



Correlation-Enhanced Heat-Bath Algorithmic Cooling

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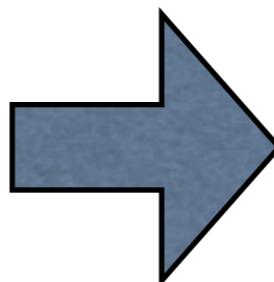
Outline



- 1) Introduction to algorithmic cooling
- 2) Improvement due to correlated relaxation processes between qubit-environment
- 3) Improvement due to correlations present in the initial state

Introduction

Cooling physical systems



Processing of information → **Entropy manipulation**

Tools of quantum information processing give us new techniques for cooling, in non conventional ways.

Algorithmic Cooling

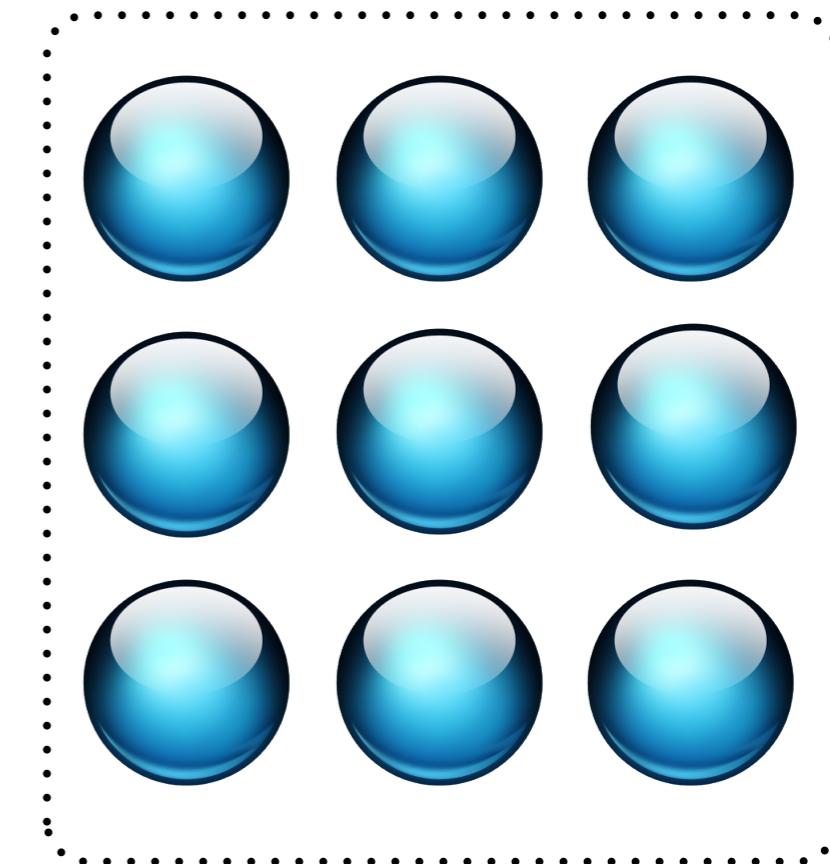


How can we cool some qubits using unitary operations and contact with a bath?

Target qubit



Qubit system

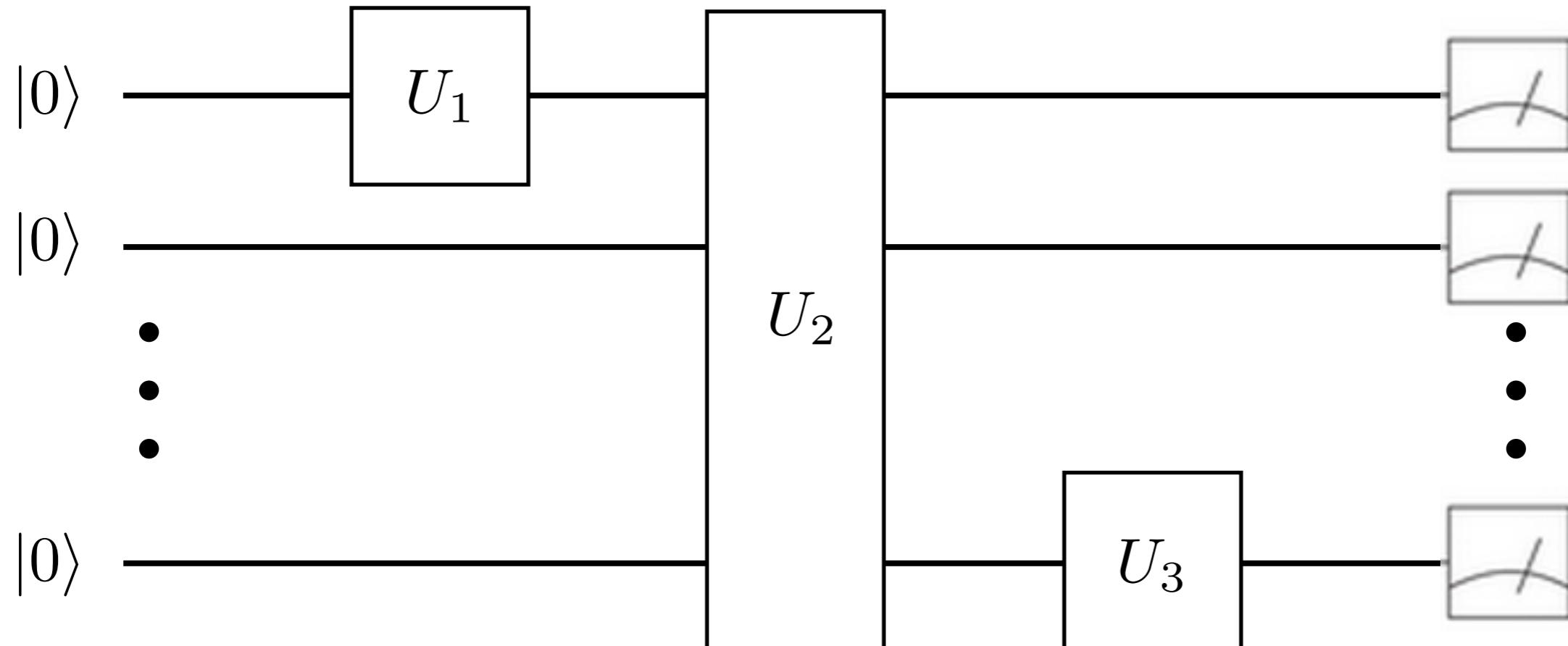


Heat-bath

(and projective measurements not allowed. *)

Motivation

Preparation of highly pure quantum states

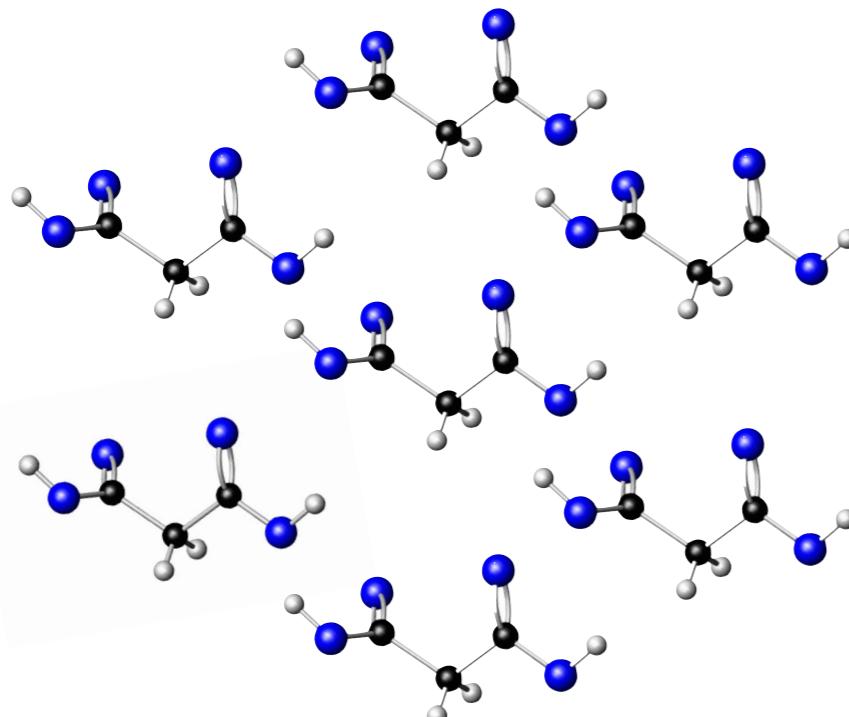


A quantum computer needs a constant supply of highly pure qubits:

- As initial state (ex. $|0000\rangle$) for quantum algorithms
- As ancillas for quantum error correction

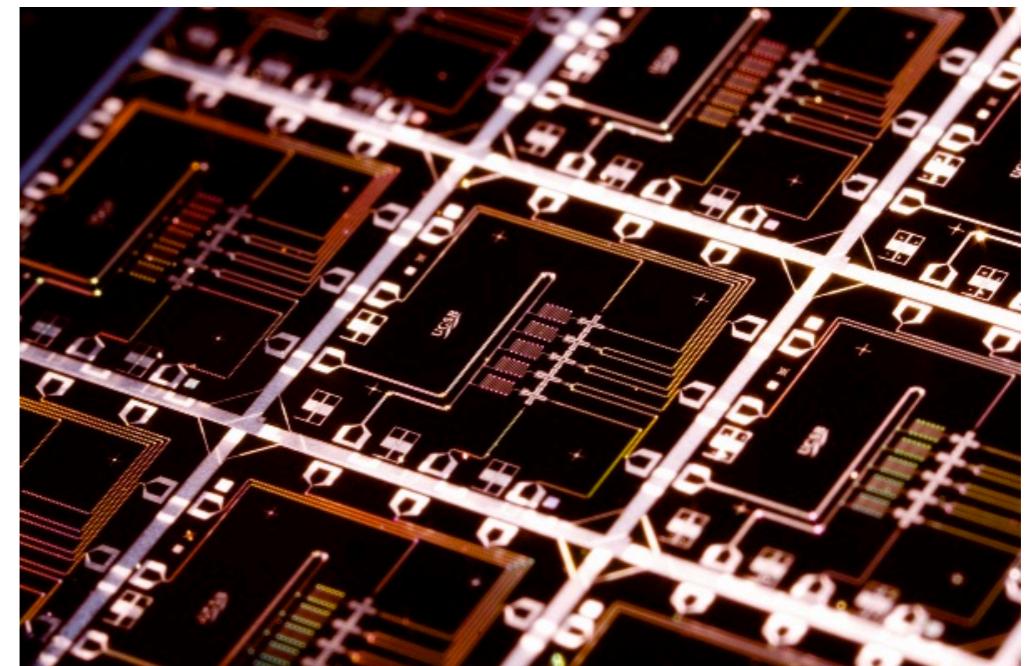
State preparation without projective measurements is a big challenge in Quantum Computing

