

Under what conditions do quantum systems thermalize?

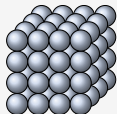
Christian Gogolin, Markus Müller, Arnau Riera, and Jens Eisert

Dahlem Center for Complex Quantum Systems, Freie Universität Berlin

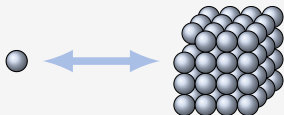
Institute for Physics and Astronomy, Universität Potsdam

Workshop “Information and Foundations of Thermodynamics”
at ETH Zürich 2011

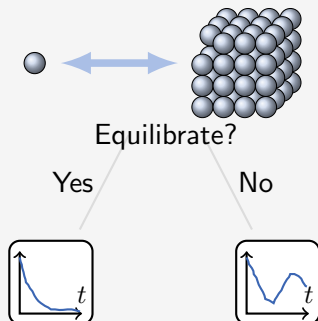
Understanding thermalization



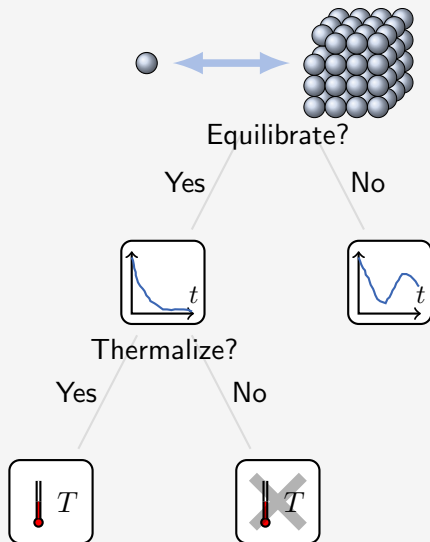
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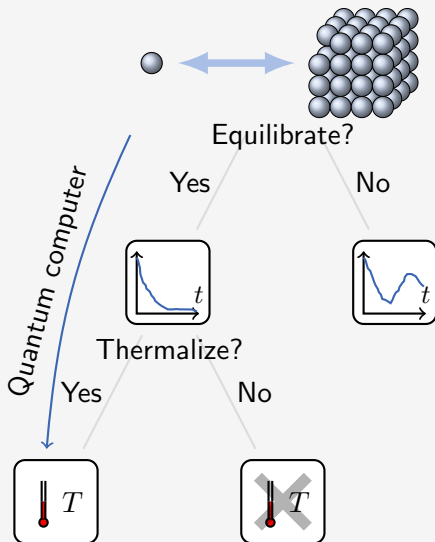
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Setup

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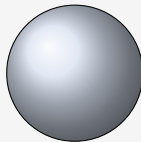
System, $\mathcal{H}_S, \mathcal{H}_S$

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



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

$$d_B \gg d_S$$



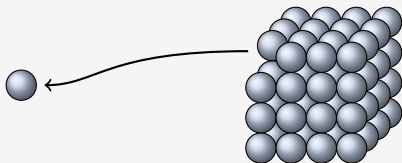
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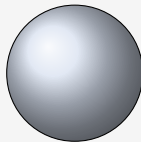
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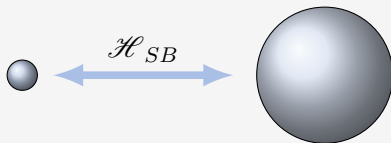
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$$\mathcal{H} = \mathcal{H}_S \otimes \mathbf{1} + \mathbf{1} \otimes \mathcal{H}_B + \mathcal{H}_{SB}$$

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

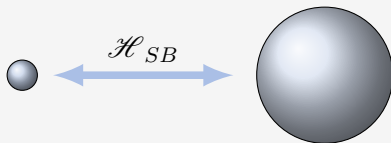
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$$\psi_t^S = \text{Tr}_B[\psi_t]$$

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Equilibration

Equilibration

Theorem 1 (Equilibration [2])

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

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

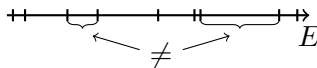
Equilibration

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

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

Theorem

If \mathcal{H} has
there exists

$\langle \psi_0 |$

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Effective dimension

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

Intuition: Dimension of supporting energy subspace

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

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

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Maximum entropy principle

Theorem 2 (Maximum entropy principle [3])

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

$$\overline{\text{Tr}[A \psi_t]} = \text{Tr}[A \overline{\psi_t}] = \text{Tr}[A \omega],$$

$$\text{where } \omega = \sum_k \pi_k \psi_0 \pi_k$$

(with π_k the energy eigen projectors) is the *dephased* state that *maximizes the von Neumann entropy, given all conserved quantities.*

