

# Quantum Mpemba effect in a quantum dot with reservoirs

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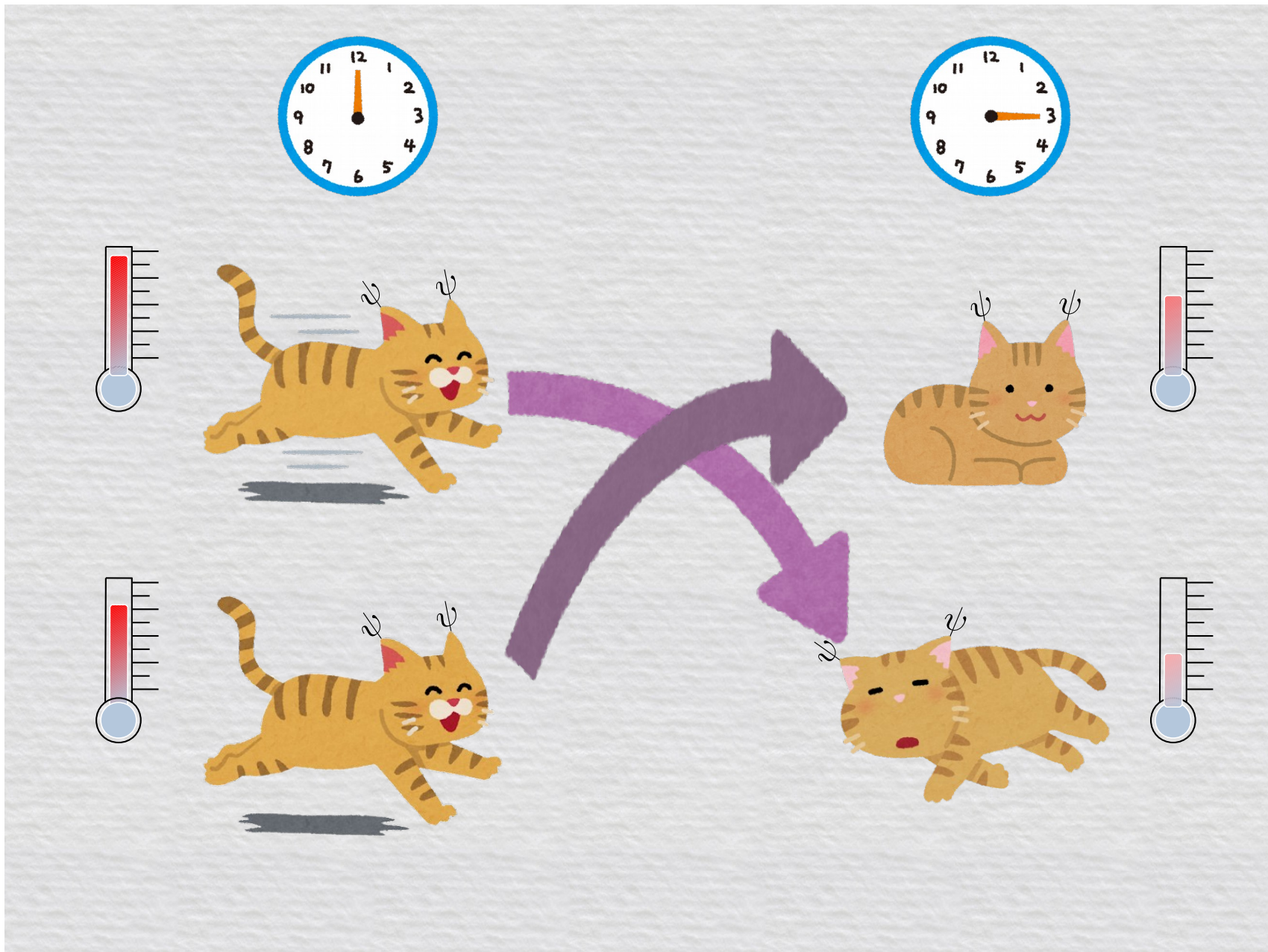
&

Satoshi Takada



[Ref: Amit Kumar Chatterjee, Satoshi Takada, and Hisao Hayakawa, arXiv:2304.02411]

(Accepted in Physical Review Letters)

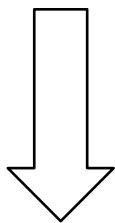


A quantum system (cat) can cool faster when starting from initially hotter rather than colder temperature.

