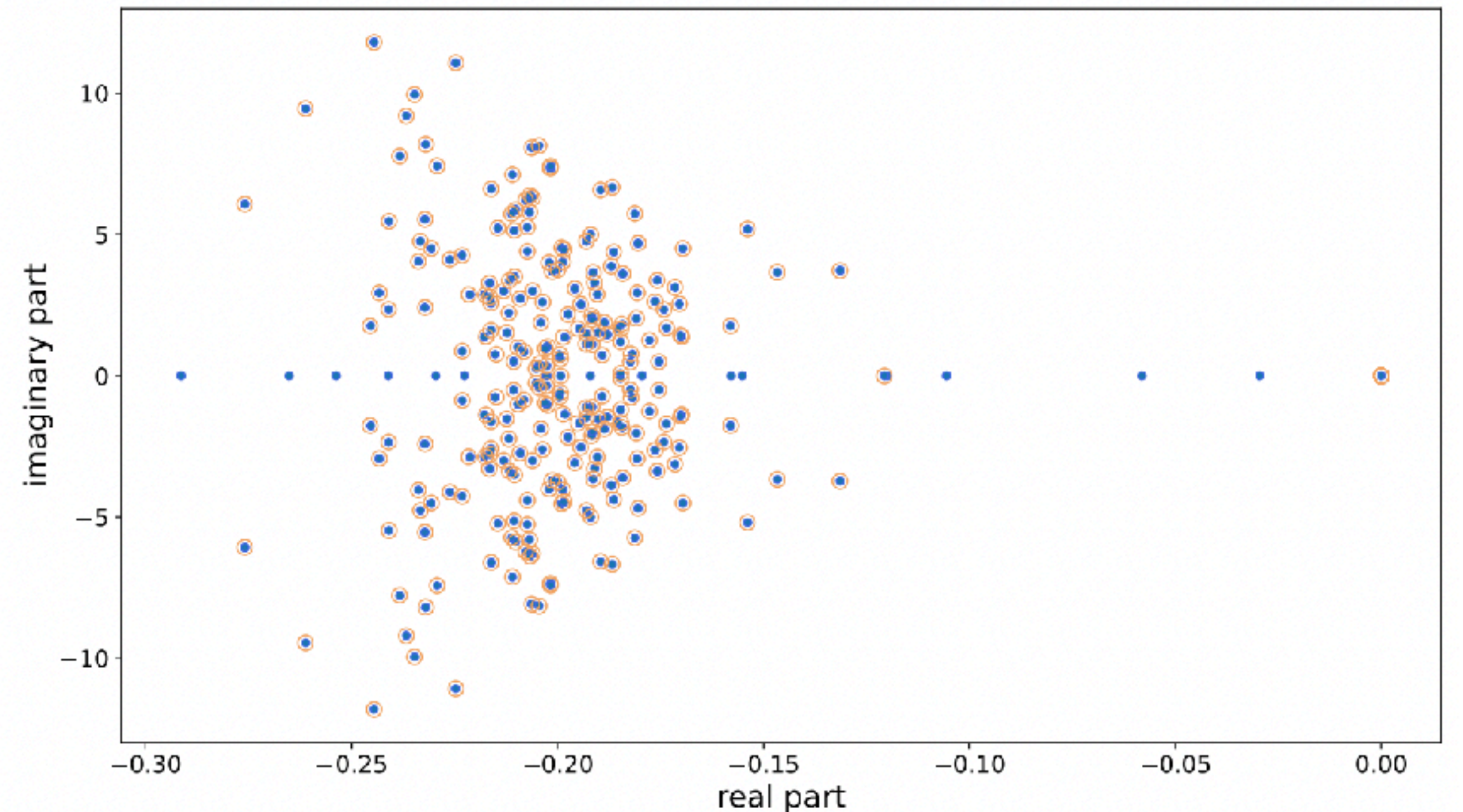
- Correlation function is $\langle 1 | M | 1 \rangle \approx \langle 1 | \tilde{M} | 1 \rangle$
- Because it concerns return probabilities it does not “feel” the change of coordinates
- Bound state vs. lack of bound state \sim Feynman trajectories dominating correlation function (Nahum et al., 2022)



adding conservation laws (**speculative**)

what we hope to see

- Operator Hilbert space separates into hydro and non-hydro subspaces
- Adiabatic elimination of non-hydro subspace: to leading order, simply project it out
- To get the long-time dynamics of a general operator, compute $P_{\text{hydro}}OP_{\text{hydro}}$, which lives in a much smaller space
- Extra slow operators show up as slow modes of \mathcal{E}
- Many practical challenges to making this work...



toy application: mazur bounds

- Suppose we have some number of exactly conserved quantities Q_i
- Then the late-time limit of any autocorrelation function is bounded by the Mazur bound

$$\lim_{t \rightarrow \infty} \langle O(t)O(0) \rangle \geq \sum_i \langle OQ_i \rangle (M^{-1})_{ij} \langle Q_j O \rangle, M_{ij} = \langle Q_i Q_j \rangle$$

- Can construct time-dependent versions (wang, ren, sg, vasseur, 2025): from the positivity of Lehmann representation, we have that

$$\int_0^t dt \langle J(t)J(0) \rangle \geq \langle J_s^2 \rangle t - O(\epsilon)$$

where $\|[J_s, H]\| \leq \epsilon$

- So far, useful mostly in cases where we already know the structure of the problem
- Similar approach has been used to find slow operators: Banuls et al. (2015, 2018)

wrapping up

- Even if exact simulations of quantum states are hard, many of the questions we are interested in might not be
- Hydrodynamics is a successful theory of some observables in some systems
- Where are its boundaries? Can we quantify non-hydrodynamic effects and compute their timescales? (Recall: bit-string distribution has specifically quantum fluctuations)
- Are there universal effects in transport beyond hydrodynamics?
- How do we turn the “hydrodynamic projection” into an efficient numerical scheme?