[3] C. Gogolin, M. P. Müller, and J. Eisert, PRL 106, 040401 (2011)

[4] M. Rigol, V. Dunjko, V. Yurovsky, and M. Olshanii, PRL 98, 050405 (2007)

Maximum Time averaging

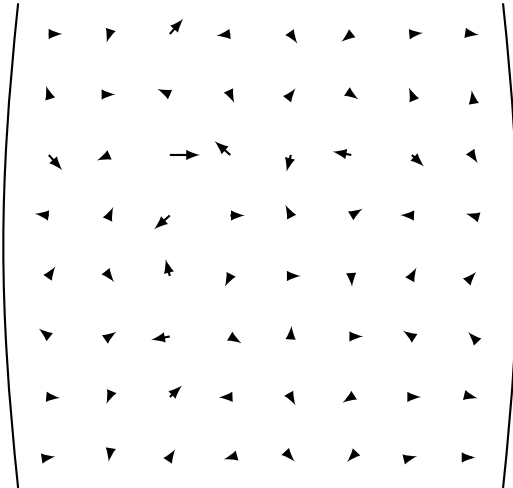
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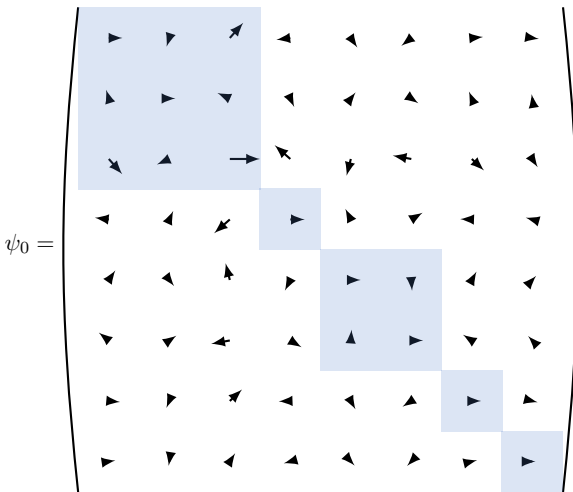
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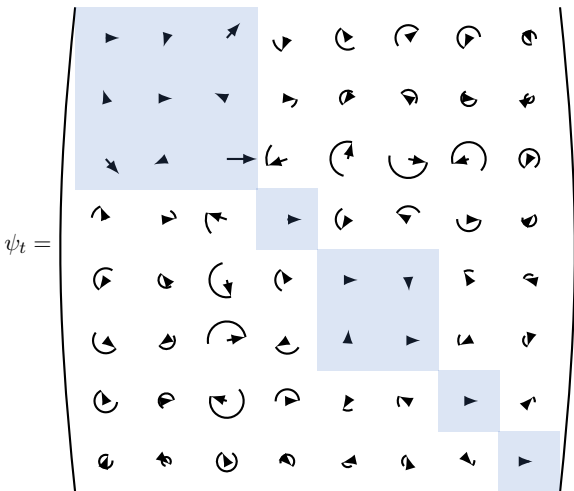
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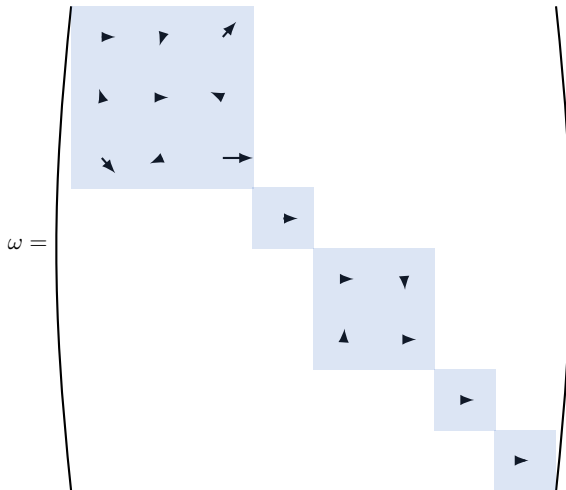
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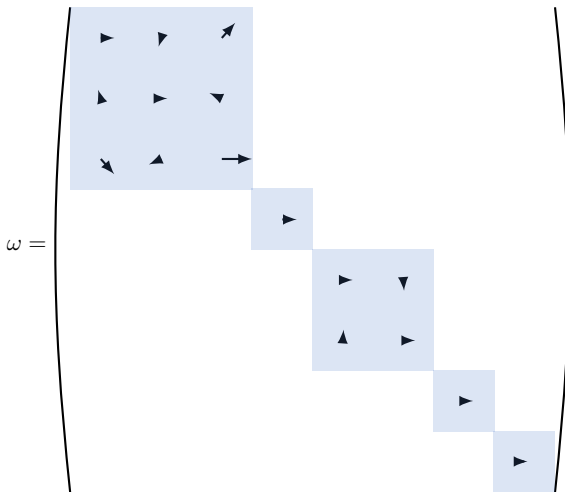
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Maximum Entropy