# Mpemba effect

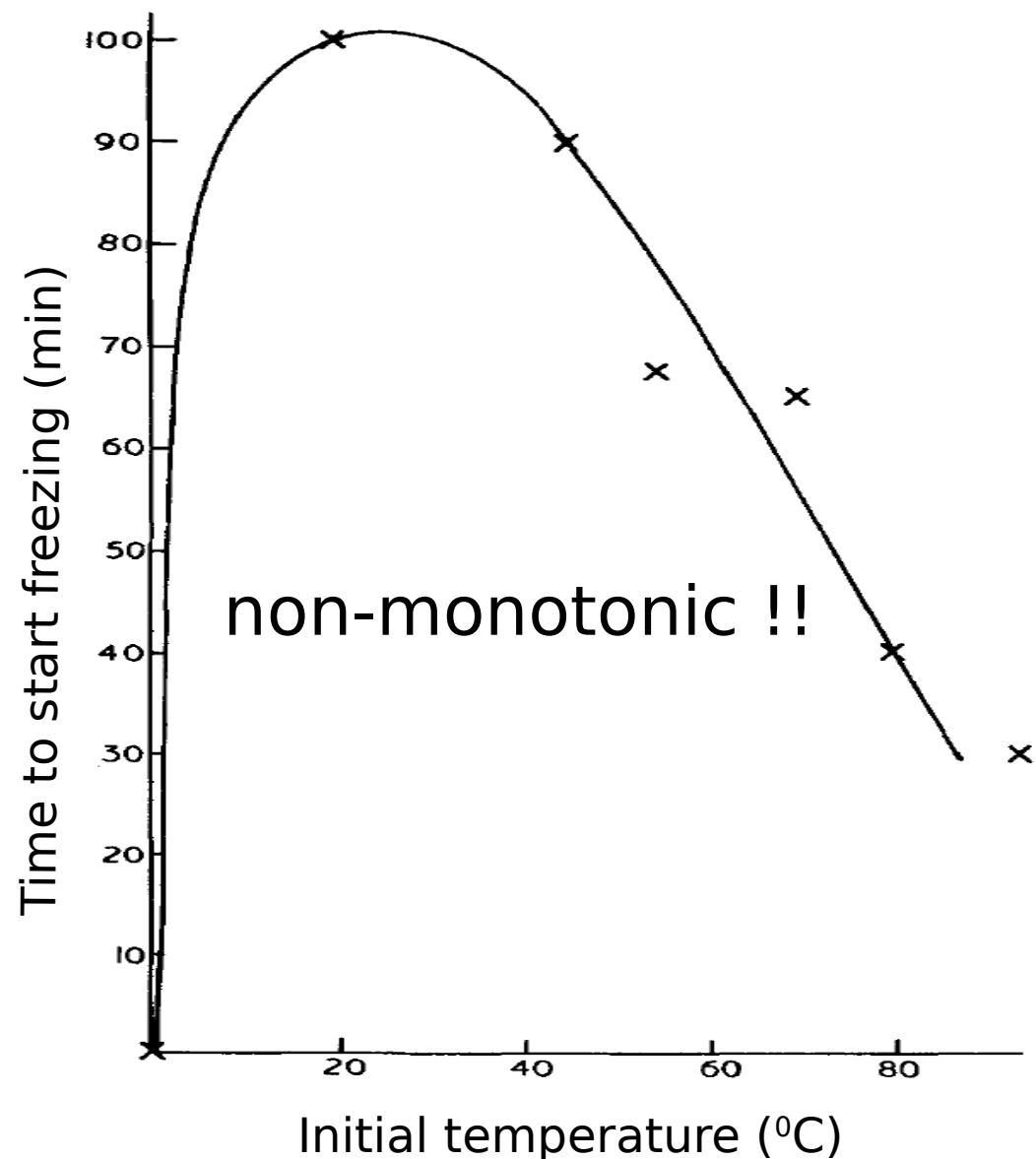
Experiment (Mpemba & Osborne):

freezing equal volumes  
of water samples at different  
initial temperatures



Hot water can freeze  
faster than cold water

[\* *thermal pheomena* \*]



# Mpemba effect as Generalized anomalous relaxation

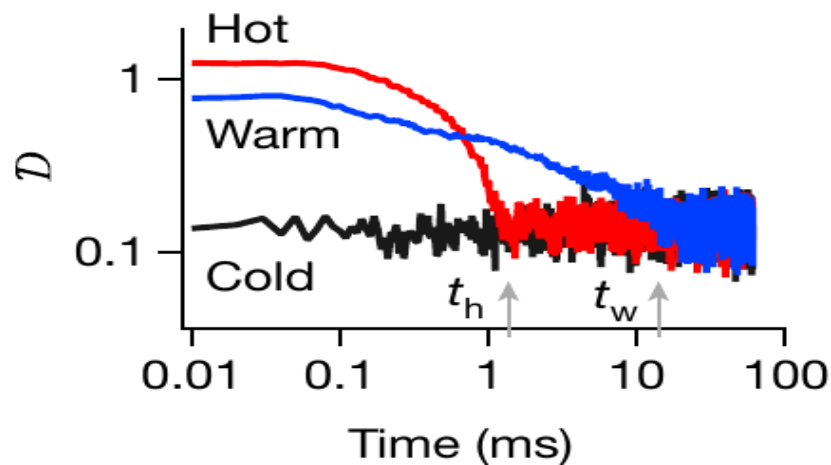
“time to start freezing”

replaced by

“crossing of trajectories” before reaching same steady state

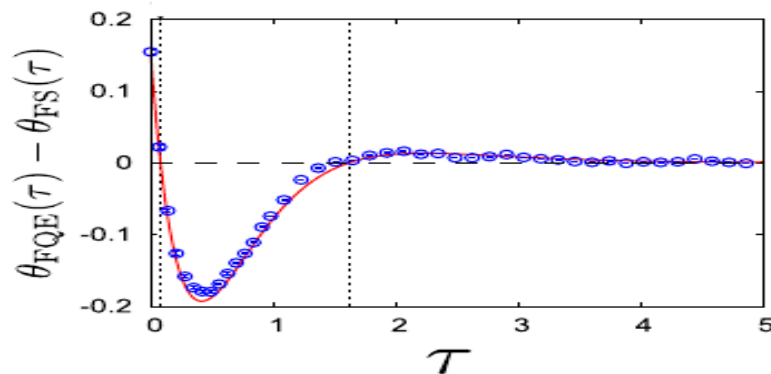
- Colloids

[Kumar and Bechhoefer,  
Nature 584, 64 (2020)]



- Inertial suspensions

[Takada, Hayakawa and Santos,  
PRE 103, 032901 (2021)]



