

Under what conditions do quantum systems thermalize?

Christian Gogolin

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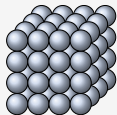
2014-01-20 COST conference

Old questions and new results

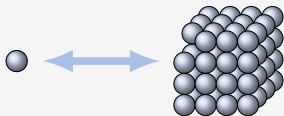
How do **quantum mechanics** and **statistical mechanics** go together?



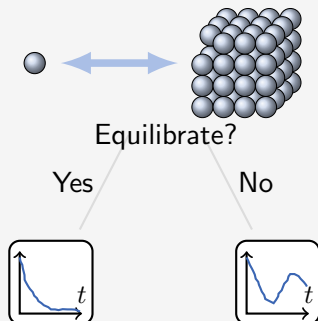
Understanding thermalization



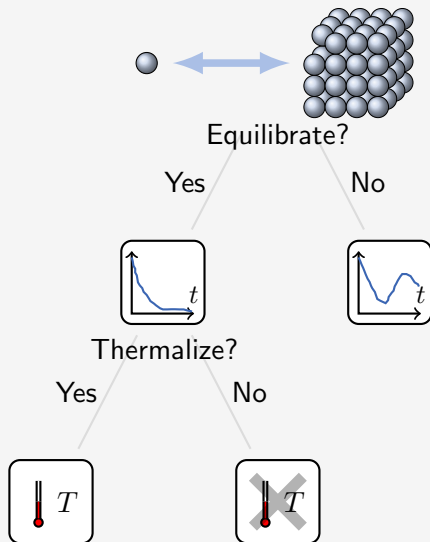
Understanding thermalization



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Setup

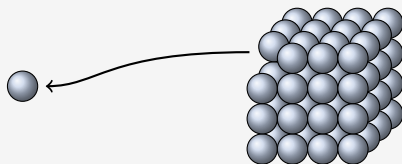
Setup

Subsystem, $\mathcal{H}_S, \mathcal{H}_S$

$$d_S = \dim(\mathcal{H}_S)$$

Bath, $\mathcal{H}_B, \mathcal{H}_B$

$$d_B \gg d_S$$



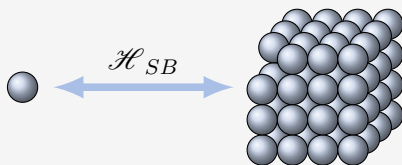
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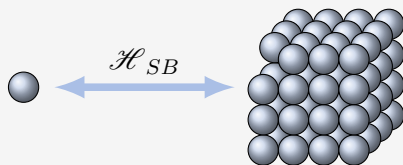
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$$\psi_t^S = \text{Tr}_B[\psi_t]$$

$$\mathcal{H} = \mathcal{H}_S \otimes \mathbf{1} + \mathbf{1} \otimes \mathcal{H}_B + \mathcal{H}_{SB}$$

$$\frac{d\psi_t}{dt} = i[\psi_t, \mathcal{H}]$$

Equilibration

Equilibration

Theorem 1 (Equilibration on average [3])

If \mathcal{H} has *non-degenerate energy gaps*, then for every $\psi_0 = |\psi_0\rangle\langle\psi_0|$ there exists a ω^S such that:

$$\overline{\mathcal{D}(\psi_t^S, \omega^S)} \leq \frac{1}{2} \sqrt{\frac{d_S^2}{d^{\text{eff}}}}.$$

[1] M. Cramer, C. M. Dawson, J. Eisert, and T. J. Osborne, PRL 100, 030602 (2008).

[2] P. Reimann, PRL 101, 190403 (2008).

[3] N. Linden, S. Popescu, A. J. Short, and A. Winter, PRE 79, 061103 (2009).

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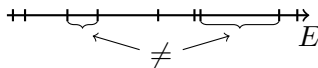
Equilibration

Non-degenerate energy gaps

\mathcal{H} has non-degenerate energy gaps iff:

$$E_k - E_l = E_m - E_n$$

$$\implies k = l \wedge m = n \quad \vee \quad k = m \wedge l = n$$



Intuition: Sufficient for \mathcal{H} to be fully interactive

$$\mathcal{H} \neq \mathcal{H}_1 \otimes \mathbb{1} + \mathbb{1} \otimes \mathcal{H}_2$$

- [1] M. Cramer, *Phys. Rev. Lett.* **102**, 080401 (2009).
- [2] P. Reimann, *Phys. Rev. Lett.* **101**, 190403 (2008).
- [3] N. Linden, S. Popescu, A. J. Short, and A. Winter, *Phys. Rev. Lett.* **79**, 061103 (2009).
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there exists

Effective dimension

$$d^{\text{eff}} = \frac{1}{\sum_k |\langle E_k | \psi_0 \rangle|^4}.$$

Intuition: Dimension of supporting energy subspace

$\langle \psi_0 |$

-
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$$\overline{\mathcal{D}(\psi_t^S, \omega^S)} \leq \frac{1}{2} \sqrt{\frac{d_S^2}{d^{\text{eff}}}}.$$

\implies If $d^{\text{eff}} \gg d_S^2$ then ψ_t^S *equilibrates on average*.