Time averaging



$\psi_0 \rightarrow \omega$ is a **pinching** $\Rightarrow \omega$ **maximizes entropy**.

[3] C. Gogolin, M. P. Müller, and J. Eisert, PRL 106, 040401 (2011)

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⇒ Maximum entropy principle from pure quantum dynamics.

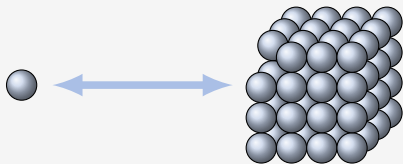
Proves a *conjecture from the condensed matter literature* from 2007 [4].

[3] C. Gogolin, M. P. Müller, and J. Eisert, PRL 106, 040401 (2011)

[4] M. Rigol, V. Dunjko, V. Yurovsky, and M. Olshanii, PRL 98, 050405 (2007)

Thermalization and realistic weak coupling

Thermalization is a complicated process



Thermalization implies:

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

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

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

[3] C. Gogolin, M. P. Müller, and J. Eisert, PRL 106, 040401 (2011)

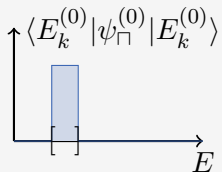
[5] A. Riera, C. Gogolin, and J. Eisert, 1102.2389

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

[7] J. Gemmer, M. Michel, and G. Mahler, Springer (2009)

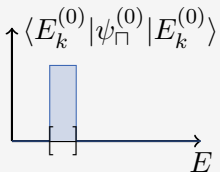
Level counting with no coupling

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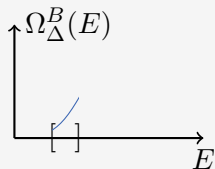
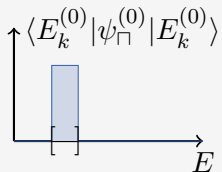
Well known fact [8]:

$$\omega_{\square}^{S(0)} \stackrel{\text{no coupling}}{\propto} \sum_k \Omega_{\Delta}^B(E - E_k^S) \underbrace{\# \text{ bath states in } [E - E_k^S, E - E_k^S + \Delta]}_{\leftarrow} |E_k^S\rangle \langle E_k^S|$$

[8] S. Goldstein, PRL 96, 050403 (2006)

Level counting with no coupling

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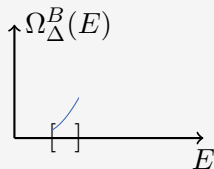
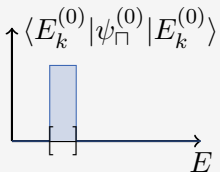


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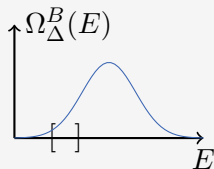
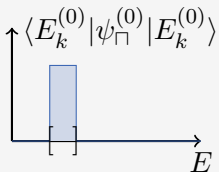


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perturbative coupling

\Downarrow [3]

effective entanglement in the eigenbasis $R(\psi_0)$ is small

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absence of initial state independence.

$$\mathcal{D}(\omega^{S(1)}, \omega^{S(2)}) \geq \mathcal{D}(\psi_0^{S(1)}, \psi_0^{S(2)}) - R(\psi_0^{S(1)}) - R(\psi_0^{S(2)})$$

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Effective entanglement in the eigenbasis

$$R(\psi_0) = \sum_k |\langle E_k | \psi_0 \rangle|^2 \mathcal{D}(\text{Tr}_B |E_k\rangle \langle E_k|, \psi_0^S)$$

Measures **how entangled the eigenbasis** feels for the given initial state.

