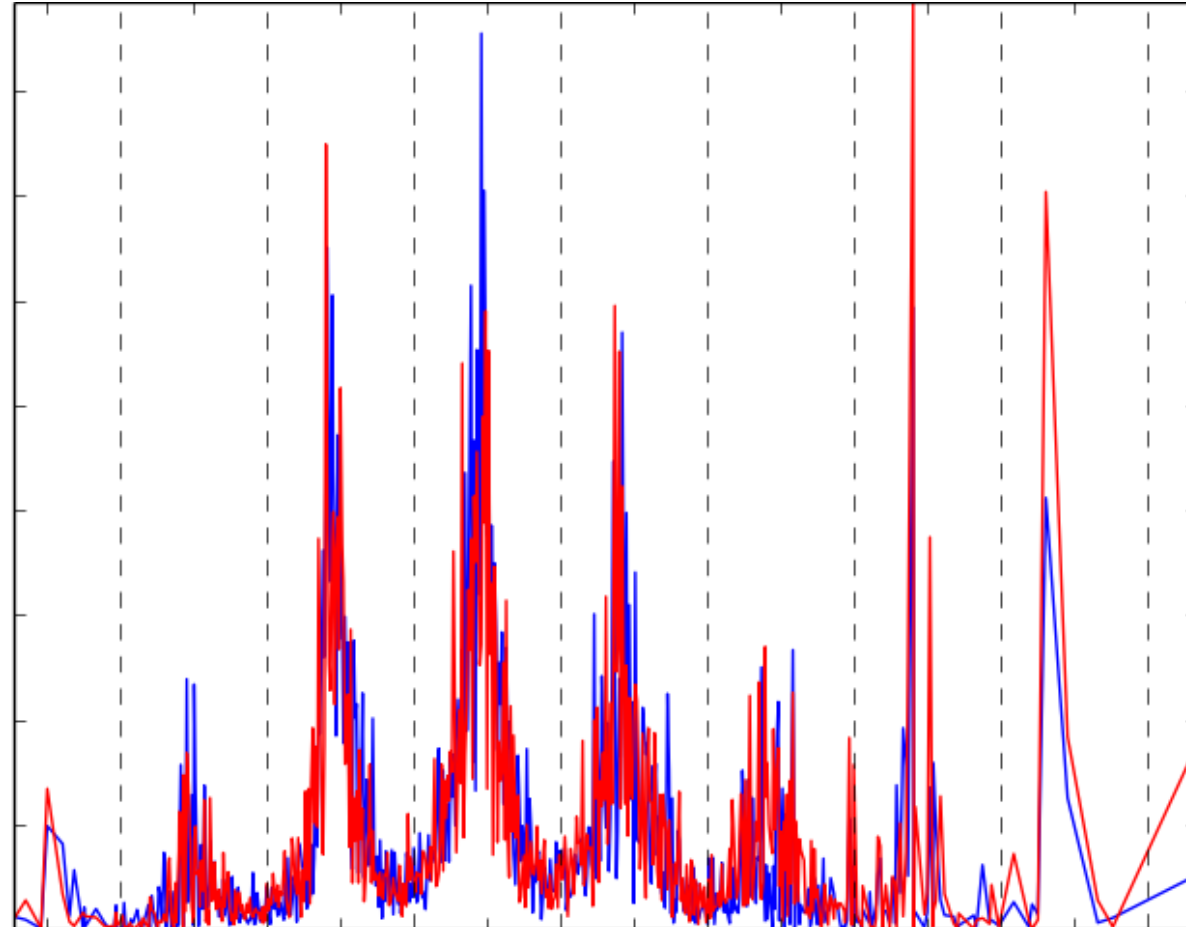


(Pre-)thermalization in Periodically-Driven Systems



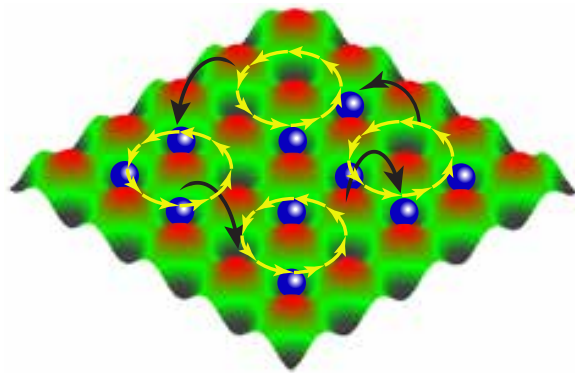
quantum or classical?