• Markovian models: 
$$\frac{dp_i(t)}{dt} = \sum_j R_{ij}(T_b)p_j(t)$$

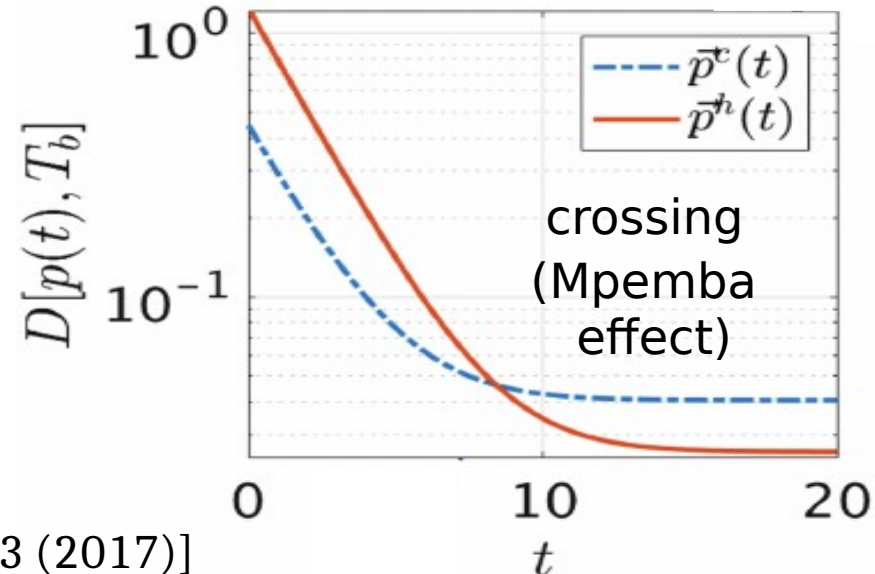
$$\vec{p}(t) = \vec{\pi}(T_b) + \underbrace{e^{\lambda_2 t} a_2 \vec{v}_2}_{\text{slowest relaxation}} + \underbrace{e^{\lambda_3 t} a_3 \vec{v}_3 + \dots + e^{\lambda_n t} a_n \vec{v}_n}_{\text{other relaxation modes}} \quad \lambda_2 > \lambda_3 > \dots > \lambda_n$$

Hot copy:  $\vec{p}^h(t) \approx \vec{\pi}(T_b) + a_2^h \vec{v}_2 e^{\lambda_2 t}$   
 Cold copy:  $\vec{p}^c(t) \approx \vec{\pi}(T_b) + a_2^c \vec{v}_2 e^{\lambda_2 t}$  } ignores initial relaxation

Distance function from equilibrium:

$$D_e[\vec{p}; T_b] = \sum_i \left( \frac{E_i \Delta p_i}{T_b} + p_i \ln p_i - \pi_i^b \ln \pi_i^b \right)$$

[Lu and Raz, PNAS 114, 5083 (2017)]



$$\vec{p}^h(t) \approx \vec{\pi}(T_b) + a_2^h \vec{v}_2 e^{\lambda_2 t}$$

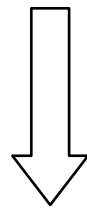
$$\vec{p}^c(t) \approx \vec{\pi}(T_b) + a_2^c \vec{v}_2 e^{\lambda_2 t}$$

Criterion for Mpemba effect (proposed by Lu & Raz):

$$|a_2^c| > |a_2^h|$$

But, what about the other relaxation modes ?

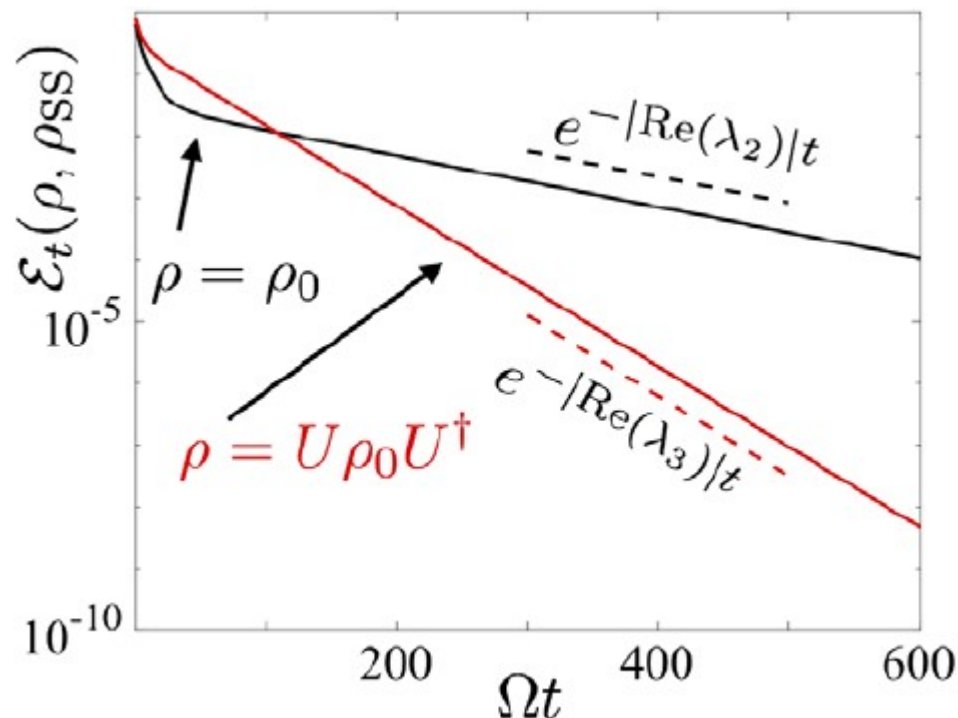
The co-efficients  $a_3, \dots, a_n$  have no role to play ?