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Perturbative coupling $\| \mathcal{H}_{SB} \|_{\infty} < \text{gaps}(\mathcal{H}_0) \dots$

- ... is *locally dense*.
- ... pro

Theorem 3 (Entanglement in eigenbasis)

For every orthonormal basis $\{|i\rangle\}$ for S and every initial product state with $\psi_0 = |j\rangle\langle j| \otimes \phi_0^B$, the effective entanglement in the eigenbasis (for non-degenerate \mathcal{H}) is on average upper bounded by

$$\mathbb{E}_{\phi_0^B} R(|j\rangle\langle j| \otimes \phi_0^B) \leq 2 \delta d_S,$$

where

$$\delta = \max_k \min_i \mathcal{D}(\text{Tr}_B |E_k\rangle\langle E_k|, |i\rangle\langle i|)$$

is the *geometric measure of entanglement* of the eigenstate $|E_k\rangle$ with respect to the basis $\{|i\rangle\}$.

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\Rightarrow Refutes wide spread believe that “non-integrable models thermalize.”

[3] C. Gogolin, M. P. Müller, and J. Eisert, PRL 106, 040401 (2011)

Realistic weak coupling

- Naive perturbation theory fails.
- More realistic weak coupling:

$$\text{gaps}(\mathcal{H}_0) \ll \| \mathcal{H}_{SB} \|_{\infty} \ll \Delta$$

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Theorem 4 (Corollary of a theorem from [5])

If $\|\mathcal{H}_{SB}\|_\infty \ll \Delta$ the dephased states $\omega_\square^{S(0)}$ and ω_\square^S are *close to each other* in the sense that

$$\mathcal{D}(\omega_\square^S, \omega_\square^{S(0)}) \lesssim 3\sqrt{\frac{\|\mathcal{H}_{SB}\|_\infty}{2\Delta}}.$$

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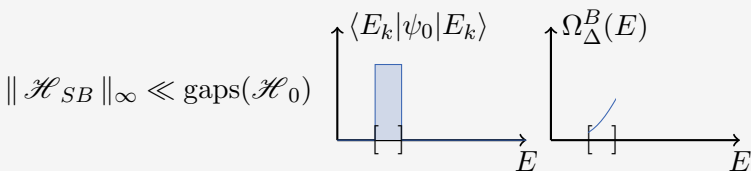
$$\mathcal{D}(\omega_\square^S, \omega_\square^{S(0)}) \lesssim \frac{\sqrt{\beta \|\mathcal{H}_{SB}\|_\infty}}{1 - e^{-\beta\Delta}}.$$

Putting everything together

Consequences

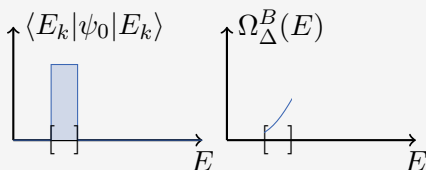
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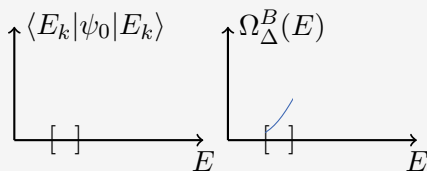
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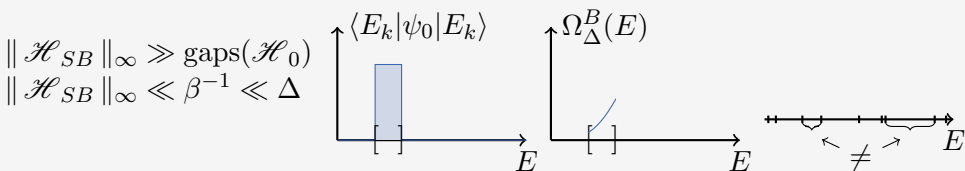
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The left diagram shows energy levels E_k on the vertical axis and energy E on the horizontal axis. A gap is indicated between two levels, and a state $|\psi_0\rangle$ is shown between them. The right diagram shows the density of states $\Omega_{\Delta}^B(E)$ on the vertical axis and energy E on the horizontal axis. A shaded region indicates a microcanonical subspace $[E, E + \Delta]$.