M.B., Heyl, Huse, Polkovnikov, PRB 93, 155132 (2016)

Howell, Weinberg, Sels, Polkovnikov, and M.B., PRL 122, 010602 (2019)

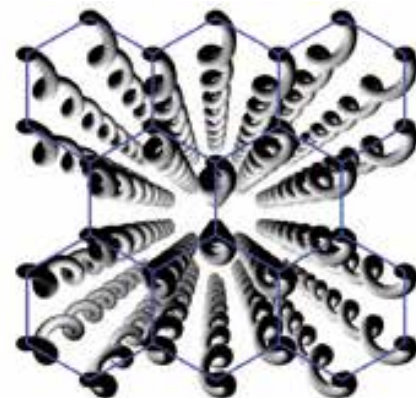
Why study periodically-driven systems?

➔ practically: Floquet engineering



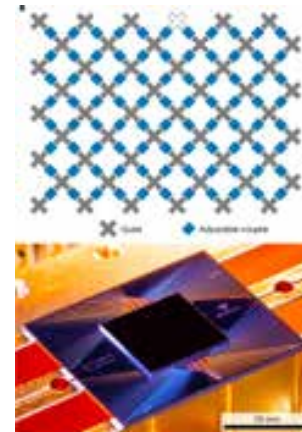
ultracold atoms

PRA, 90:043613, 2014



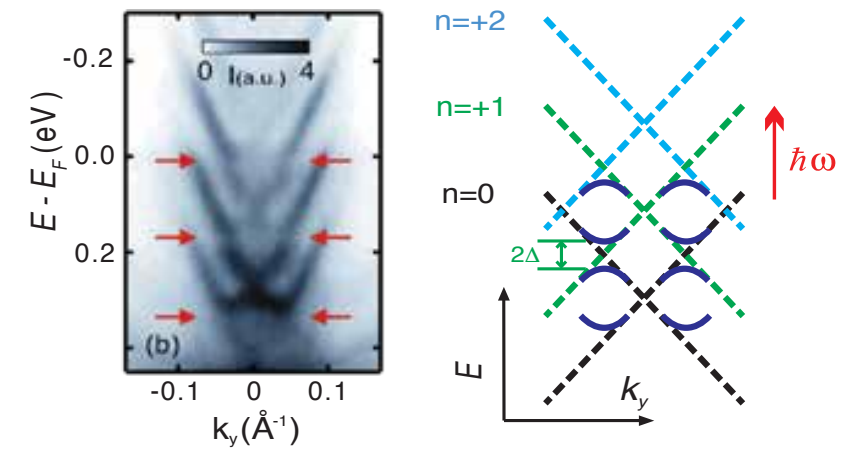
photonic insulators

Nature, 496:196, 2013



superconducting qubits

Nature, 574:505 2019

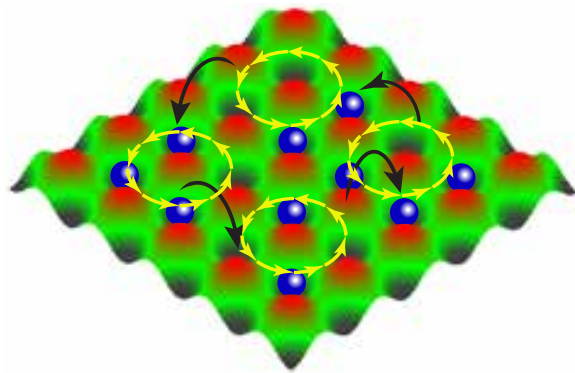


quantum materials: Bi_2Se_3

Nature Physics, 12(4):306, 2016