Ensemble implementations



Ex. NMR, ESR
Highly mixed states

In the lab

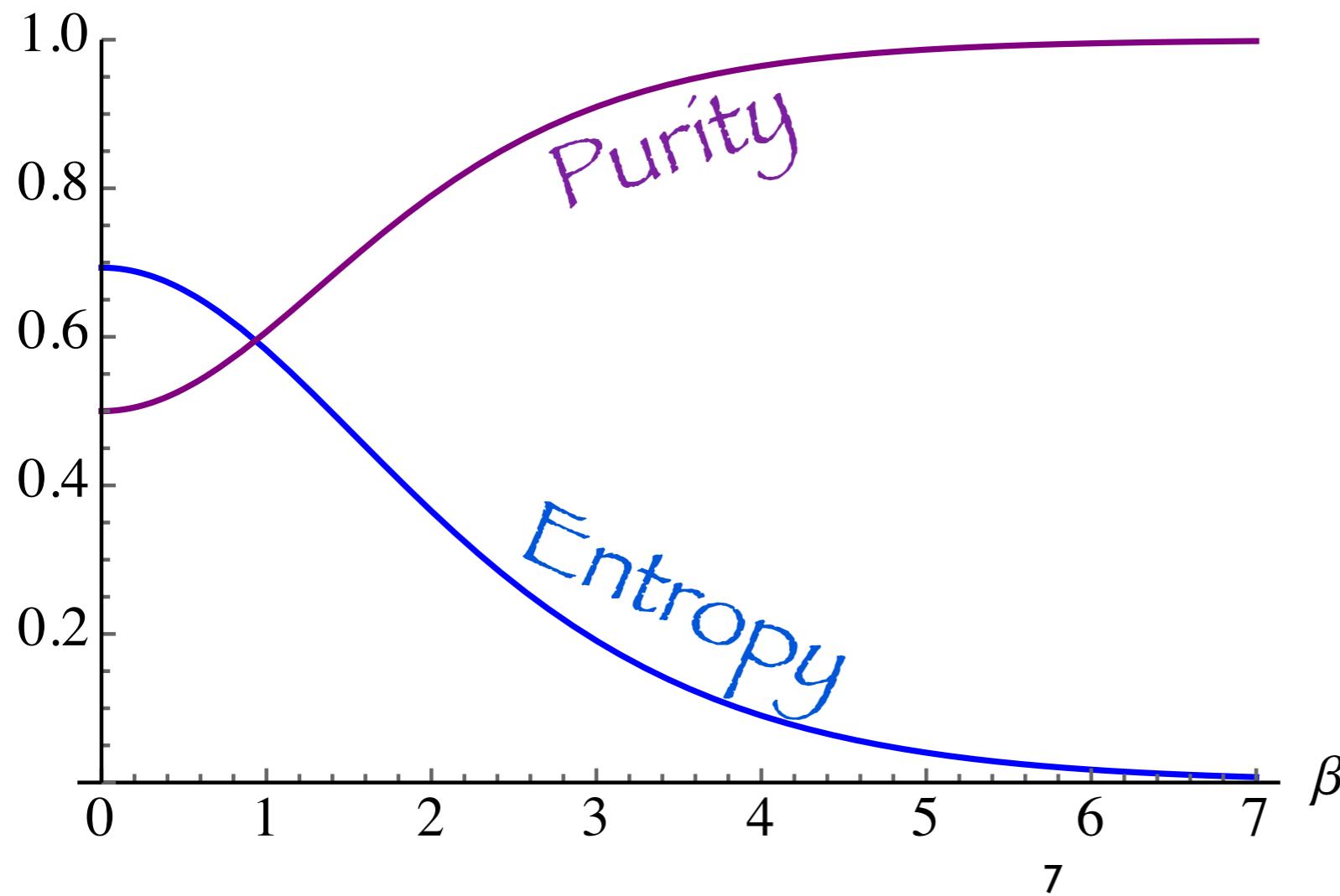


Non-perfect projective measurements

For the target qubit:



$$\rho = \begin{pmatrix} \rho_{00} & \rho_{01} \\ \rho_{10} & \rho_{11} \end{pmatrix}$$



Entropy $S(\rho) = -\text{Tr}(\rho \log \rho)$
Purity $\mathcal{P} = \text{Tr}(\rho^2)$

Local thermal state

$$\rho = \frac{e^{-\beta H}}{Z}, \quad H = \frac{1}{2}\omega\sigma_z$$

$$\mathcal{P} = \frac{1}{2}(1 + \tanh^2[\omega\beta/2])$$

$$T = \beta^{-1}$$

Chapter one

Introduction to algorithmic cooling

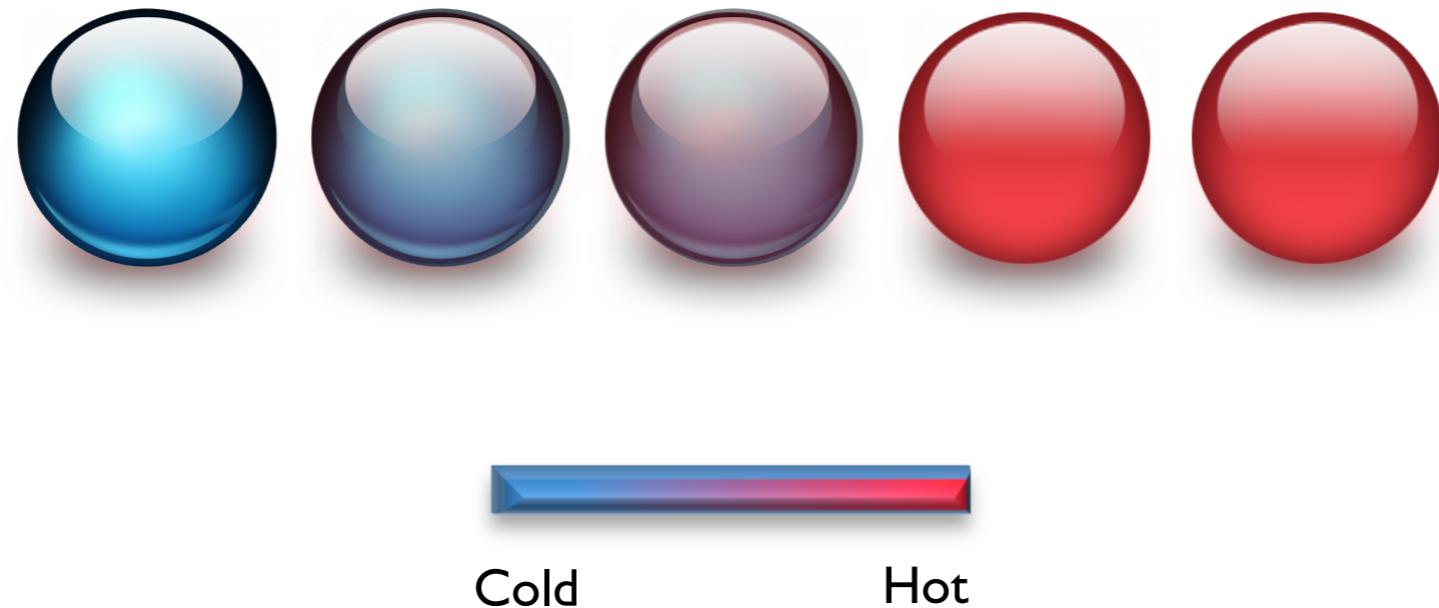
Algorithmic Cooling (AC)



“A quantum mechanical heat engine”

* Schulman and Vazirani, 1990

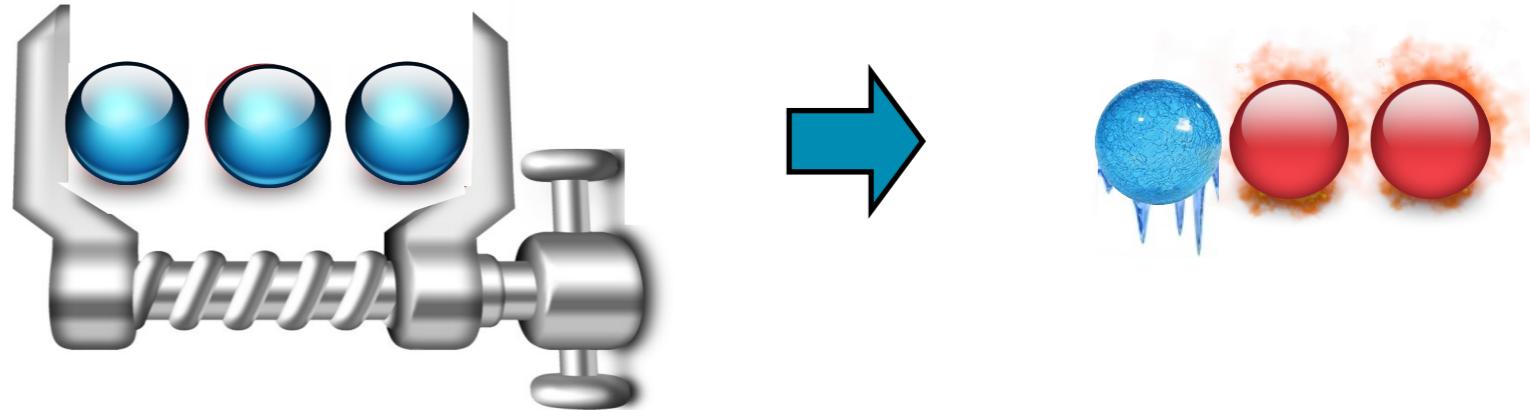
Algorithmic Cooling (AC)



The idea is to re-distribute the entropy within a group of qubits using a unitary operation U .

$$\rho_{1\text{st}} = \text{Tr}_{\overline{1\text{st}}} (U \rho_{\text{total}} U^\dagger) \quad S(\rho_{1\text{ST}}) \rightarrow 0$$

Simple example



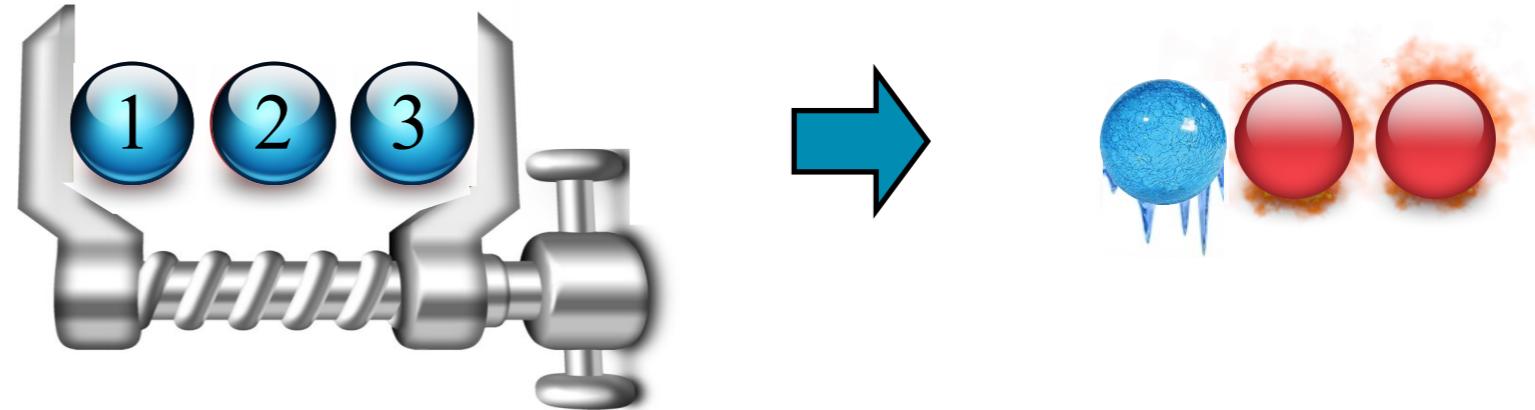
$$\rho_{\epsilon_0} = \frac{1}{2} \begin{pmatrix} 1 + \epsilon_0 & 0 \\ 0 & 1 - \epsilon_0 \end{pmatrix}$$

$$\text{diag}(\rho_{\epsilon}^{\otimes 3}) = \frac{1}{2^3} \begin{bmatrix} (1 + \epsilon_0)^3 \\ (1 + \epsilon_0)^2 (1 - \epsilon_0) \\ (1 + \epsilon_0)^2 (1 - \epsilon_0) \\ (1 + \epsilon_0) (1 - \epsilon_0)^2 \\ (1 + \epsilon_0)^2 (1 - \epsilon_0) \\ (1 + \epsilon_0) (1 - \epsilon_0)^2 \\ (1 + \epsilon_0) (1 - \epsilon_0)^2 \\ (1 - \epsilon_0)^3 \end{bmatrix}$$

000
001
010
011
100
101
110
111

A pink circle highlights the first two entries of the vector, corresponding to the first two rows of the matrix. A black arrow points from the bottom row to the rightmost column.