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Maximum entropy principle

Theorem 2 (Maximum entropy principle [6])

If $\text{Tr}[A \psi_t]$ equilibrates on average, it *equilibrates towards its time average*

$$\overline{\text{Tr}(A \psi_t)} = \text{Tr}(A \overline{\psi_t}) = \text{Tr}(A \omega),$$

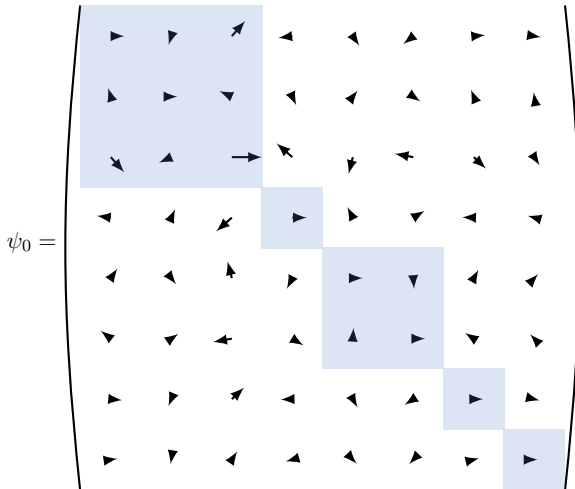
where

$$\omega := \sum_k \Pi_k \psi_0 \Pi_k$$

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Time averaging



Theorem

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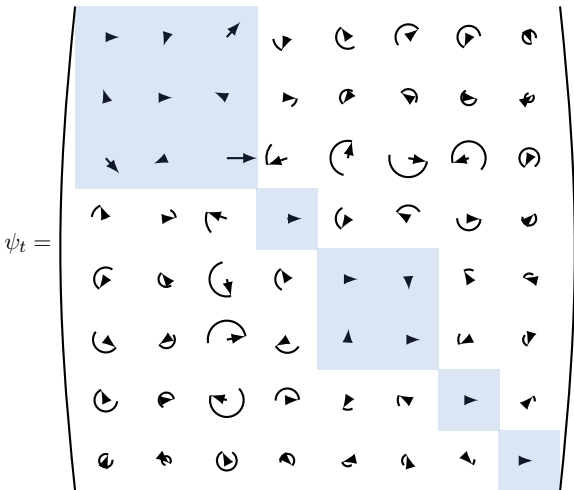
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is the dep
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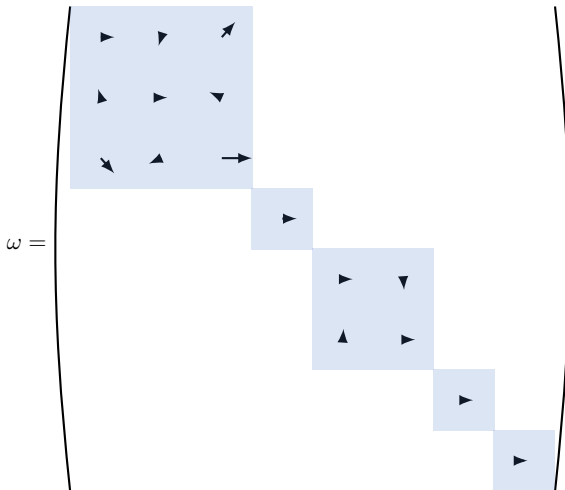
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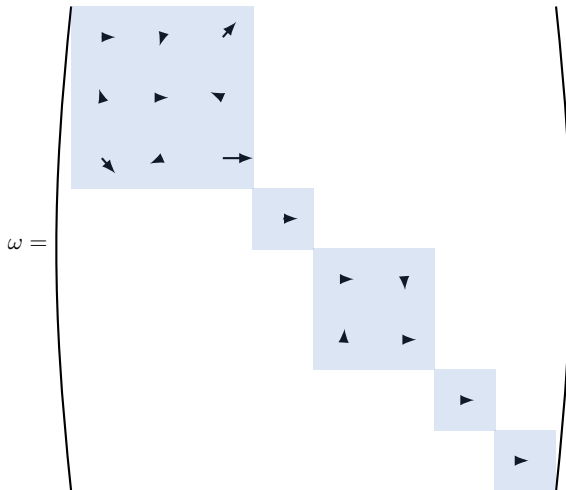
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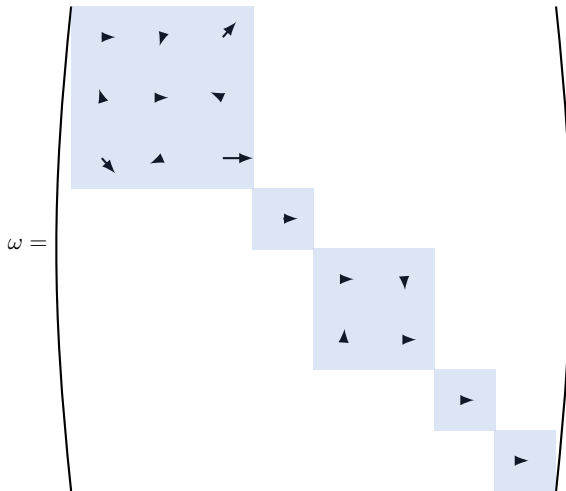
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⇒ Maximum entropy principle from pure quantum dynamics.