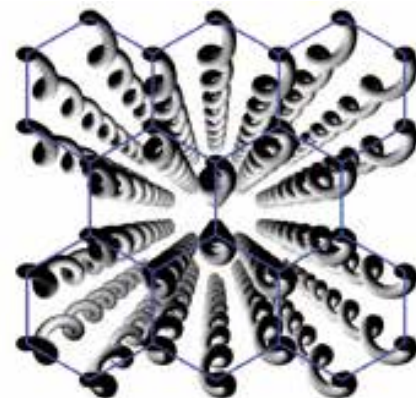
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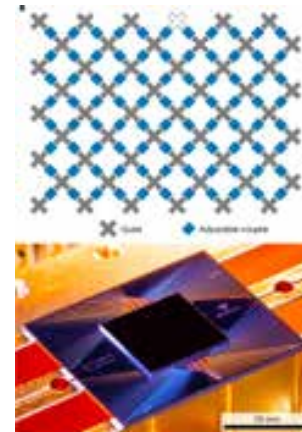
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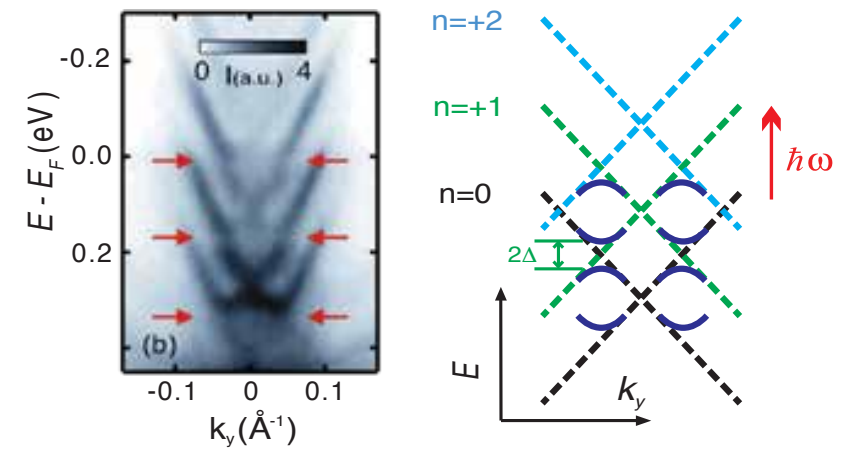
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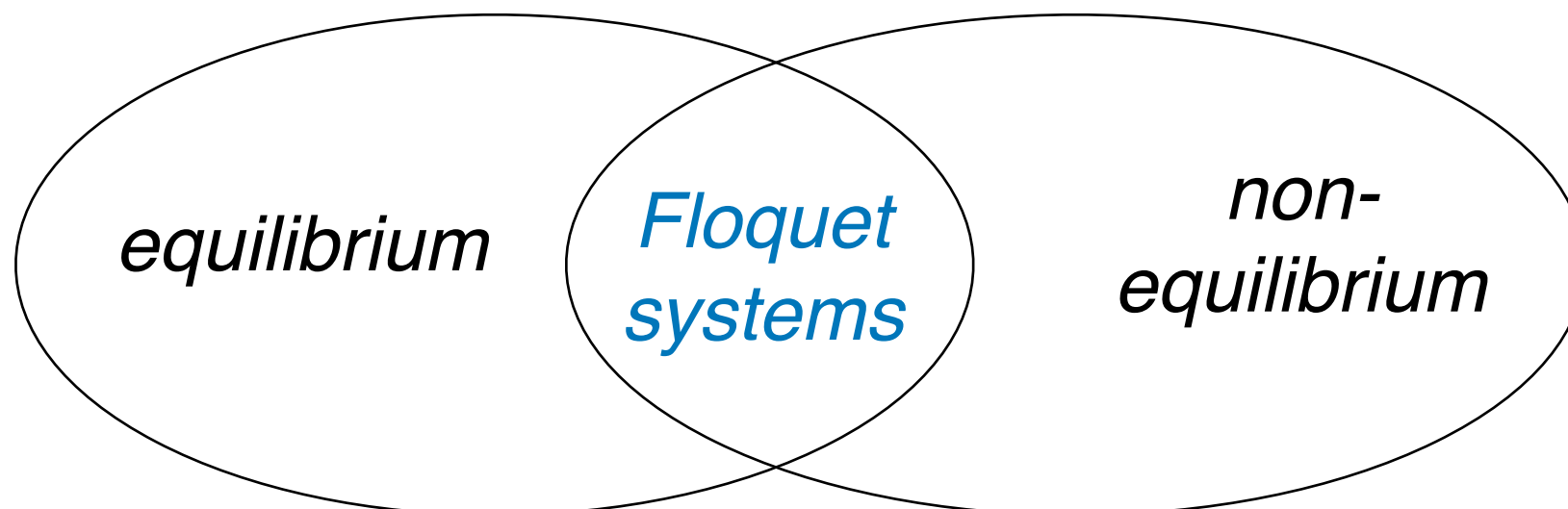
Nature, 574:505 2019



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Nature Physics, 12(4):306, 2016

→ fundamentally:



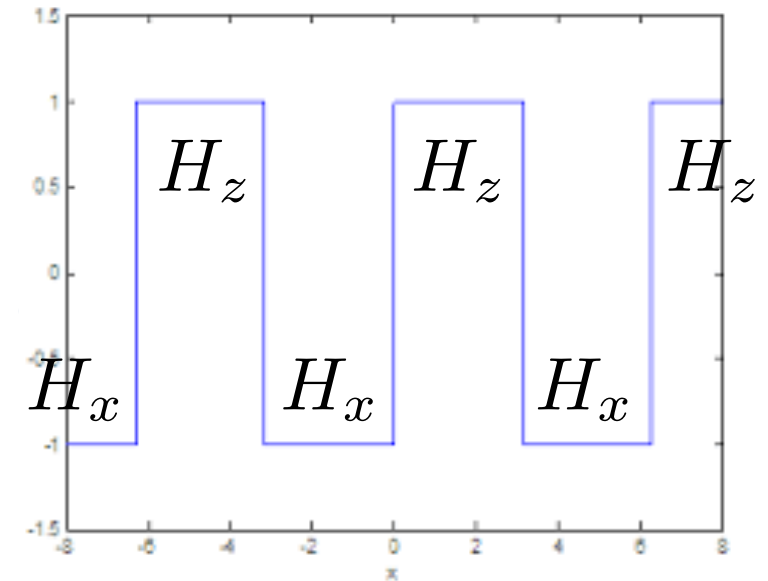
Energy absorption in quantum Floquet systems