Simple example

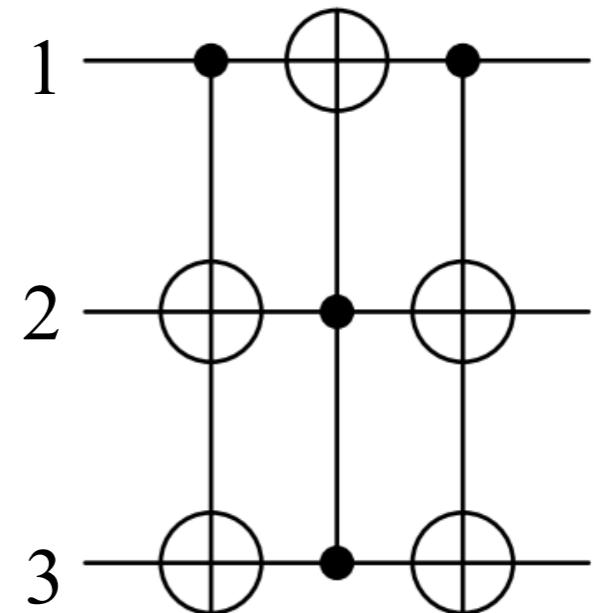


$$U = |011\rangle\langle100| + |100\rangle\langle011|$$

Example, when $\beta = \frac{1}{k_B T_0} \ll 1$:

Target qubit
effective temperature

$$T_0 \rightarrow \frac{2}{3}T_0$$



Second and third qubit effective temperatures

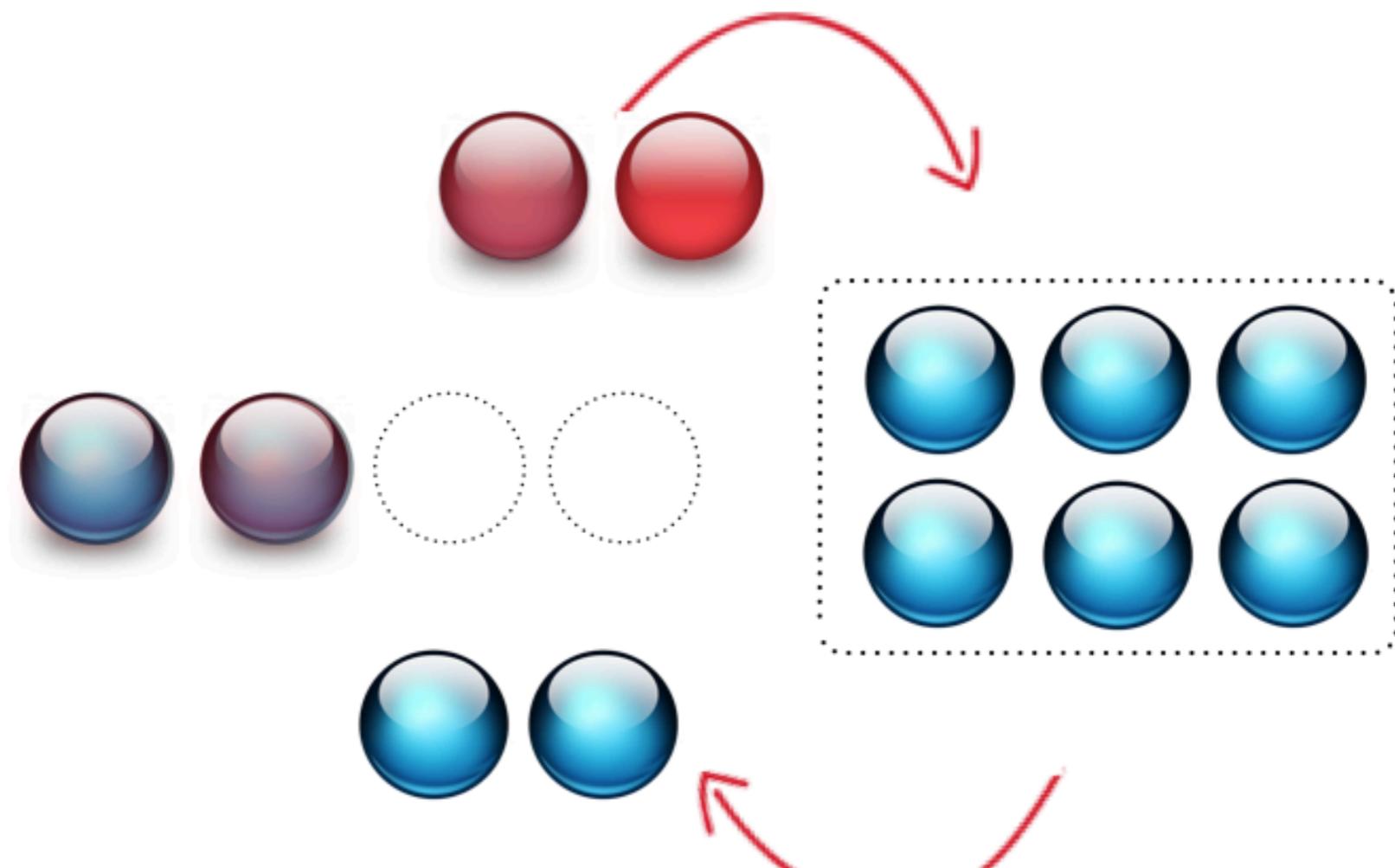
$$T_0 \rightarrow 2T_0$$

Limits of entropy compression using unitary dynamics for a close system

- Shannon Entropy bound
(Limited by unitary dynamics)
- Conservation of eigenvalues of the total system
- Still uses a lot of resources

Ex. $\epsilon_b \sim 10^{-5} \rightarrow 10^{12}$ qubits

Heat-Bath Algorithm Cooling



Refreshing step

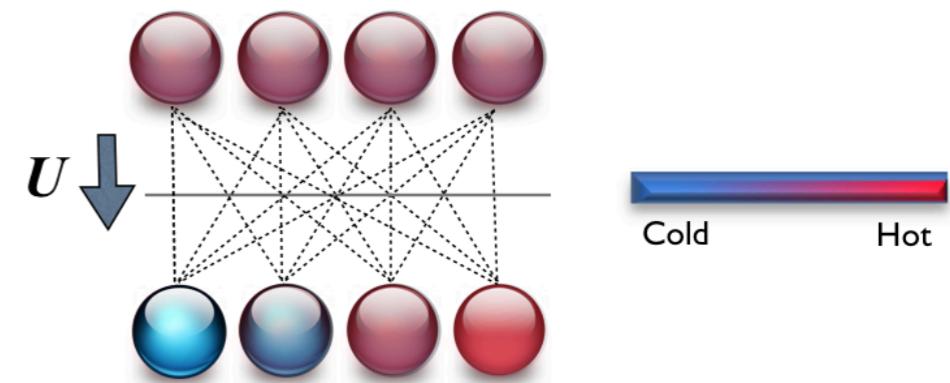
$$\rho' \rightarrow \rho'' = \text{Tr}_m (\rho) \otimes \rho_{\epsilon_b}^{\otimes m}$$

The Partner Pairing Algorithm (PPA-HBAC)

PPA-HBAC is the iteration of the two steps:

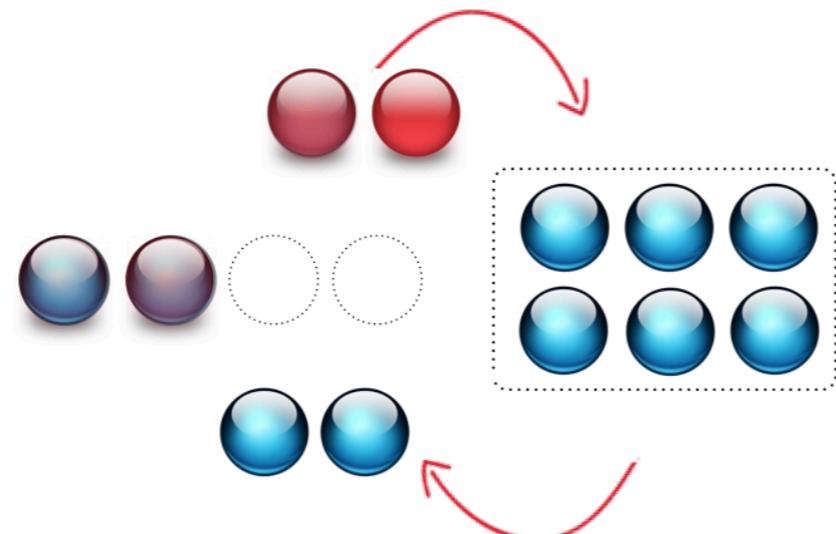
① **Entropy Compression Step.**

$$\rho \rightarrow \rho' = U\rho U^\dagger$$



$U \equiv$ Descending SORT of the diagonal elements.

② **Refresh Step.**

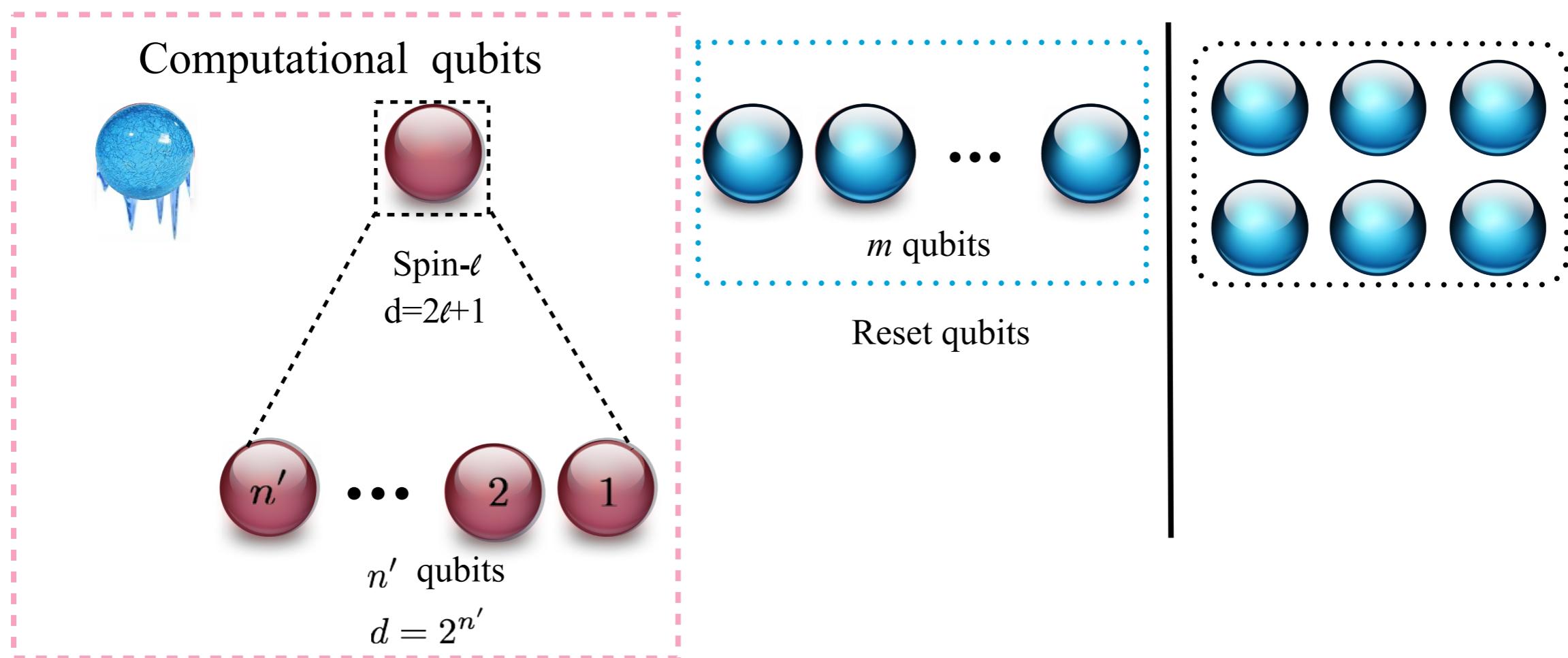


$$\rho' \rightarrow \rho'' = \text{Tr}_m(\rho) \otimes \rho_{\epsilon_b}^{\otimes m}$$

Cooling Limit

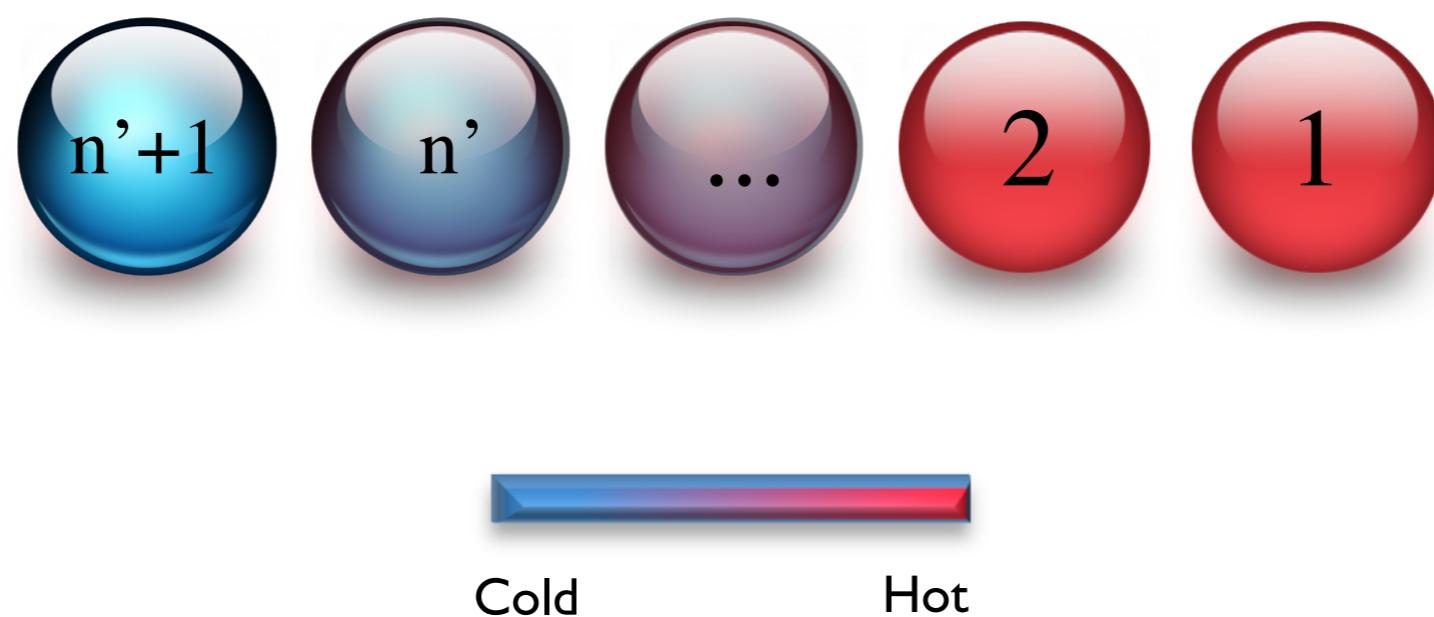
$$\Phi(\rho) := \text{Tr}_m(U(\rho)\rho U^\dagger(\rho)) \otimes \rho_b^{\otimes m}$$