\implies “Theorem” 5 (Theorem 2 in [5])

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

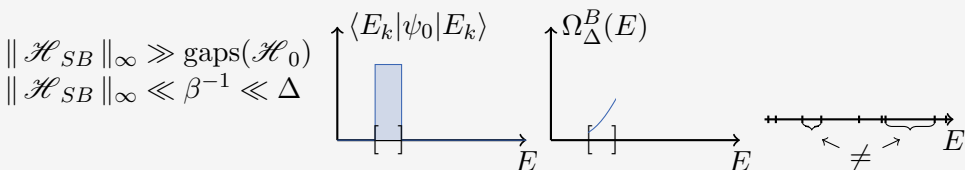
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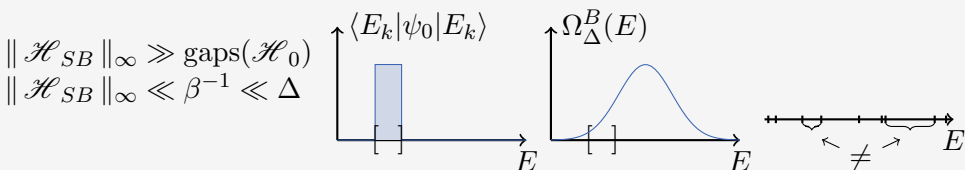


\Rightarrow “Theorem” 5 (Theorem 2 in [5])

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

(Dynamic) All initial states $\psi_{\square,0}$ locally equilibrate towards a Gibbs state, even if they are initially far from equilibrium.

Consequences

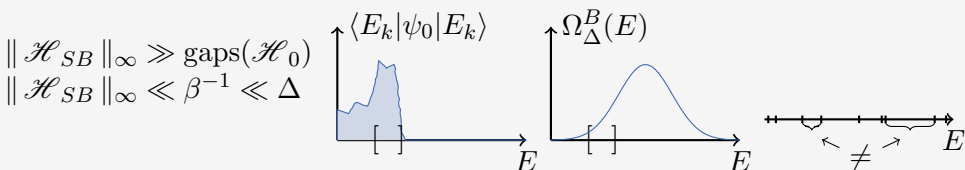


\Rightarrow “Theorem” 5 (Theorem 2 in [5])

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

(Dynamic) All initial states $\psi_{\square,0}$ locally equilibrate towards a Gibbs state, even if they are initially far from equilibrium.

Consequences



\Rightarrow “Theorem” 5 (Theorem 2 in [5])

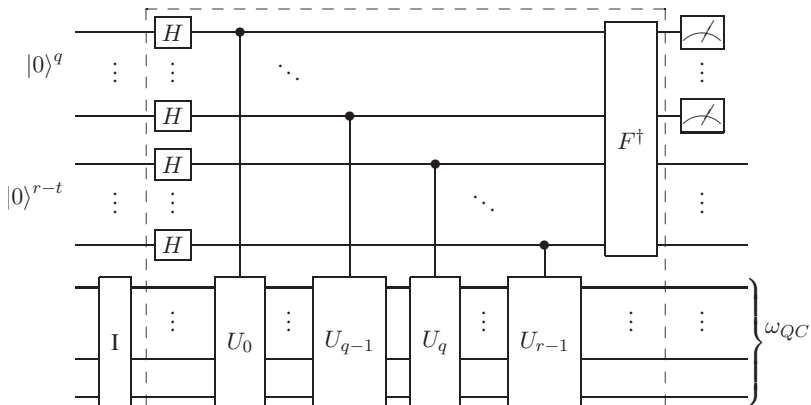
(Kinematic) Almost all pure states from a microcanonical subspace $[E, E + \Delta]$ are locally close to a Gibbs state.

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A quantum algorithm for Gibbs state preparation

A quantum algorithm for Gibbs state preparation

Quantum circuit



[5] A. Riera, C. Gogolin, and J. Eisert, 1102.2389

[9] D. Poulin and P. Wocjan, PRL 103, 220502 (2009)

A quantum algorithm for Gibbs state preparation

Quantum circuit

- No detailed knowledge about \mathcal{H}_S is required.

$|0\rangle^q$

$|0\rangle^{r-t}$

ω_{QC}

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A quantum algorithm for Gibbs state preparation

Quantum circuit

- No detailed knowledge about \mathcal{H}_S is required.
 - The algorithm uses partial phase estimation to prepare ω_Π .
 - Complements quantum Metropolis
 - Trace distance error bound
 - Explicit runtime
- (also compare [9])

ω_{QC}

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A quantum algorithm for Gibbs state preparation

Quantum circuit

 $|0\rangle^q$
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- The algorithm uses partial phase estimation to prepare ω_Π .
- Complements quantum Metropolis
 - Trace distance error bound
 - Explicit runtime
 (also compare [9])
- Complexity for fixed trace distance error:
 - polynomially many ancilla qubits
 - exponential runtime
 ($\Omega(e^{\beta(E_2^S - E_1^S)})$ is necessary)

 ω_{QC}

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Collaborators



Arnau Riera



Martin Kliesch



Jens Eisert



Markus P. Müller



Andreas Winter

References

Thank you for your attention!

→ slides: www.cgogolin.de

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