$$\Omega = \frac{2\pi}{T}$$

→ generic local interacting driven system:

$$H(t) = \begin{cases} \sum_{j=1}^L J S_j^z S_{j+1}^z + h S_j^z & \text{for } t \in [0, T/2] \bmod T \\ \sum_{j=1}^L g S_j^x & \text{for } t \in [T/2, T] \bmod T \end{cases}$$

$$[S_i^\alpha, S_j^\beta] = i\delta_{ij}\varepsilon^{\alpha\beta\gamma} S_j^\gamma$$



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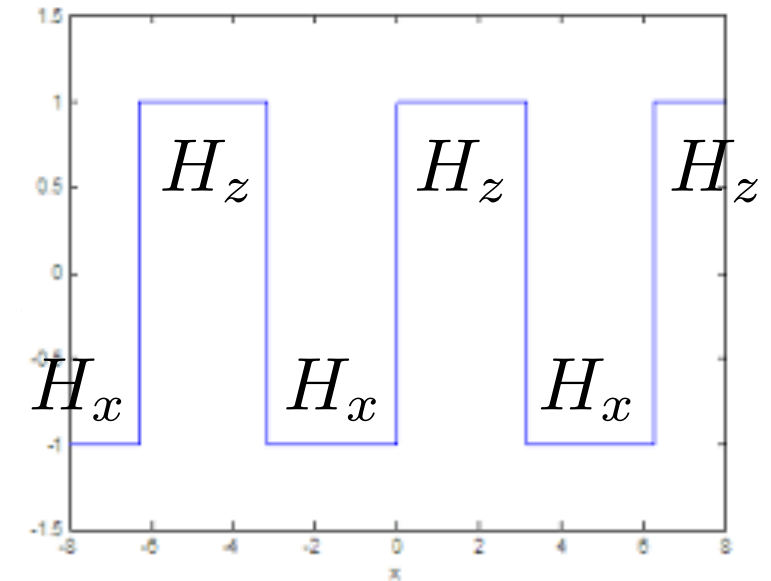
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→ **large** switching frequency: $\Omega \gg J, h, g$

average energy:

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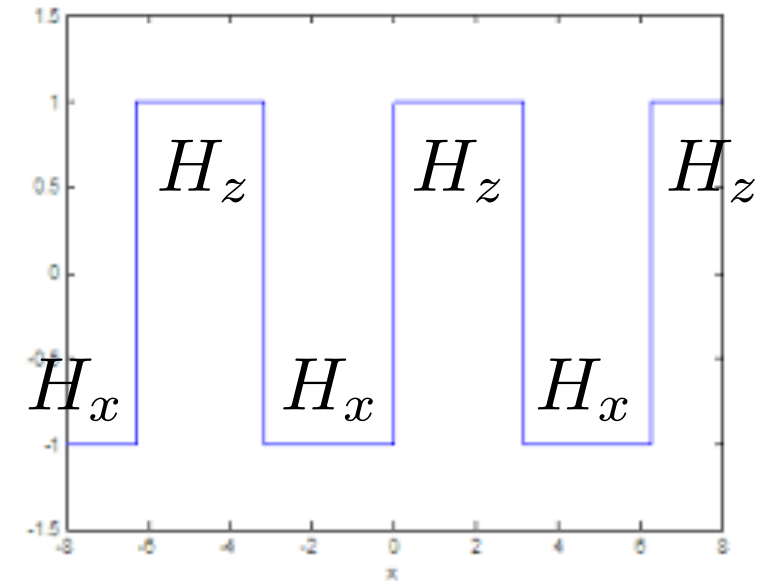
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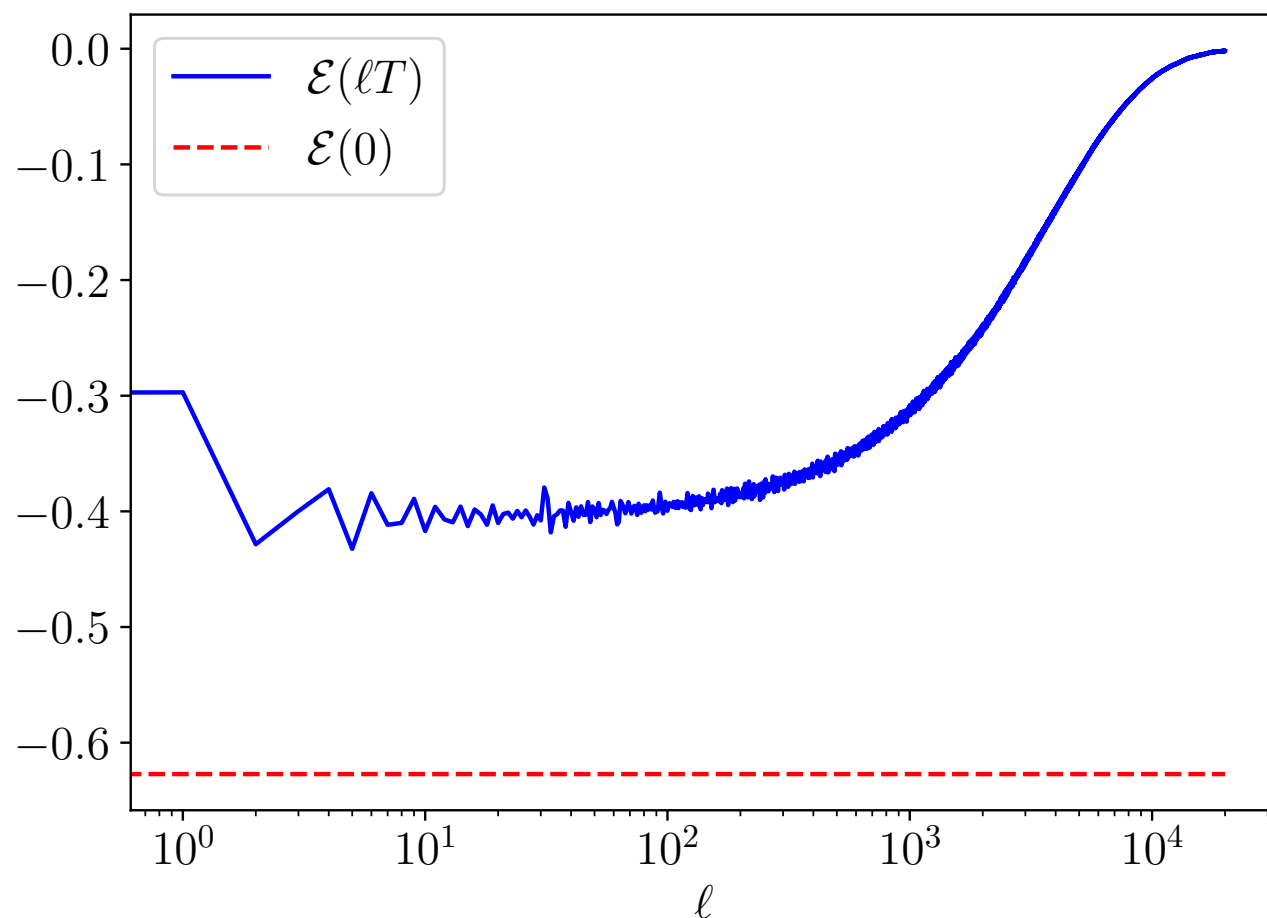
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$$|\psi(t=0)\rangle = |\text{GS}(H_{\text{ave}})\rangle$$

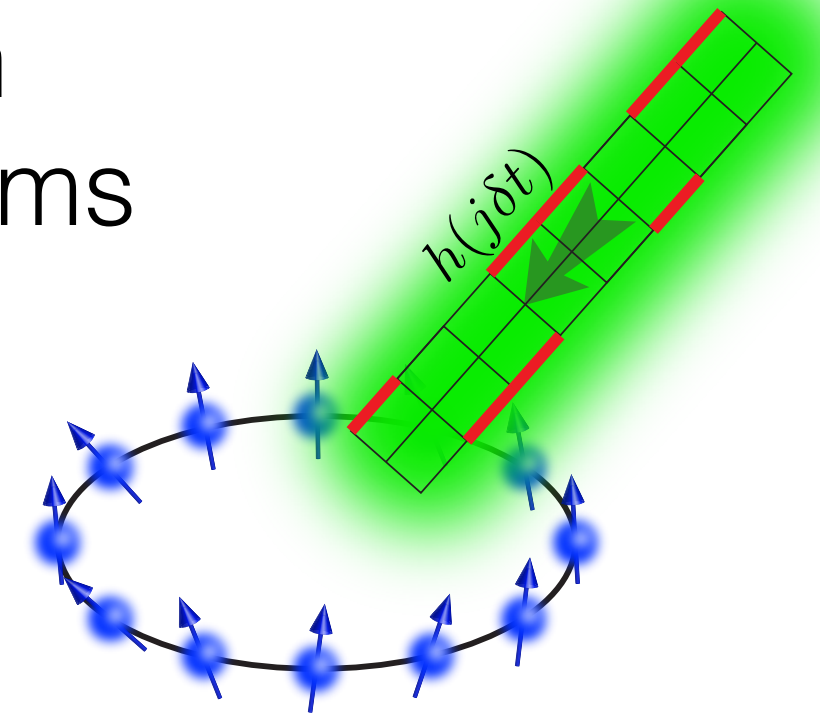
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D'Alessio and Rigol, PRX 4, 041048 (2014)

Lazarides et al, PRL 112, 150401 (2014), PRE 90, 012110 (2014).