In the limit: $\rho = \Phi(\rho)$



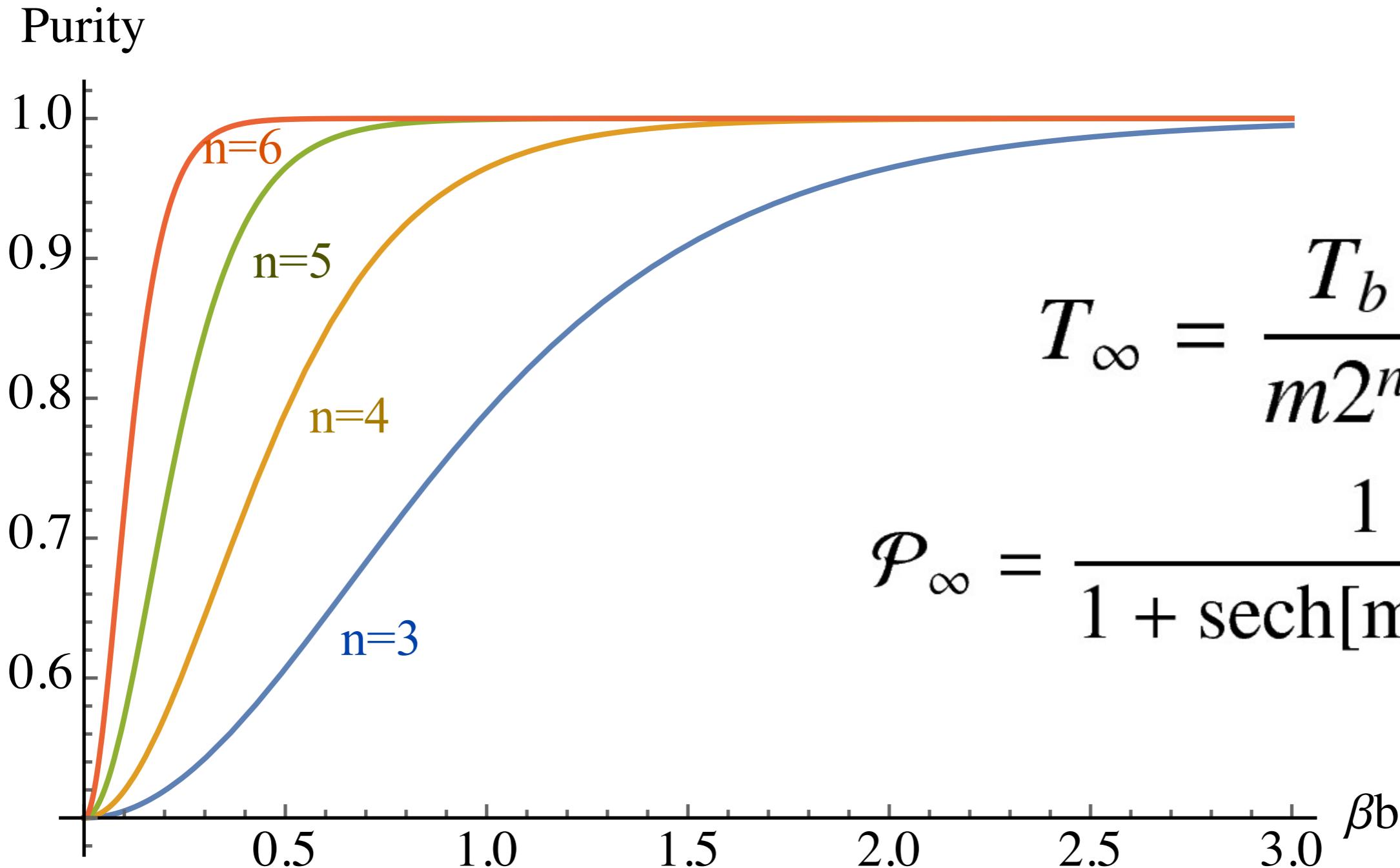
Steady State

$$\rho_T^\infty = \bigotimes_{j=1}^{n'+1} \rho_{j^{th}}^\infty = \bigotimes_{j=1}^{n'+1} \frac{1}{2} \begin{pmatrix} 1 + \epsilon_j^\infty & 0 \\ 0 & 1 - \epsilon_j^\infty \end{pmatrix}$$

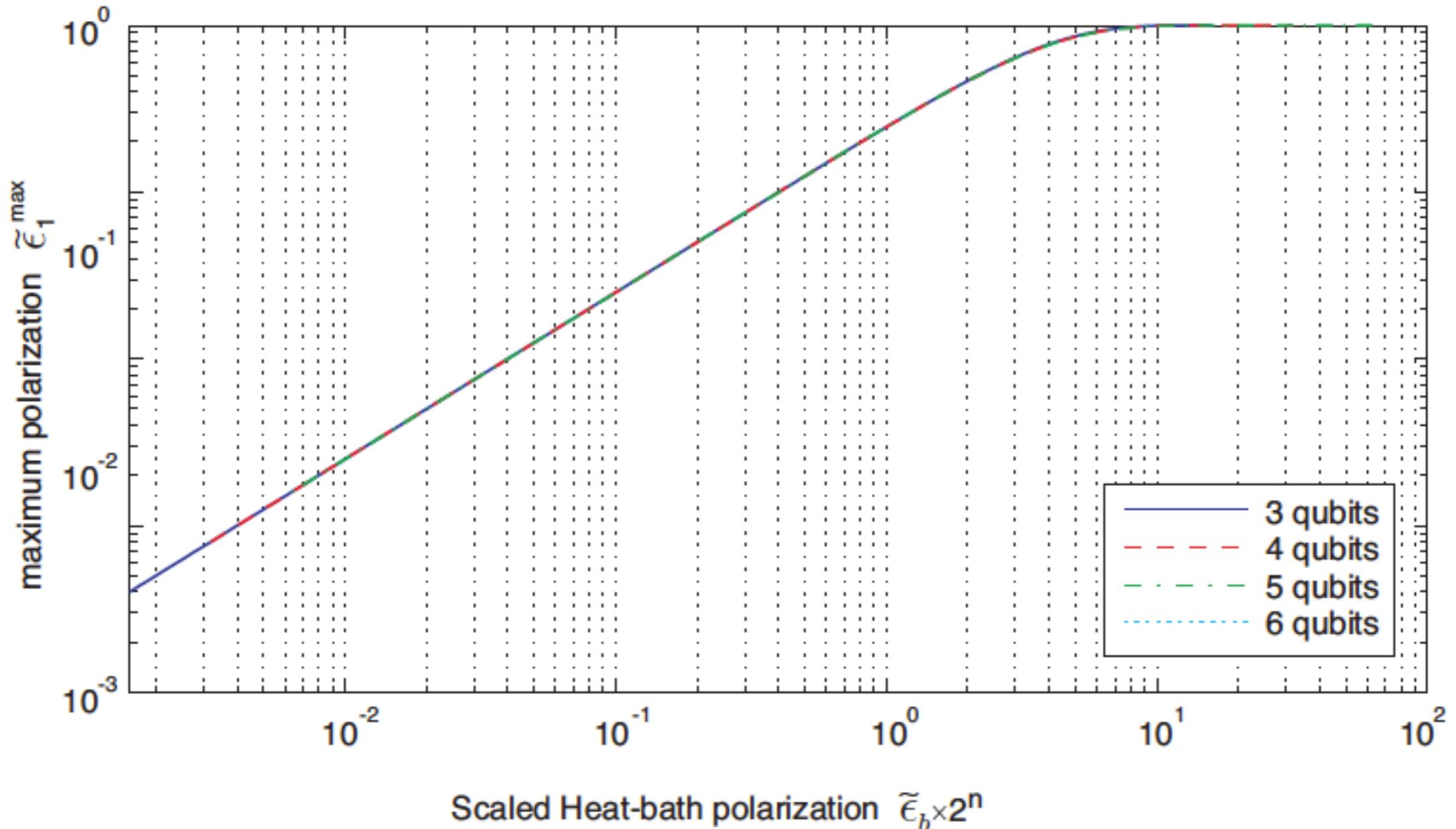


$$\epsilon_j^\infty = \frac{(1 + \epsilon_b)^{m2^{j-1}} - (1 - \epsilon_b)^{m2^{j-1}}}{(1 + \epsilon_b)^{m2^{j-1}} + (1 - \epsilon_b)^{m2^{j-1}}}$$

Cooling limits of PPA-HBAC



Threshold effect



$\epsilon \ll \frac{1}{2^n}$ Can cool to only $2^{n-2}\epsilon$

$\epsilon \gg \frac{1}{2^n}$ Can purify to order unity.

Is this the fundamental cooling
limit?

Some people claim it is....

S. Raesi and M. Mosca, PRL 114, 100404 (2016)

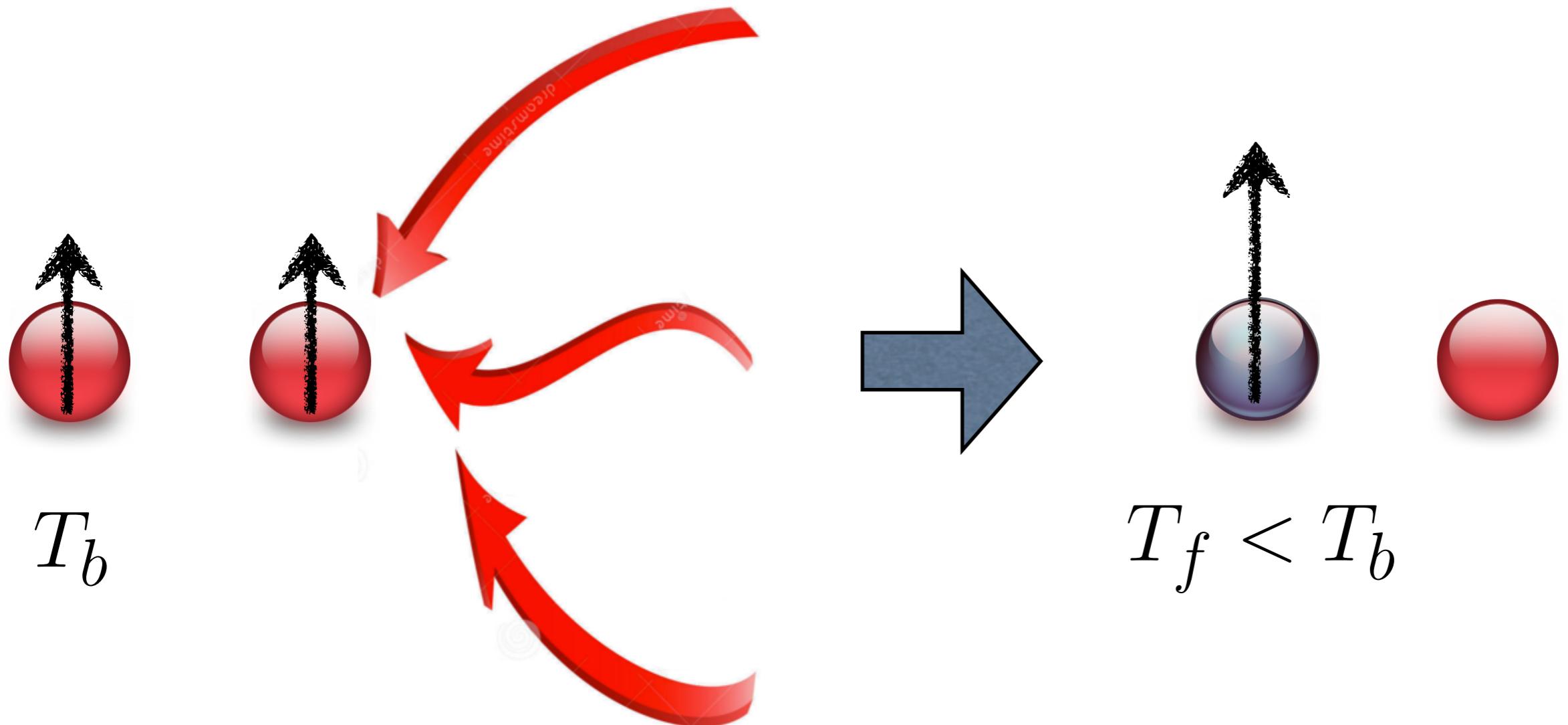
L. Schulman, T. Mor, and Y. Weinstein, PRL 94, 120501 (2005)