Question: Role of initial relaxation on Mpemba effect ?

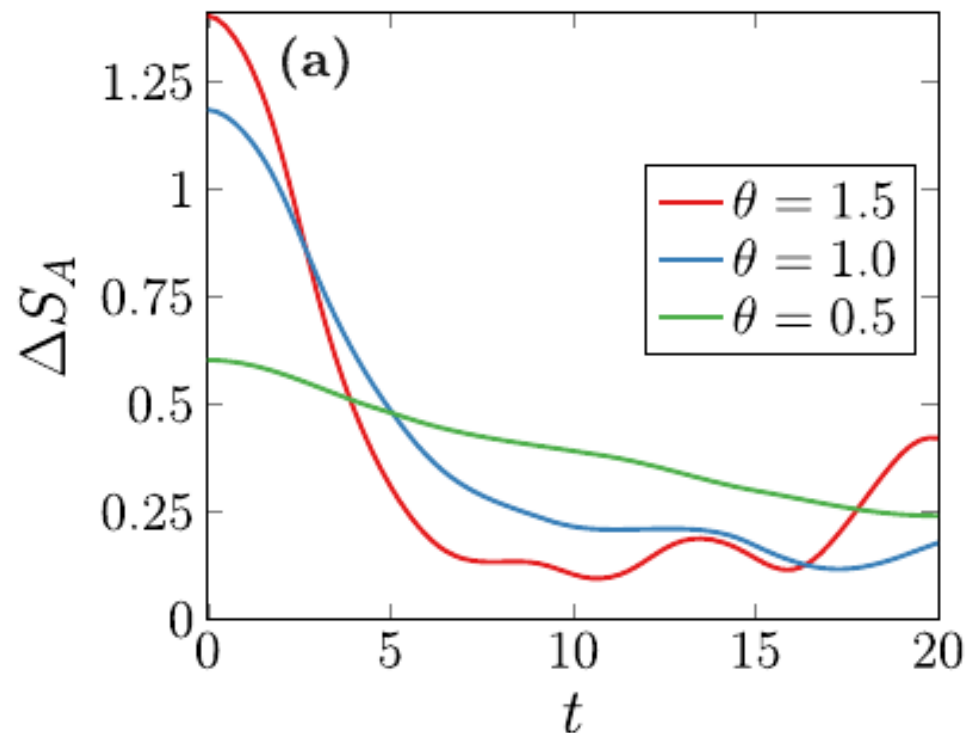
# Quantum Mpemba effect (QMPE)

vastly unexplored field



Distance from equilibrium in dissipative Dicke model

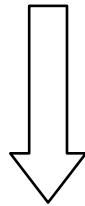
[Carollo, Lasanta and Lesanovsky, PRL 127, 060401 (2021)]



Entangle asymmetry in XXZ spin chain

[Ares, Murciano and Calabrese, Nature Communications 14, 2036 (2023)]

Analysis of temperature missing



Question

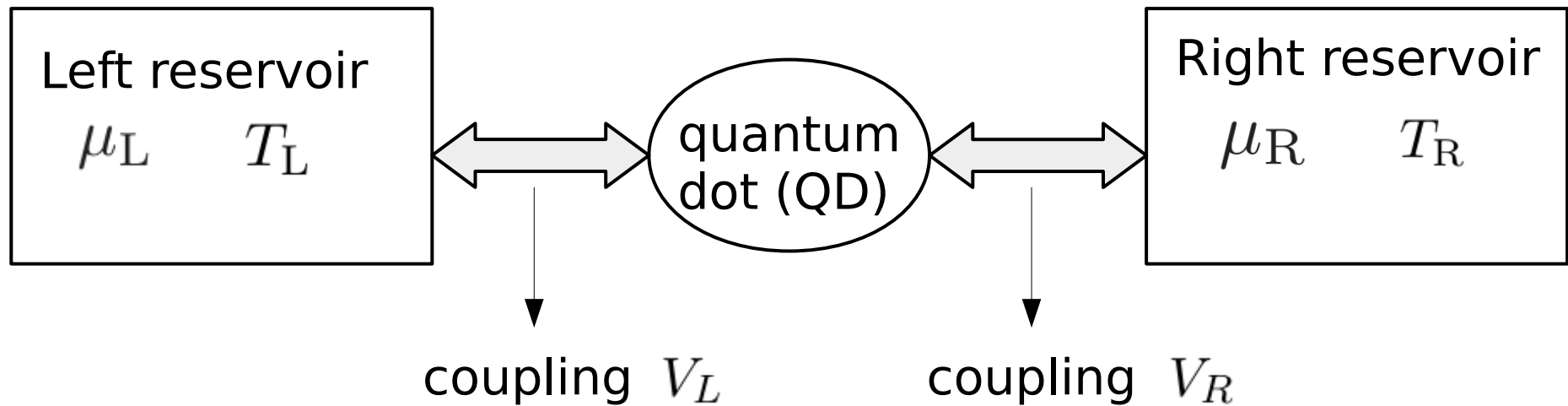
What about “thermal” Quantum Mpemba effect ?  
(*closer to original classical Mpemba effect*)



# Our motivations

- Show the existence of *thermal* quantum Mpemba effect
- Explore the roles of relaxation modes other than the slowest mode