Pre-thermalization in quantum Floquet systems

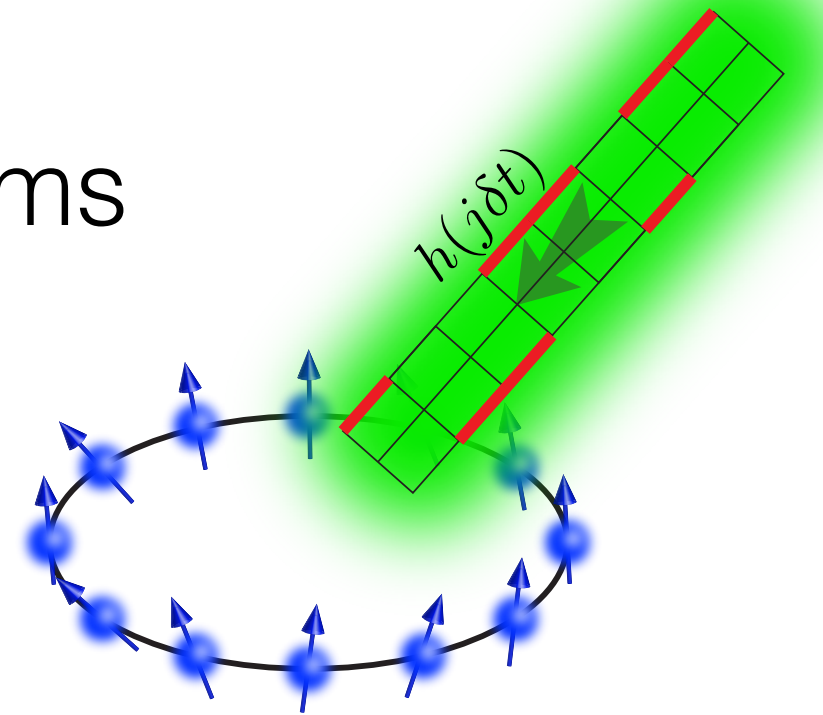
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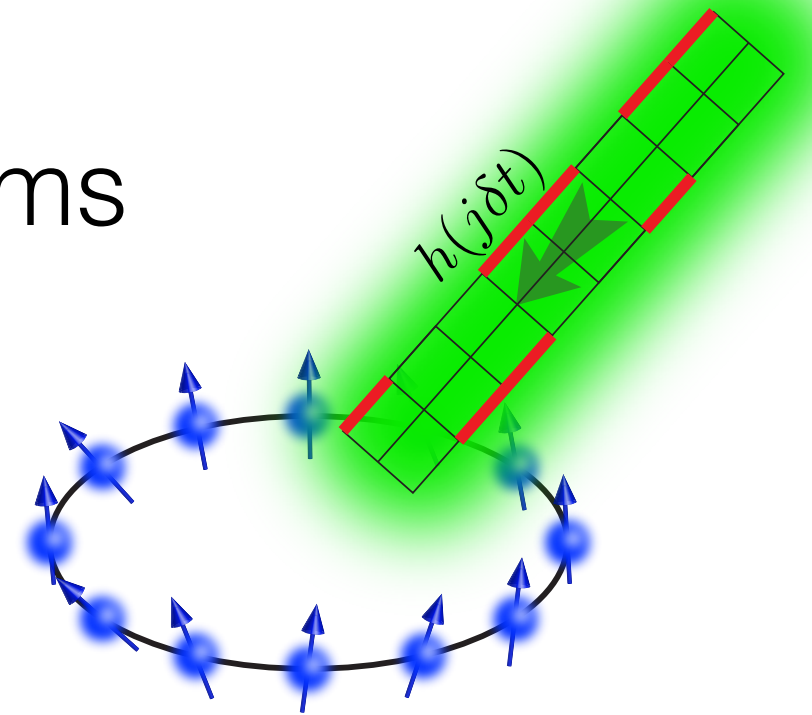