However, we found new techniques to circumvent this cooling limit by taking advantage of correlations

Chapter two

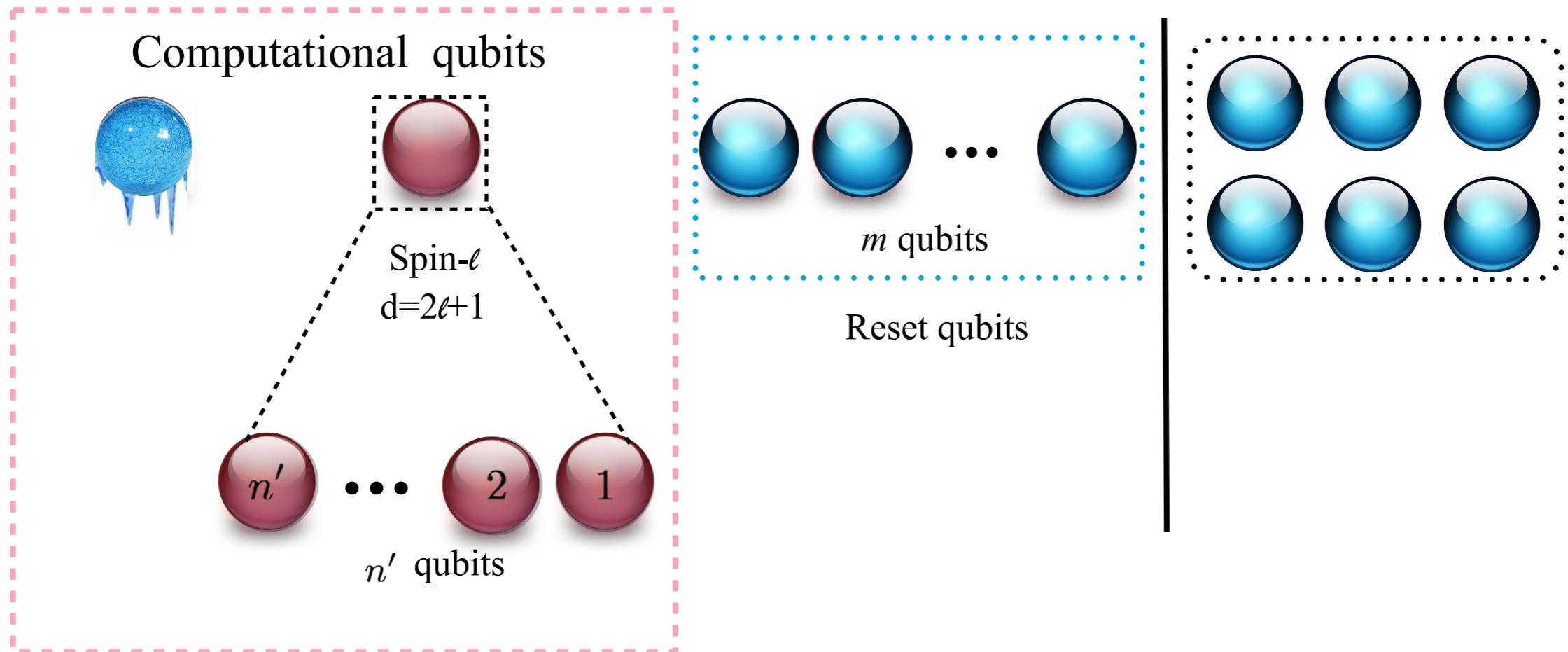
Correlated relaxation

Nuclear Overhauser Effect (NOE), 1953

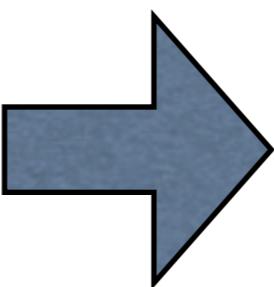


A. Overhauser, Phys. Rev. 89, 689 (1953).

Cooling limits of PPA-HBAC



$$T_\infty = \frac{T_b}{m2^{n'}}$$



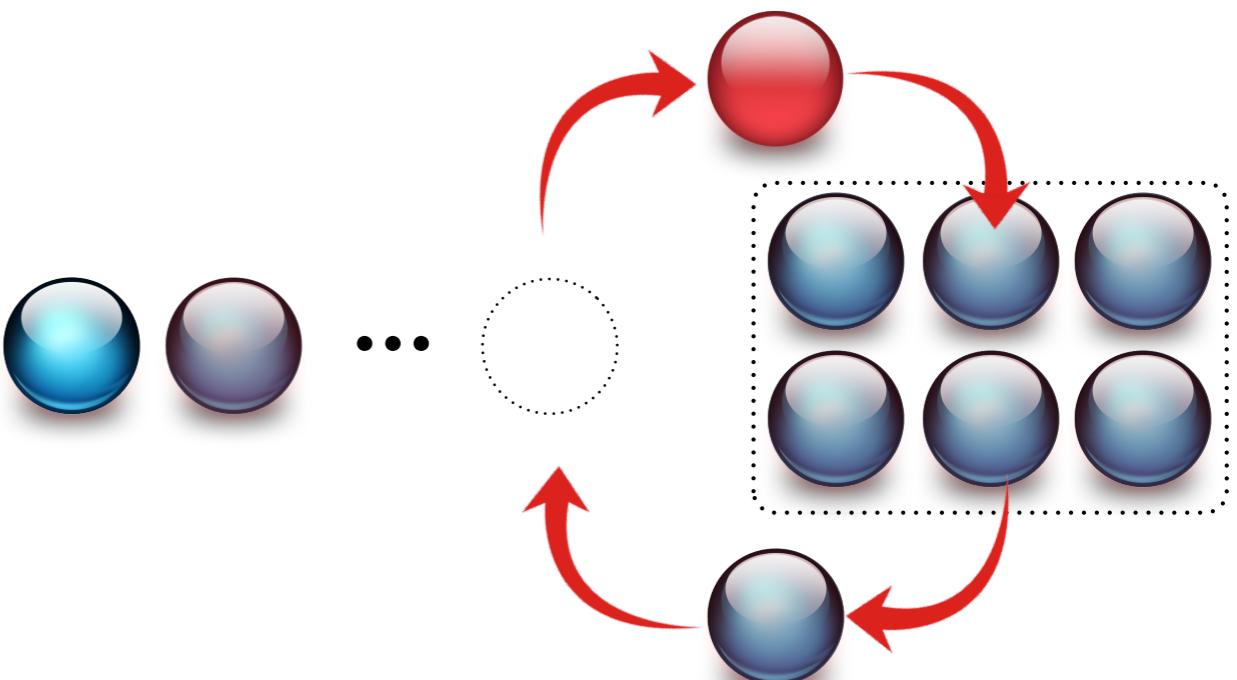
For 2-qubit system:
PPA-HBAC: $T_\infty = T_b$

Thermalization of qubits

PPA-HBAC

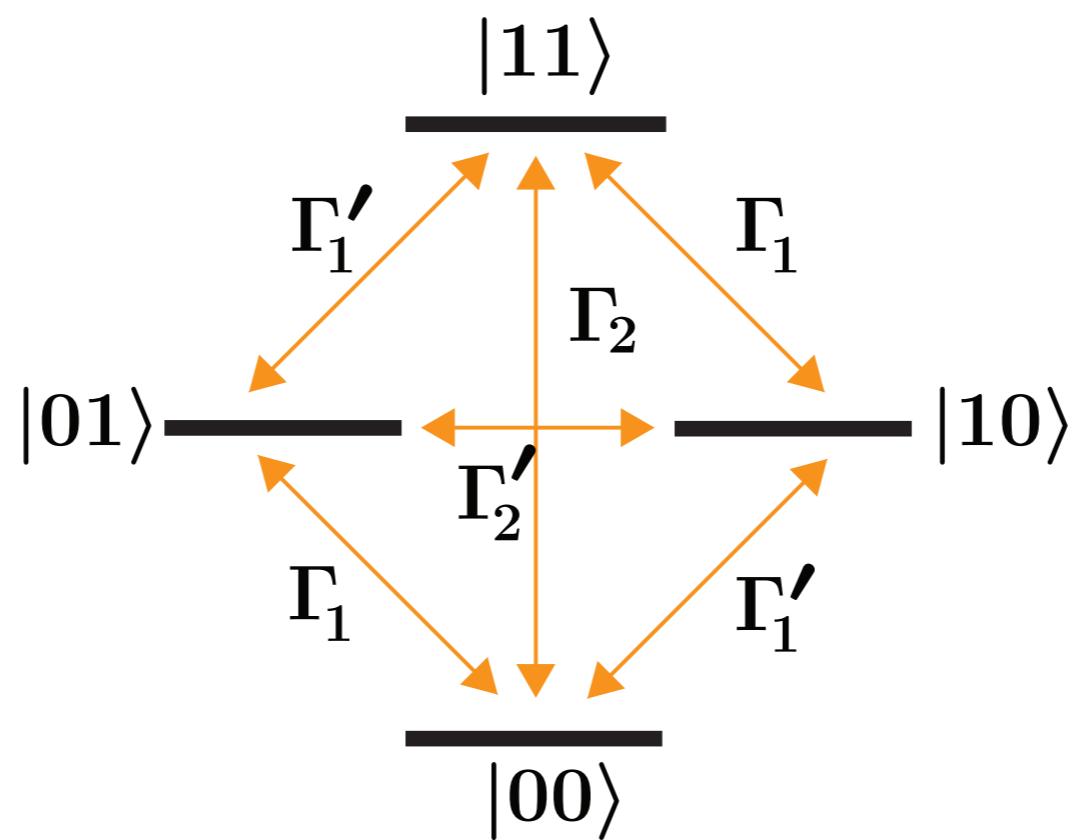
$$\rho_{total} \rightarrow \text{Tr}_m(\rho_{total}) \otimes \rho_b^{\otimes m}$$

$m \equiv$ number of refreshed qubits

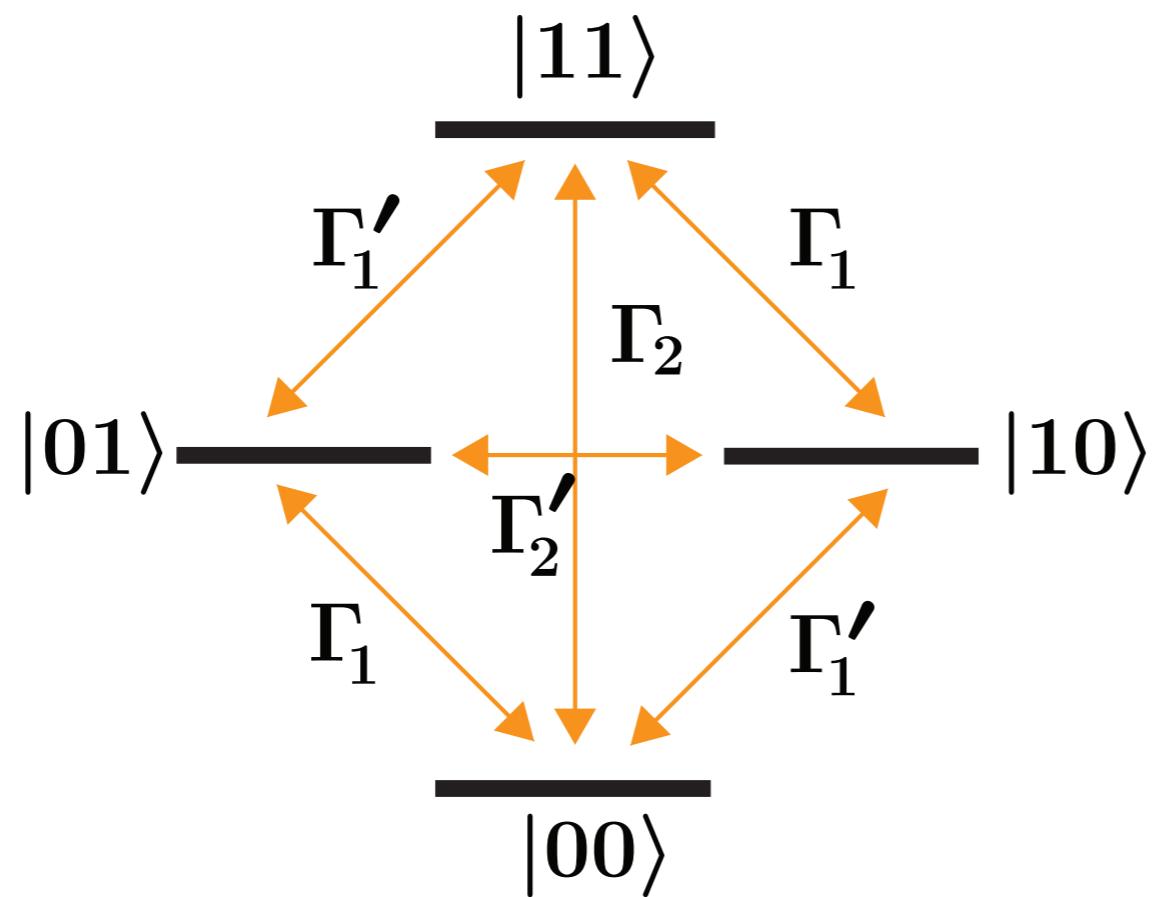


NOE

Crossed-relaxation



Crossed-relaxation



$$\begin{aligned} \Gamma_2 &\neq 0 \\ \& \quad \& \\ \Gamma_1 = \Gamma'_1 = \Gamma'_2 &= 0 \end{aligned}$$