# Quantum dot coupled to two reservoirs



QD states:  $\uparrow\downarrow, \uparrow, \downarrow, \text{vacant}$

Total Hamiltonian:  $\hat{H}_{\text{tot}} = \hat{H}_s + \hat{H}_r + \hat{H}_{\text{int}}$

System Hamiltonian      Reservoir Hamiltonian      System-reservoirs interaction

## Anderson model:

$$\hat{H}_s = \sum_{\sigma} \epsilon_0 \hat{d}_{\sigma}^{\dagger} \hat{d}_{\sigma} + U \hat{n}_{\uparrow} \hat{n}_{\downarrow}$$

$$\hat{H}_r = \sum_{\gamma, k, \sigma} \epsilon_k \hat{a}_{\gamma, k, \sigma}^{\dagger} \hat{a}_{\gamma, k, \sigma}$$

$$\hat{H}_{\text{int}} = \sum_{\gamma, k, \sigma} V_{\gamma} \hat{d}_{\sigma}^{\dagger} \hat{a}_{\gamma, k, \sigma} + \text{h.c.}$$

$\epsilon_0$ : energy of electron in quantum dot

$\epsilon_k$ : energy of electron corresponding to wave number  $k$  in reservoirs

$U$ : electron-electron interaction in quantum dot

$V_L, V_R$ : coupling strength between quantum dot and reservoirs

$\hat{d}^{\dagger}, \hat{d}$ : creation and annihilation operators in quantum dot

$\hat{a}^{\dagger}, \hat{a}$ : creation and annihilation operators in reservoirs

$\hat{n}$ : number operator ( $= \hat{d}^{\dagger} \hat{d}$ )

$\gamma$ : reservoir indices  $L, R$

$\sigma$ : up-spin ( $\uparrow$ ) or down-spin ( $\downarrow$ )

# Quantum Master equation:

$$\frac{d}{d\tau} \hat{\rho} = \hat{K} \hat{\rho}$$

$\hat{\rho}$  : four possible states:  $\uparrow\downarrow, \uparrow, \downarrow, \text{vacant}$   
(wide band approximation)

Transition matrix: 
$$\hat{K} = \begin{pmatrix} -2f_-^{(1)} & f_+^{(1)} & f_+^{(1)} & 0 \\ f_-^{(1)} & -f_-^{(0)} - f_+^{(1)} & 0 & f_+^{(0)} \\ f_-^{(1)} & 0 & -f_-^{(0)} - f_+^{(1)} & f_+^{(0)} \\ 0 & f_-^{(0)} & f_-^{(0)} & -2f_+^{(0)} \end{pmatrix}$$

where  $f_+^{(j)} := f_L^{(j)}(\mu_L, U, \epsilon_0, T) + f_R^{(j)}(\mu_R, U, \epsilon_0, T), \quad j = 0, 1$

$$f_-^{(j)} = 2 - f_+^{(j)}$$

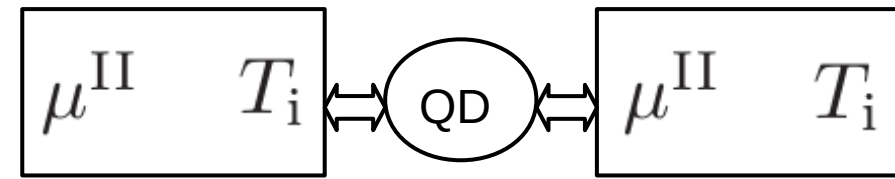
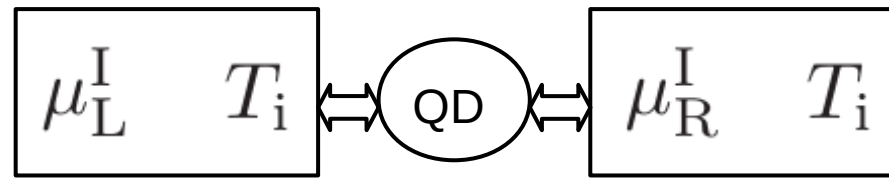
$$f_\gamma^{(j)}(\mu_\gamma, U, \epsilon_0, T) := \frac{1}{1 + e^{(\epsilon_0 + jU - \mu_\gamma)/T}} : \text{Fermi-Dirac distribution}$$