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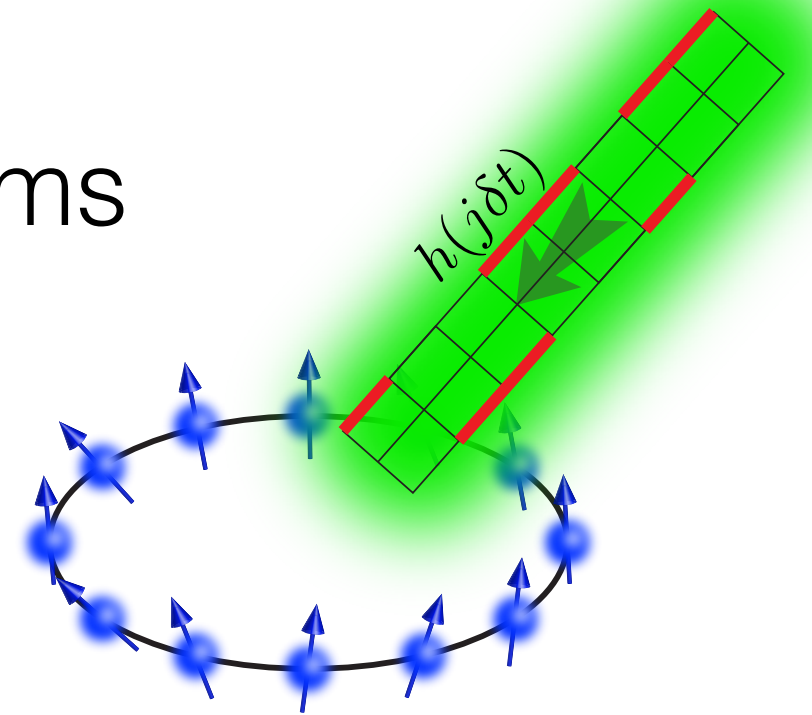


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long pre-thermal window: ideal for Floquet engineering

Floquet theory

→ Floquet's theorem

$$U(t, 0) = P(t) \exp(-itH_F) P^\dagger(0)$$

time-periodic $P(t + T) = P(t)$

Floquet Hamiltonian,
time-**in**dependent

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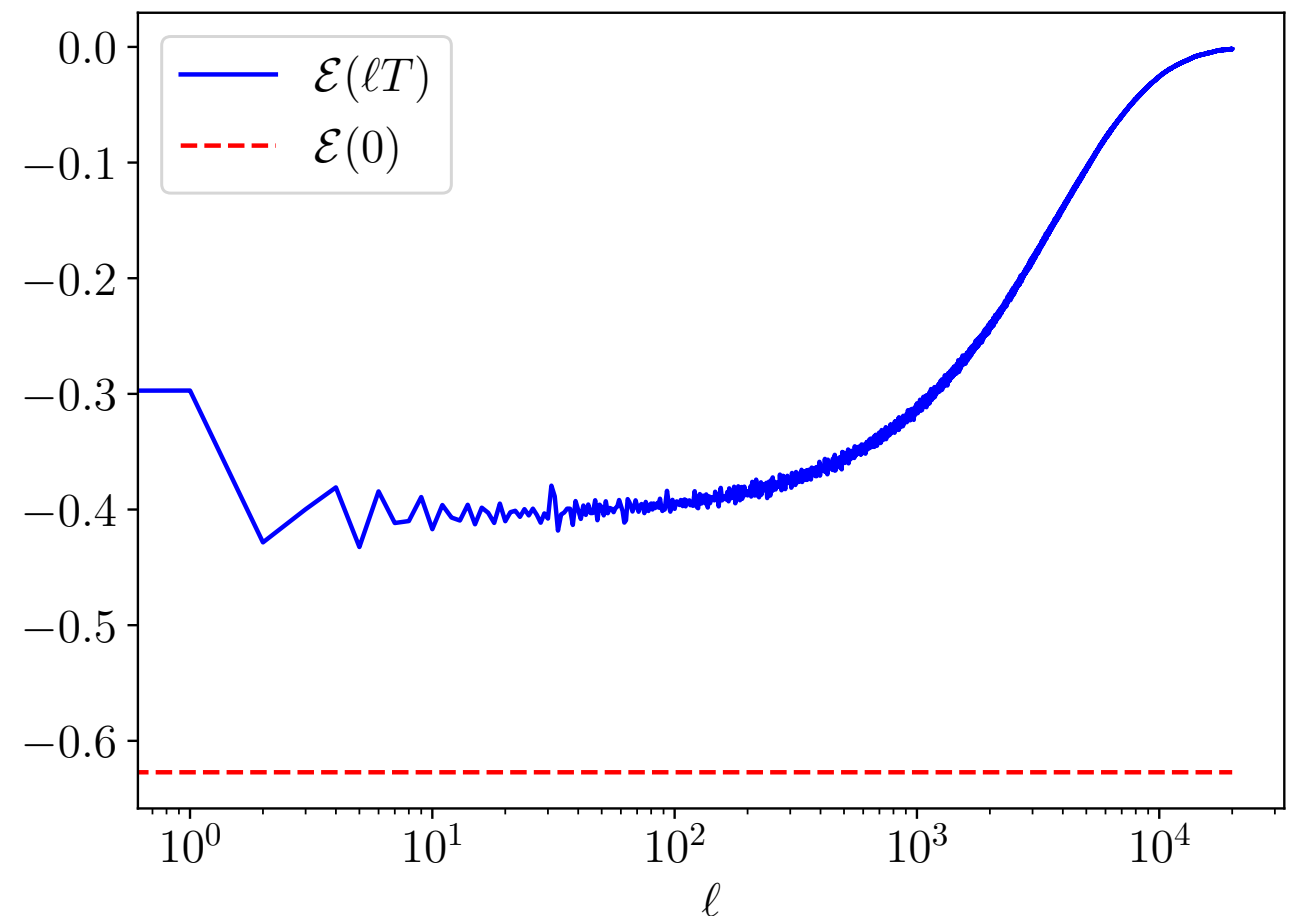
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Understanding the pre-thermal plateau