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Interesting open questions:

- Do we really need all (exponentially many) conserved quantities?
- If not, then which?
- Does this depend on integrability of the model?
- What is the relation to the GGE?

⇒ Maximum entropy principle from pure quantum dynamics.

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Thermalization

Thermalization is a complicated process

Thermalization implies:

- 1 Equilibration [1, 2, 3, 7]
- 2 Subsystem initial state independence [6, 8]
- 3 Weak bath state dependence [9]
- 4 Diagonal form of the subsystem equilibrium state [10]
- 5 Gibbs state $\omega^S = \text{Tr}_B(\omega) \approx e^{-\beta \mathcal{H}_S}$ [9]

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[9] A. Riera, C. Gogolin, and J. Eisert, PRL 108, 080402 (2012).

[10] C. Gogolin, PRE 81, 051127 (2010).

Thermalization and quantum integrability

There is a common belief in the literature [11, 12, 13, 14, 15] ...

Non-integrable	\implies	Thermalization
Integrable	\implies	No thermalization

[11] C. Kollath et. al PRL 98, 180601 (2007).

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[13] M. Rigol, V. Dunjko, and M. Olshanii, Nature 452, 854 (2008).

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Notions of (non-)integrability

A system with n degrees of freedom is **integrable** if

- there exist n (local) **conserved** mutually commuting linearly/algebraically **independent operators**.
- the system is integrable by the **Bethe ansatz**.
- the system exhibits nondiffractive scattering.
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Lack of imagination?

Absence of thermalization in non integrable systems

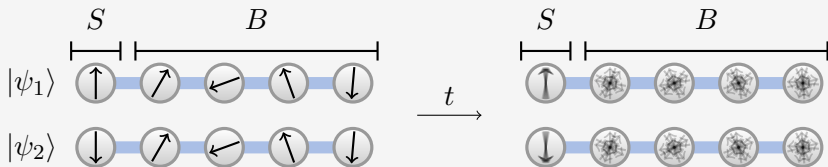
Result (Theorem 1 and 2 in [6]):

- Too little (geometric) entanglement in the energy eigenbasis prevents subsystem initial state independence.
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Absence of thermalization in non integrable systems

The model:

Spin-1/2 XYZ chain with random coupling and on-site field.

$$\mathcal{H} = \sum_{i=1}^n h_i \sigma_i^Z + \sum_{i=1}^{n-1} \vec{b}_i \cdot \vec{\sigma}_i^{\text{NN}}$$

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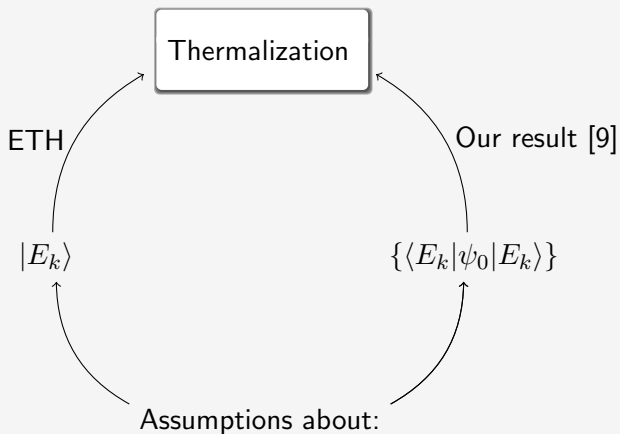
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Interesting open questions:

- What is the relation to **Anderson localization**?
- Can this also happen in **translation invariant** systems?
- Why is integrability a **useful** concept despite the ambiguity?