$$\gamma = \text{L, R}$$

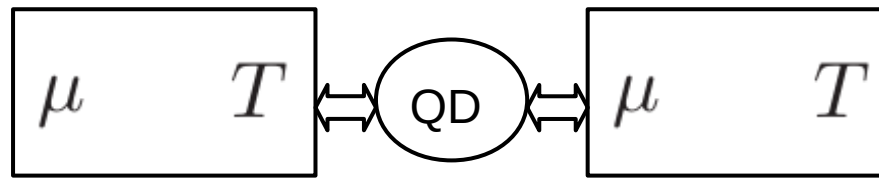
# Protocol

initial condition: I

initial condition: II



instantaneous quench



both copies  
evolve with time

If they cross  $\Rightarrow$  QMPE

If they don't cross  $\Rightarrow$  no QMPE

reach same steady state

# Thermal Quantum Mpemba effect

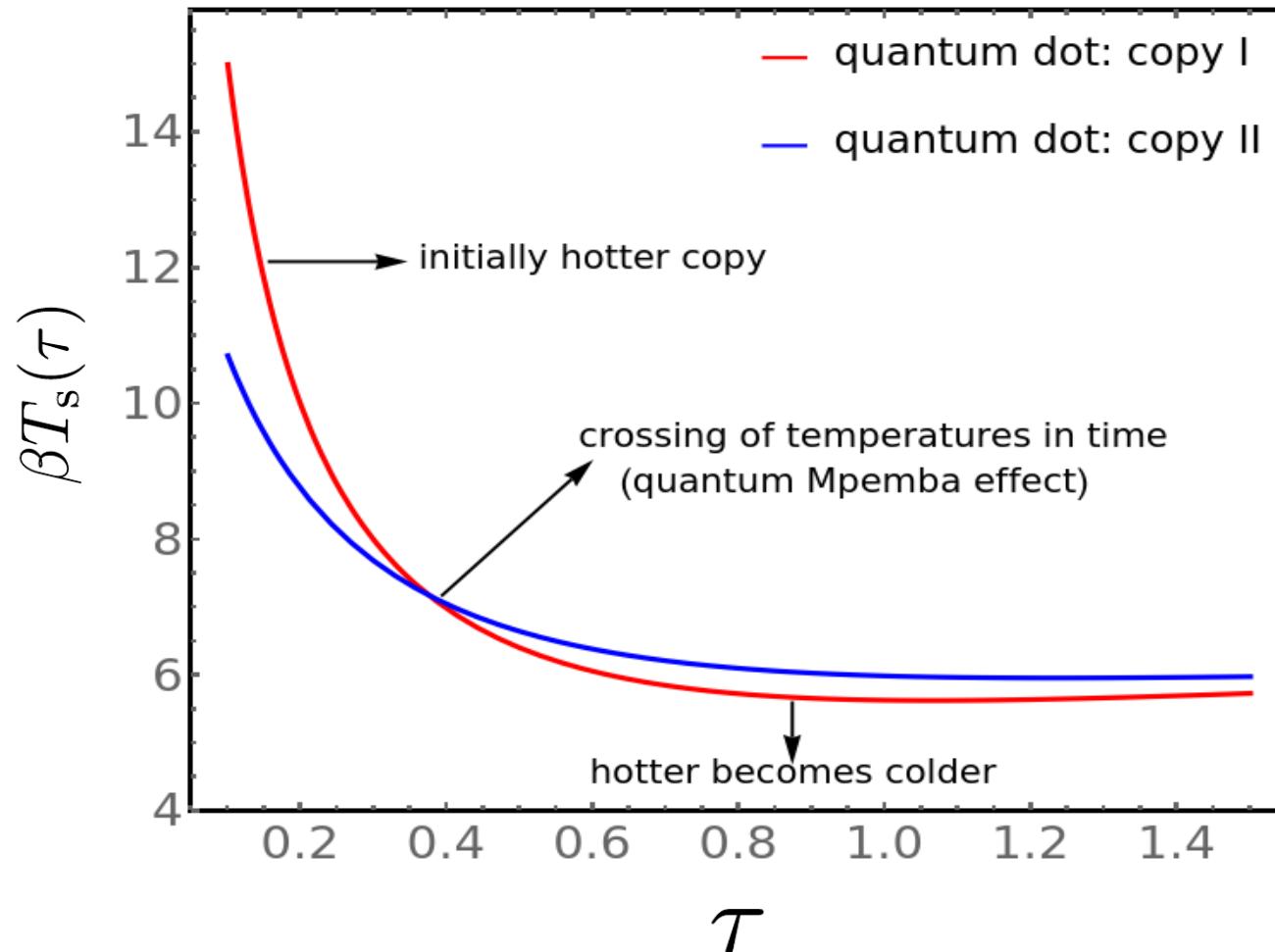
Temperature:

$$T_s(\tau) := \frac{\partial E_s(\tau)}{\partial S_{\text{vN}}(\tau)} = \frac{\partial E_s(\tau)}{\partial \tau} \bigg/ \frac{\partial S_{\text{vN}}(\tau)}{\partial \tau}$$

$$E_s(\tau) = \text{Tr}[\hat{\rho}(\tau)\hat{H}_s]$$

$$S_{\text{vN}}(\tau) = -\sum_{\alpha} \rho_{\alpha}(\tau) \ln(\rho_{\alpha}(\tau))$$