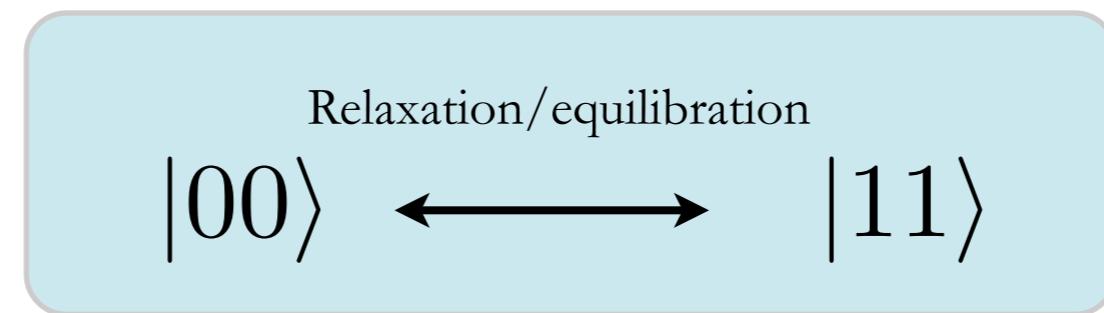
$$\text{diag}(\rho) = \begin{bmatrix} A_{00} \\ A_{01} \\ A_{10} \\ A_{11} \end{bmatrix} \xrightarrow{\Gamma_2} \begin{bmatrix} (A_{00} + A_{11}) P_2 \\ A_{01} \\ A_{10} \\ (A_{00} + A_{11})(1 - P_2) \end{bmatrix}$$

$$\begin{aligned} P_2 &= \frac{e^{2\xi}}{2 \cosh(2\xi)} \\ \xi &= \tanh^{-1}(\epsilon_b) \end{aligned}$$

Wrong implicit assumption in all previous work:

- The reset step with the bath is optimal when swapping qubits to pump entropy out of the system.

Reset of states Γ_2



Kraus Operators

$$A_1^{(n=2)} = \sqrt{p_2}|00\rangle\langle 00|$$

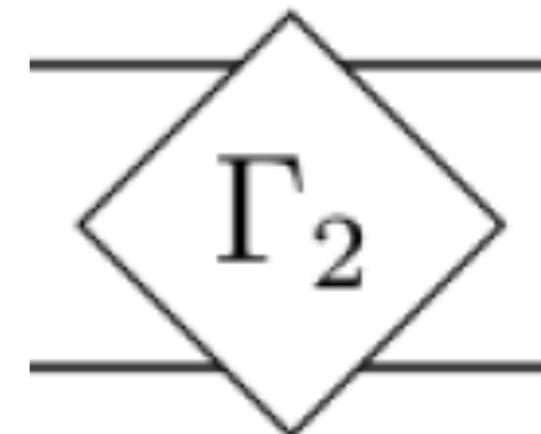
$$A_2^{(n=2)} = \sqrt{p_2}|00\rangle\langle 11|$$

$$A_3^{(n=2)} = \sqrt{1-p_2}|11\rangle\langle 11|$$

$$A_4^{(n=2)} = \sqrt{1-p_2}|11\rangle\langle 00|$$

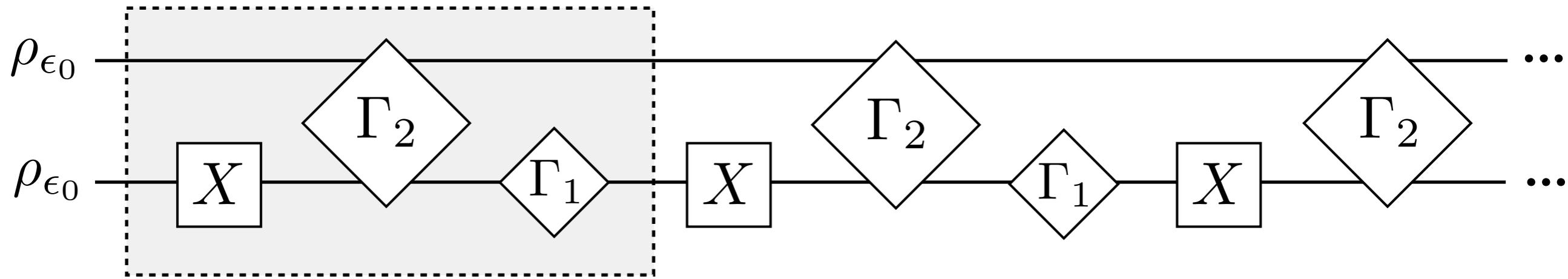
$$A_5^{(n=2)} = |01\rangle\langle 01|$$

$$A_6^{(n=2)} = |10\rangle\langle 10|$$



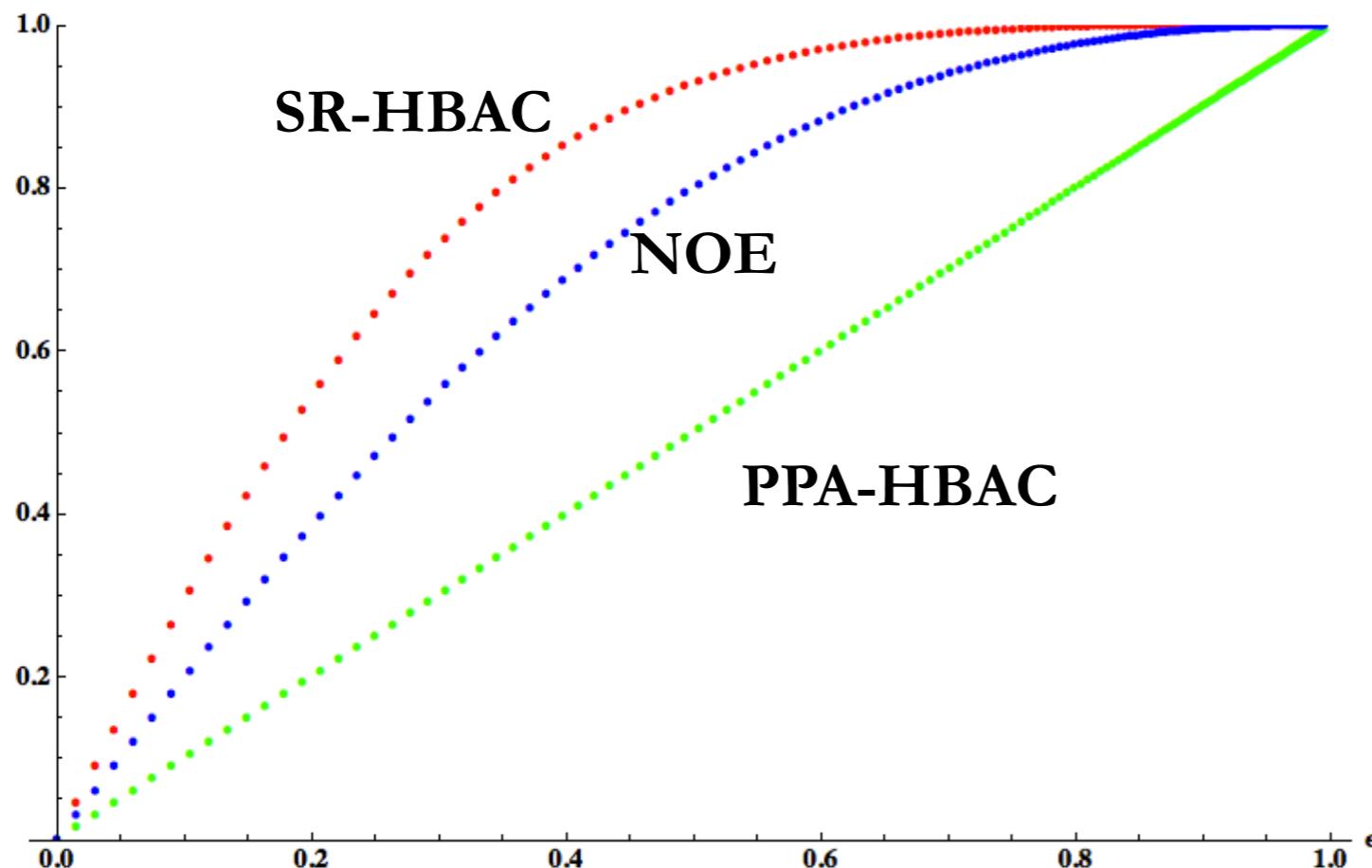
$$P_2 = \frac{e^{2\xi}}{2 \cosh(2\xi)}$$
$$\xi = \tanh^{-1}(\epsilon_b)$$

SR-HBAC algorithm, for n=2



Round_($n=2$)

$$\epsilon_{max}^{(n=2)}$$

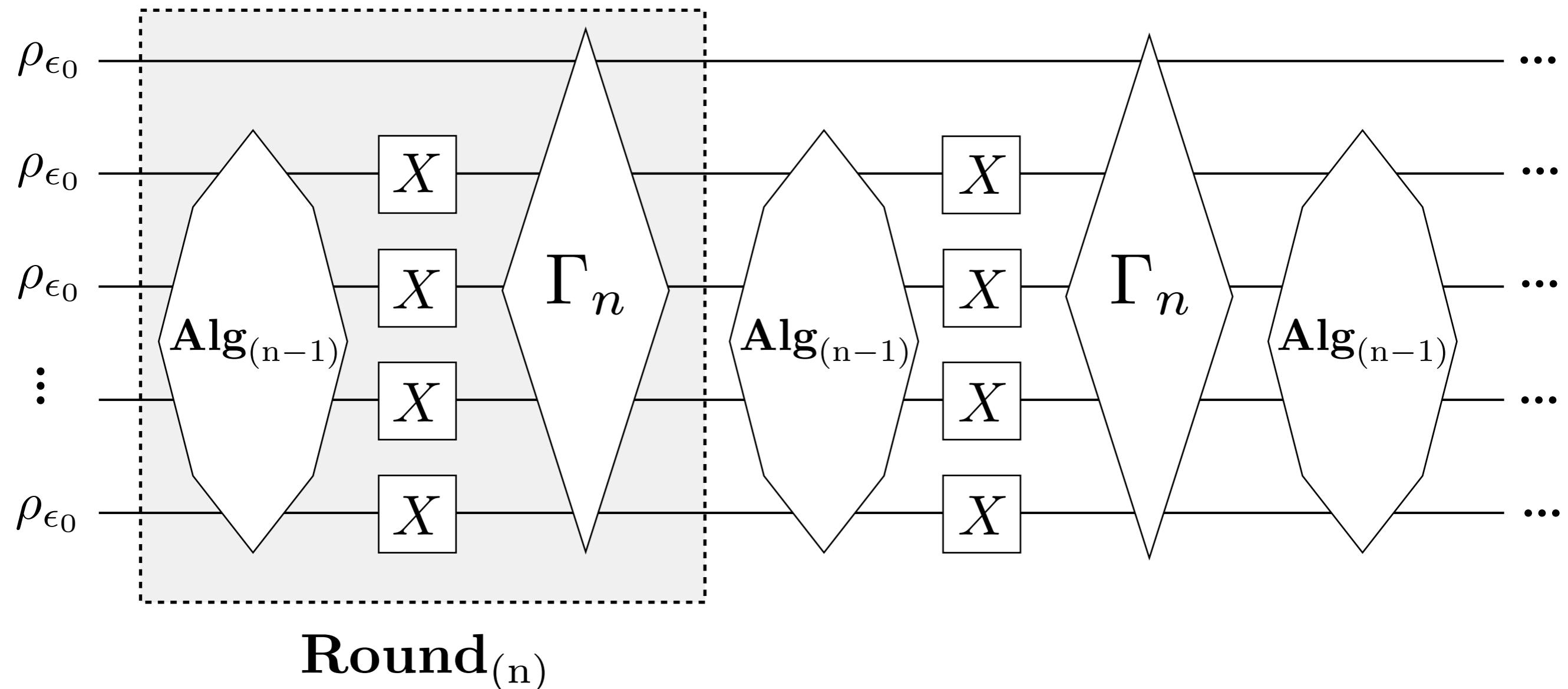


$$\epsilon_{max}^{(n=2)} = \tanh(3\xi)$$

$$\xi = \tanh^{-1}(\epsilon_b)$$

$$\epsilon_k^{(n=2)} = \left(3 - \frac{1}{2^{k-1}}\right) \epsilon_b$$

SR-HBAC algorithm for n qubits



$$\epsilon_{max}^{(n)} = \tanh [(2^n - 1) \xi]$$

$$\xi = \tanh^{-1} (\epsilon_b)$$

Final temperatures

PPA-HBAC

$$T_{\infty} = \frac{T_b}{md} \cdot \frac{\Delta E_t}{\Delta E_r}$$

using string of qubits with only one reset qubit:

$$T_{\infty} = \frac{T_b}{2^{n-2}} \cdot \frac{\Delta E_t}{\Delta E_r}$$

SR-HBAC

$$T_{\infty} = \frac{T_b}{2^n - 1} \cdot \frac{\Delta E_t}{\Delta E_r}$$

Chapter three

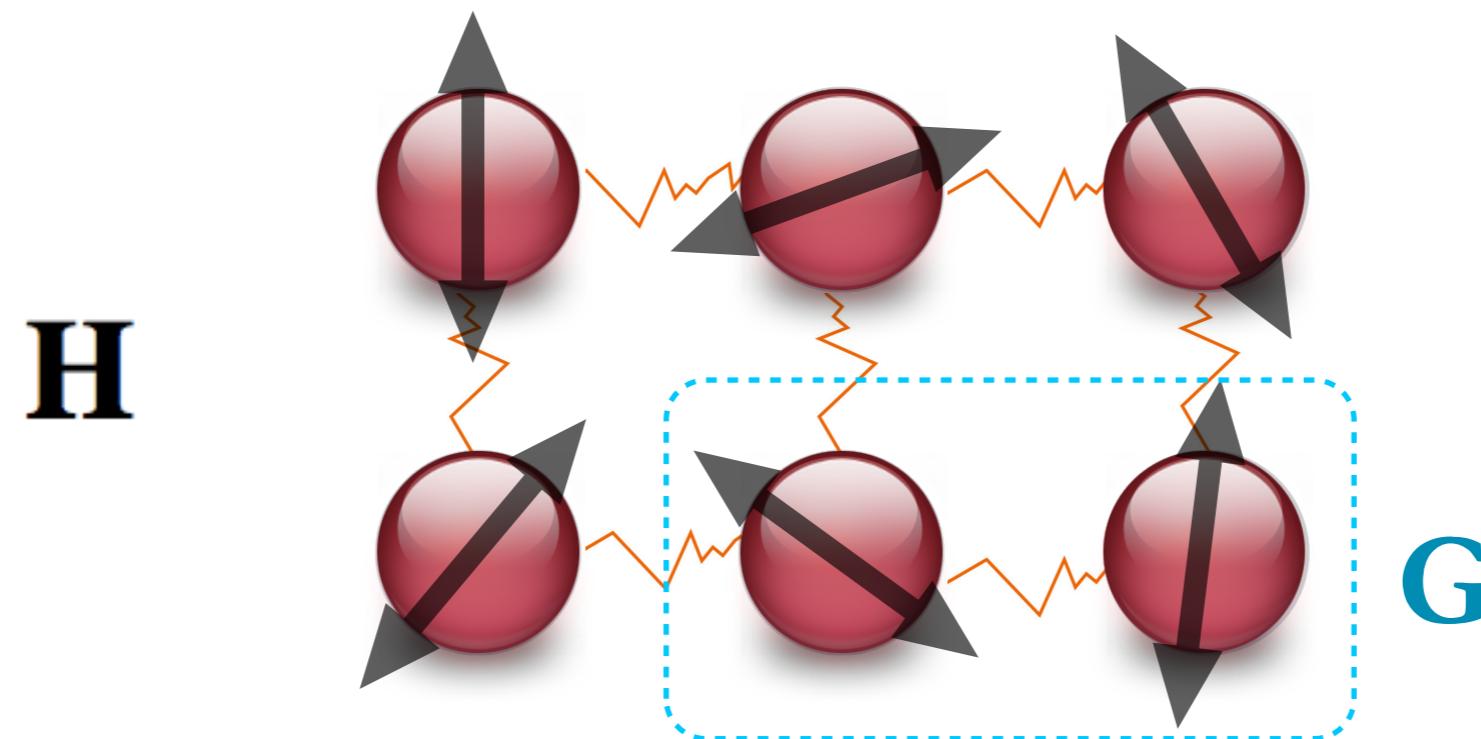
Correlations in the initial state

PPA-HBAC assumes that the initial state of the system is in a product state.

Idea: Remove this assumption, and use initial correlations for cooling

Motivation:

Purifying qubits from low temperature **interacting** quantum systems.
I.e., cooling local quantum systems subject to many-body interactions



Bad news for local cooling of interacting systems?

- Cool down the system:

Ground state is entangled \longrightarrow Subsystems are mixed

- Break entanglement through interaction with environment:

System in a Gibbs state \longrightarrow Subsystems are mixed

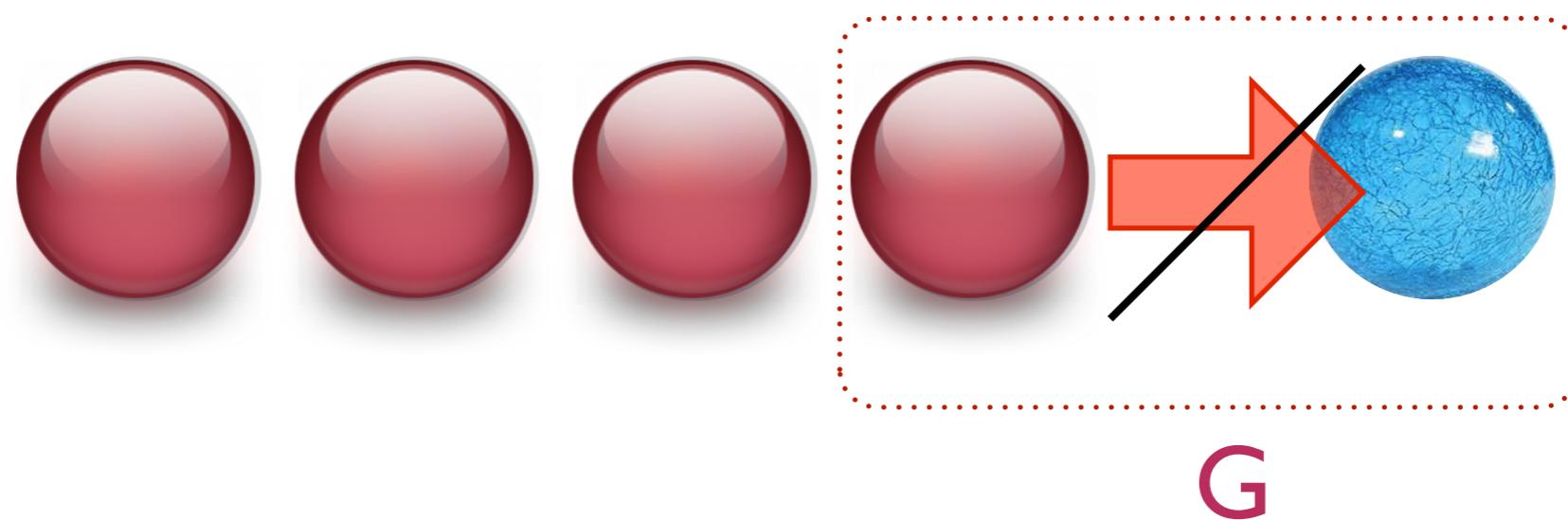
- Apply local operations on the subsystem:

Strong local passivity \longrightarrow Subsystem's energy increases

How do we cool down a part of an interacting system?

Strong local passive states

A system state is defined to be SL passive if no **general** quantum operation **G** applied **locally** to a system can extract positive energy from the system.



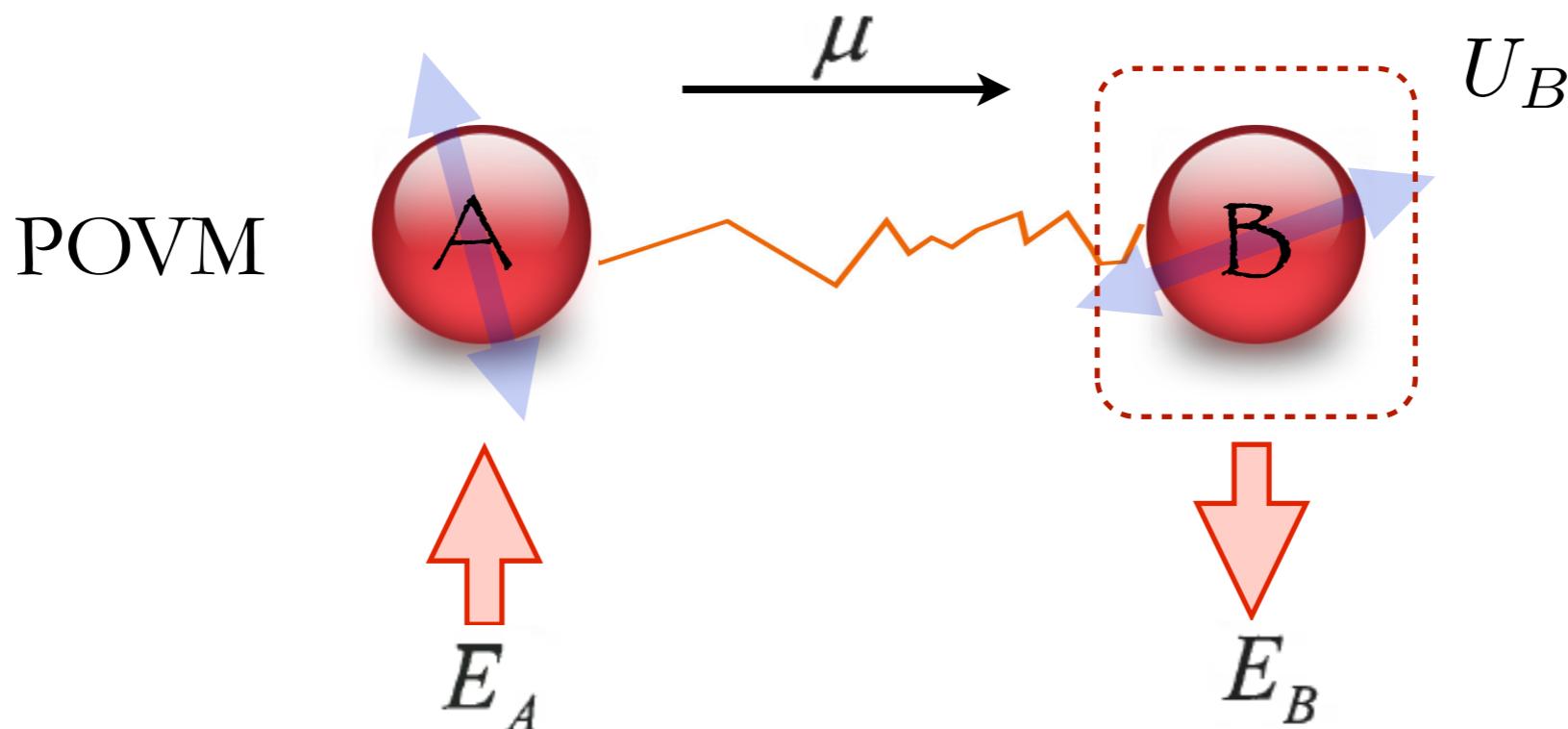
$$\Delta E (\rho) = \text{Tr} [H (I \otimes G) \rho] - \text{Tr} [H \rho]$$