Conditions for thermalization

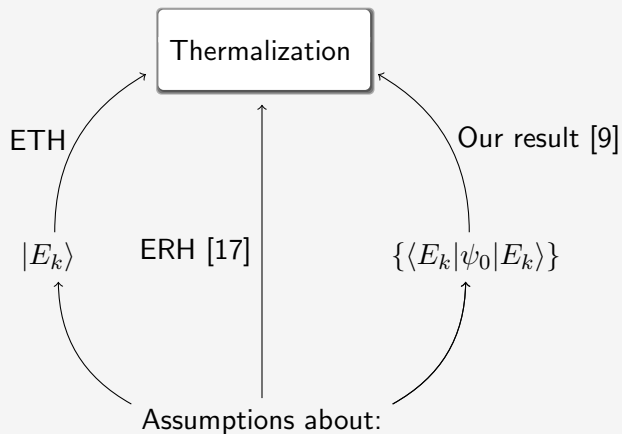
Two ways to prove thermalization



[9] A. Riera, C. Gogolin, and J. Eisert, PRL 108, 080402 (2012).

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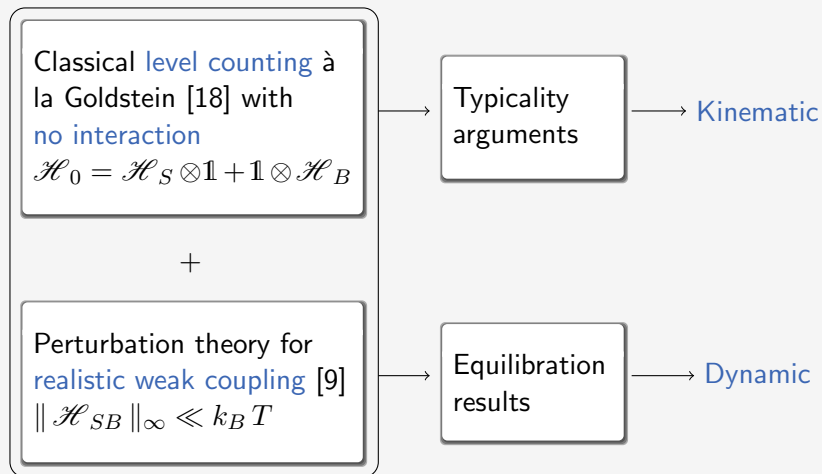
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Structure of the argument

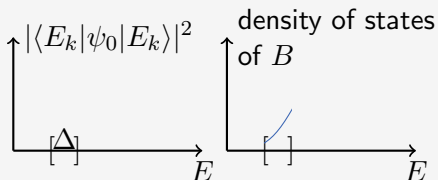


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The result

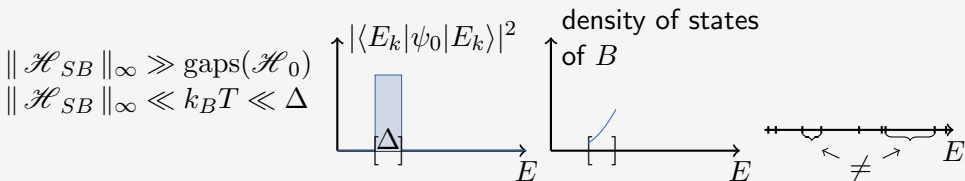
$$\begin{aligned} \|\mathcal{H}_{SB}\|_{\infty} &\gg \text{gaps}(\mathcal{H}_0) \\ \|\mathcal{H}_{SB}\|_{\infty} &\ll k_B T \ll \Delta \end{aligned}$$



\implies "Theorem" 2 (Theorem 2 in [9])

(Kinematic) Almost all pure states from a microcanonical subspace $[E, E + \Delta]$ are locally close to a Gibbs state.

The result



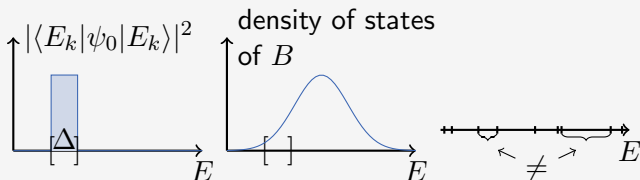
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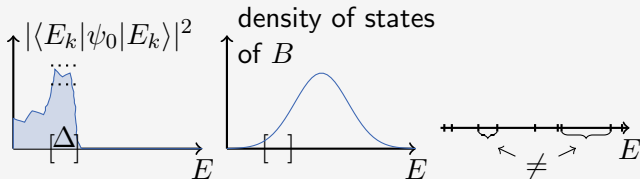
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Collaborators



Arnau Riera



Jens Eisert



Markus P. Müller

References

Thank you for your attention!

→ slides: www.cgogolin.de

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Journal of Statistical Mechanics: Theory and Experiment **2011** (2011) P02023.
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