$$\beta = 1/T$$



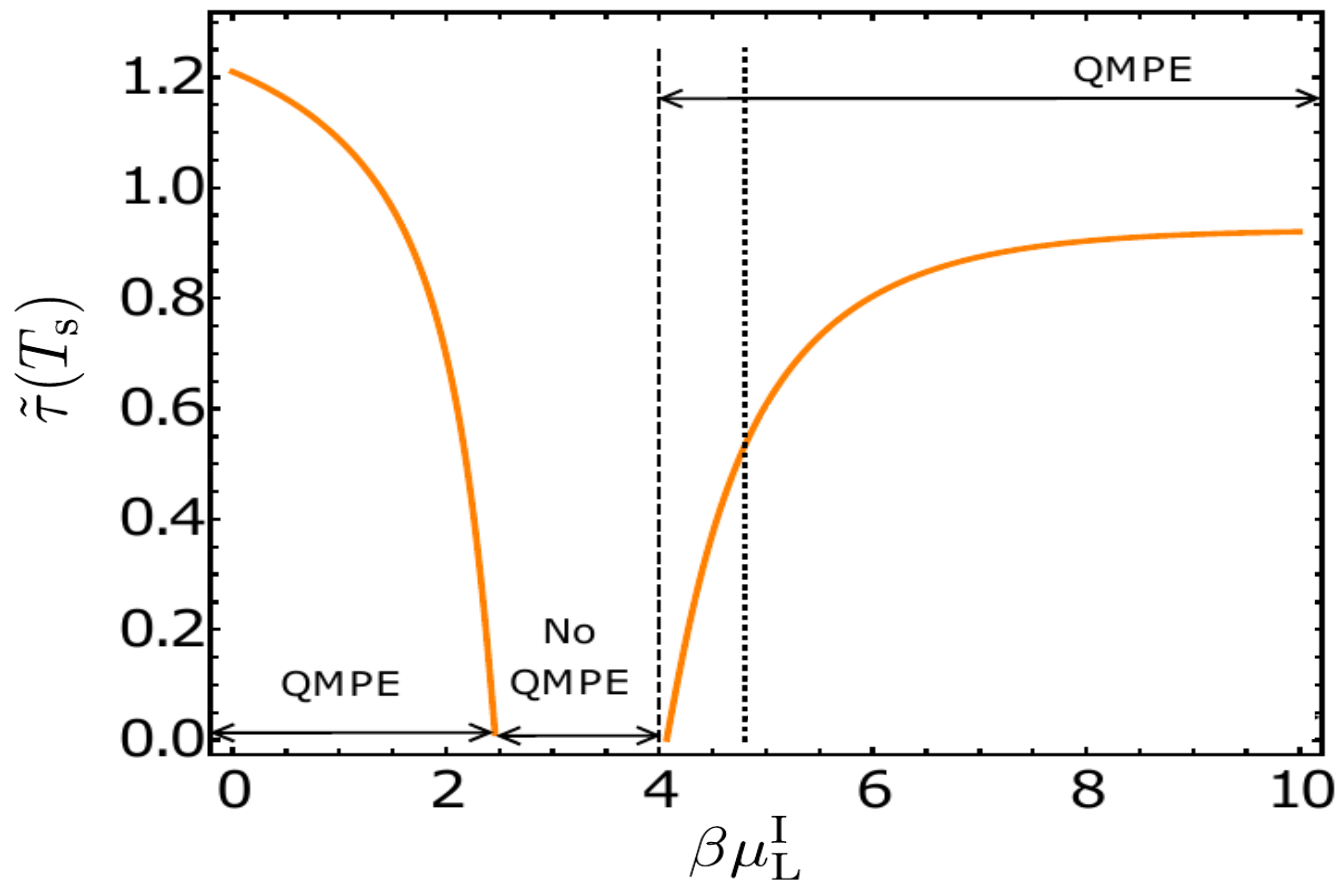
$$\beta\epsilon_0 = 2.0, \beta U = 1.25, \beta\mu_L^I = 4.5, \beta\mu_R^I = 1.0, \beta\mu^{II} = 2.43, \beta T_i = 1.15, \beta\mu = 2.0.$$

$$\Delta T_s := T_s^I - T_s^{II} \quad \tau = \tilde{\tau}(T_s) : \text{solution for } \Delta T_s = 0$$

$\tau$  is the **crossing time** (indicator of QMPE)

$0 < \tilde{\tau}(T_s) < \infty$  : thermal QMPE

$\tilde{\tau}(T_s) \rightarrow \infty$  : no QMPE



thermal QMPE  
is rather  
generic than  
occasional

$$\beta\epsilon_0 = 2.0, \beta U = 1.25, \beta\mu_R^I = 1.0, \beta\mu^{II} = 2.43, \beta T_i = 1.15, \beta\mu = 2.0.$$

# Role of relaxation modes

- $a_2$  is zero  $\implies$  No contribution from slowest relaxation mode
- To show QMPE in density matrix elements:

$$\begin{aligned}\Delta\rho_\alpha(\tau) &:= \rho_\alpha^{\text{I}}(\tau) - \rho_\alpha^{\text{II}}(\tau), \quad \alpha = 1, 2, 3, 4 \quad (\equiv \uparrow\downarrow, \uparrow, \downarrow, \text{vacant}) \\ &= e^{\lambda_3\tau} \hat{R}_{\alpha,4} \Delta a_4 \left[ S_\alpha + e^{-(\lambda_3 - \lambda_4)\tau} \right]\end{aligned}$$

$$S_\alpha := (\hat{R}_{\alpha,3} \Delta a_3) / (\hat{R}_{\alpha,4} \Delta a_4)$$