$$\forall G, \Delta E \geq 0$$

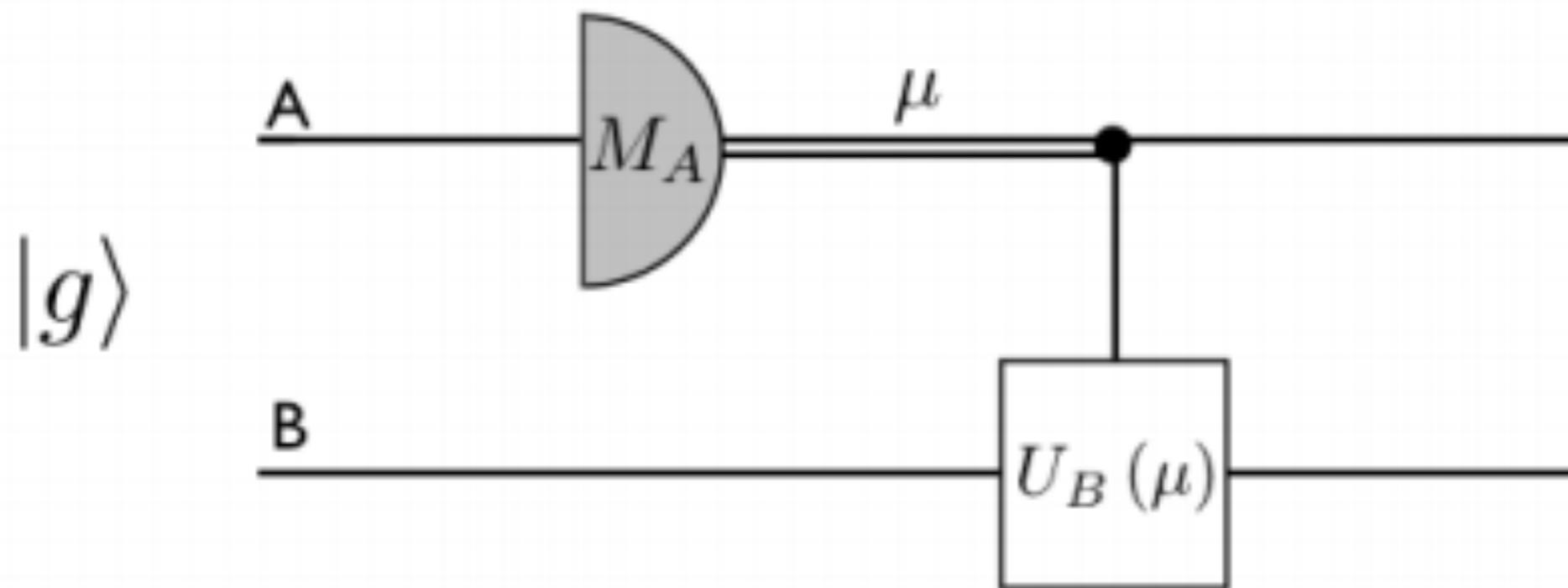
Taking advantage of pre-existing correlations

How?

Using an idea from quantum field theory, a method called quantum energy teleportation to consume correlations while extracting work locally from the target qubit.

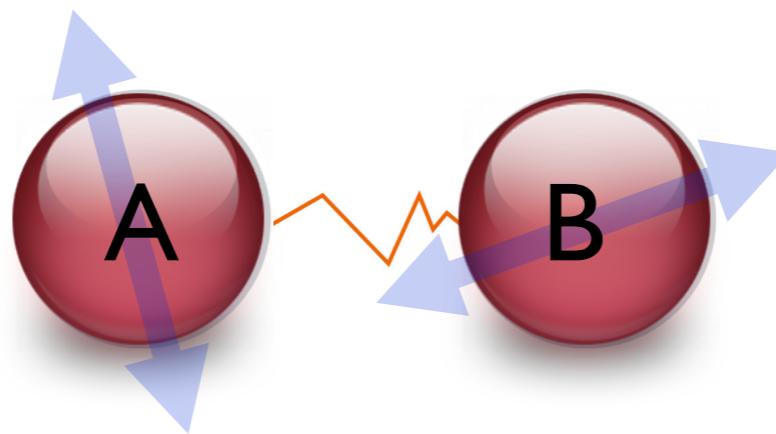


QET Circuit



POVM case with measurement operators

Two interacting qubits



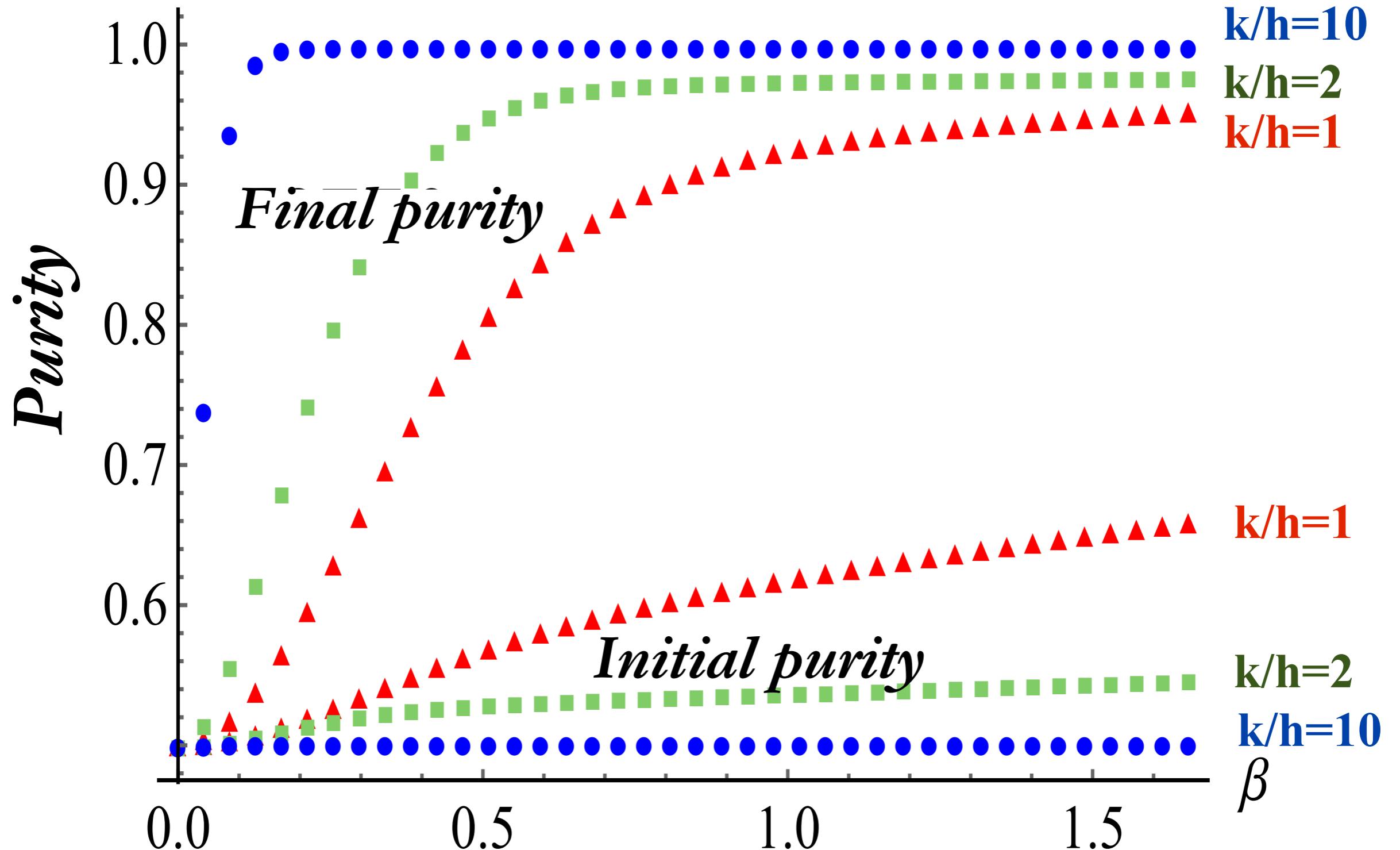
$$\hat{H} = \hat{H}_A + \hat{H}_B + \hat{V} \geq 0$$

$$\hat{H}_A = h\hat{\sigma}_A^z + \frac{h^2}{\sqrt{h^2 + k^2}},$$

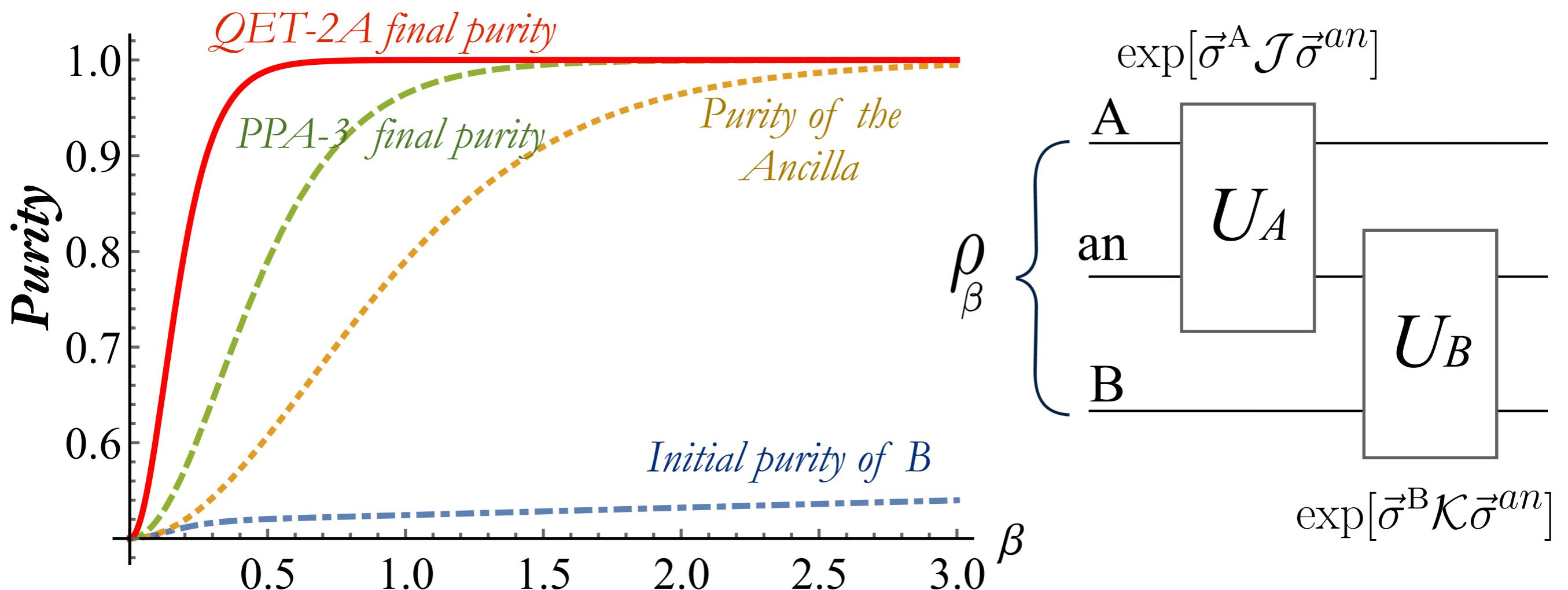
$$\hat{H}_B = h\hat{\sigma}_B^z + \frac{h^2}{\sqrt{h^2 + k^2}}, \quad (h > 0, k > 0)$$

$$\hat{V} = 2k\hat{\sigma}_A^x \hat{\sigma}_B^x + \frac{2k^2}{\sqrt{h^2 + k^2}}.$$

Results for Minimal QET



Fully unitary picture



Conclusions:

1) We give analytical results for the achievable cooling limits in the conventional HBAC.

We circumvented previous limits by using correlations:

2) Taking advantage of correlations that can be created during the re-thermalization step with the heat-bath.

3) Using correlations present in the initial state induced by the internal interactions of the system.

[1] NA Rodríguez-Briones, R Laflamme. PRL 116 (17), 170501

[2] DK Park, NA Rodriguez-Briones, G Feng, R Rahimi, J Baugh, R Laflamme. Electron Spin Resonance (ESR) Based Quantum Computing, 227-255

[3] NA Rodriguez-Briones, J Li, X Peng, T Mor, Y Weinstein, R Laflamme. NJP 19 (11), 113047

[4] NA Rodríguez-Briones, E Martín-Martínez, A Kempf, R Laflamme. PRL 119 (5), 050502