combined role of the surviving relaxation modes

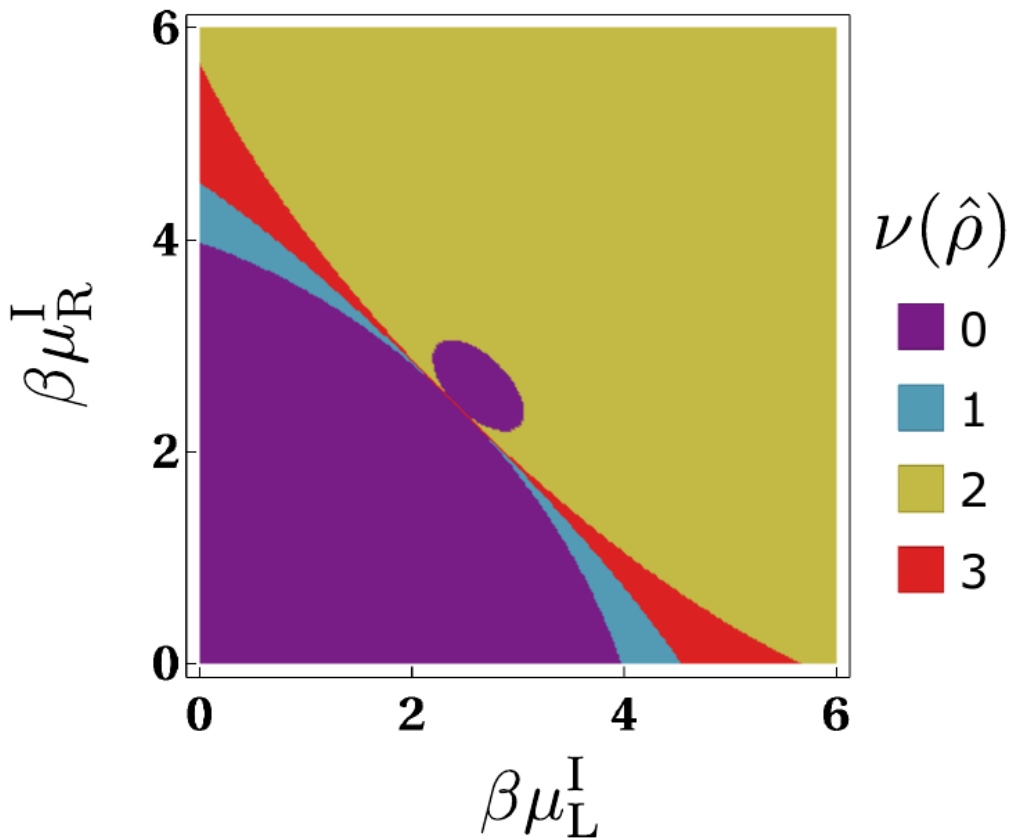


Necessary criterion for QMPE:

$$S_\alpha < 0 \quad \& \quad |S_\alpha| < 1$$



$\nu(\hat{\rho})$  : Number of density matrix elements showing QMPE



$$\beta\epsilon_0 = 2.0, \beta U = 1.25,$$

$$\beta\mu^{II} = 2.43, \beta T_i = 1.15, \beta\mu = 2.0.$$

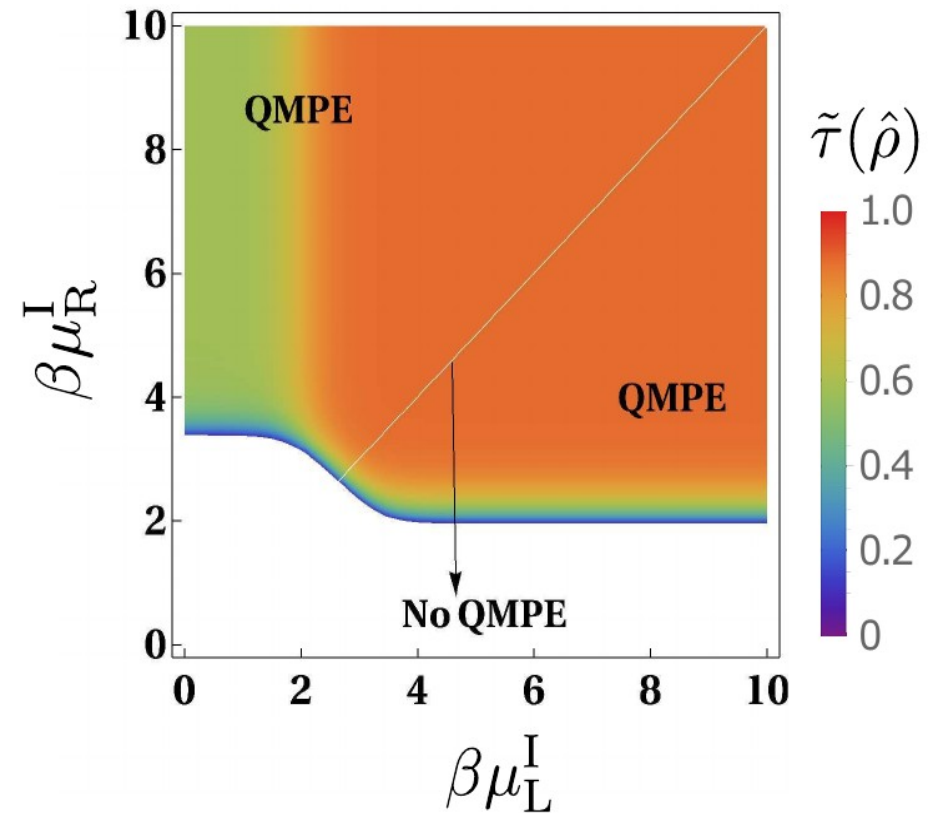
To get QMPE in density matrix elements:

$$\Delta\rho_\alpha(\tau_\alpha) = 0$$

$$\tilde{\tau}(\hat{\rho}) = \max[\tau_1, \tau_2, \tau_3, \tau_4]$$

if  $0 < \tau_\alpha < \infty$

$$\tilde{\tau}(\hat{\rho}) \rightarrow \infty \text{ if no finite } \tau_\alpha \text{ exists } \forall \alpha$$



$$\beta\epsilon_0 = 2.0, \beta U = 1.25,$$

$$\beta T_i = 0.25, \beta\mu = 2.0 \text{ and } \beta\mu^{II} = \beta\mu_R^I$$

# Summary

- A quantum system can cool faster when it starts from hotter initial temperature rather than colder initial temperature. This is the *thermal quantum Mpemba effect*.

(demonstrated in a single level quantum dot with reservoirs)

- The criterion for quantum Mpemba effect explicitly depends on the combined effect of relaxation modes other than the slowest relaxation mode.

(different from previously proposed criterion for classical Markovian systems)

[**Ref:** Amit Kumar Chatterjee, Satoshi Takada, and Hisao Hayakawa, arXiv:2304.02411]

[Accepted in Physical Review Letters]

[Selected for American Physical Society's outreach to the press]

THANK YOU