$$\begin{aligned}\mathcal{E}_{\text{ave}}(\ell T) &= L^{-1} \langle \psi_0 | (U_F^\dagger)^\ell H_{\text{ave}} (U_F)^\ell | \psi_0 \rangle \\ &\approx L^{-1} \langle \psi_0 | (U_F^{[n]\dagger})^\ell H_{\text{ave}} (U_F^{[n]})^\ell | \psi_0 \rangle\end{aligned}$$

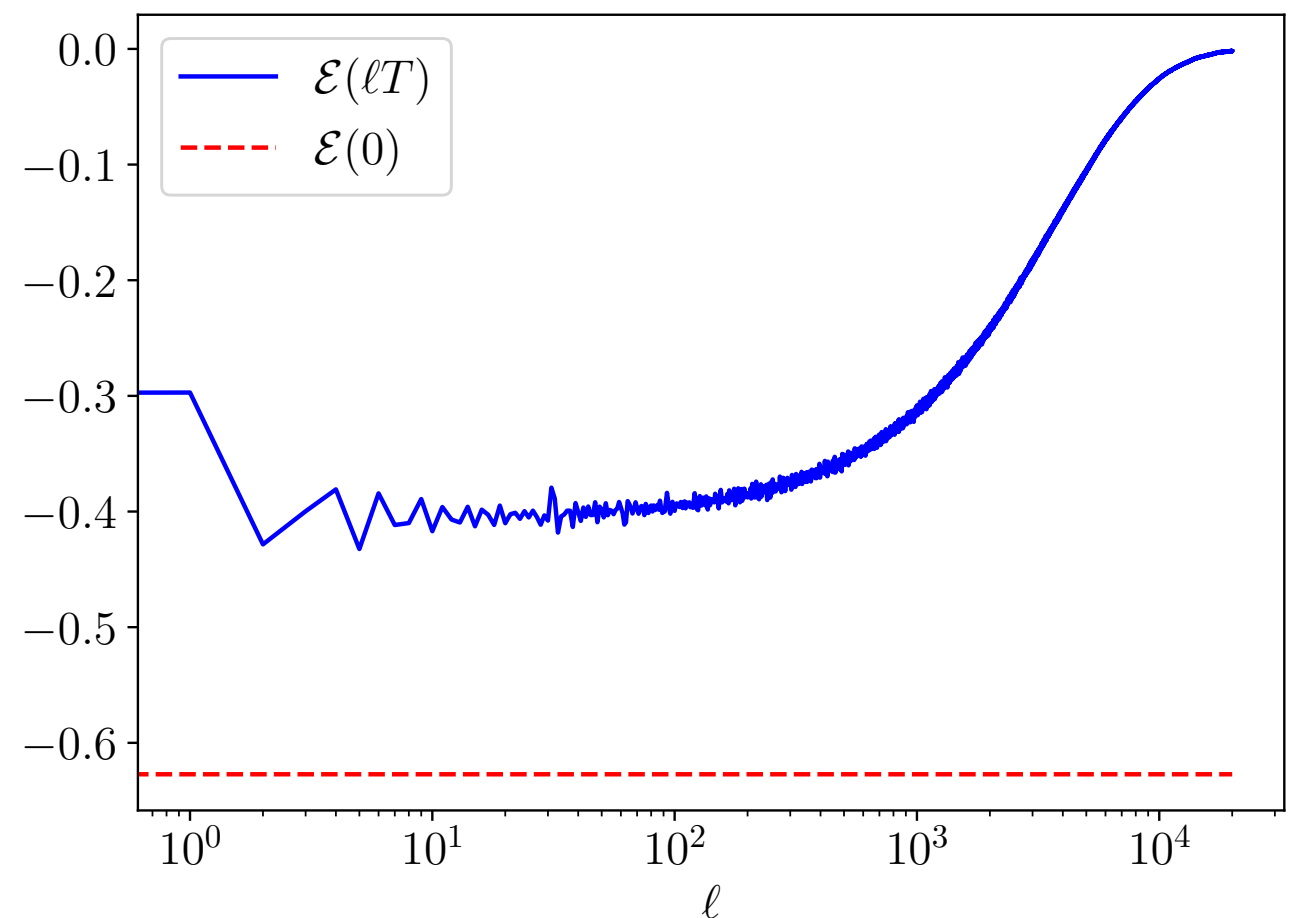


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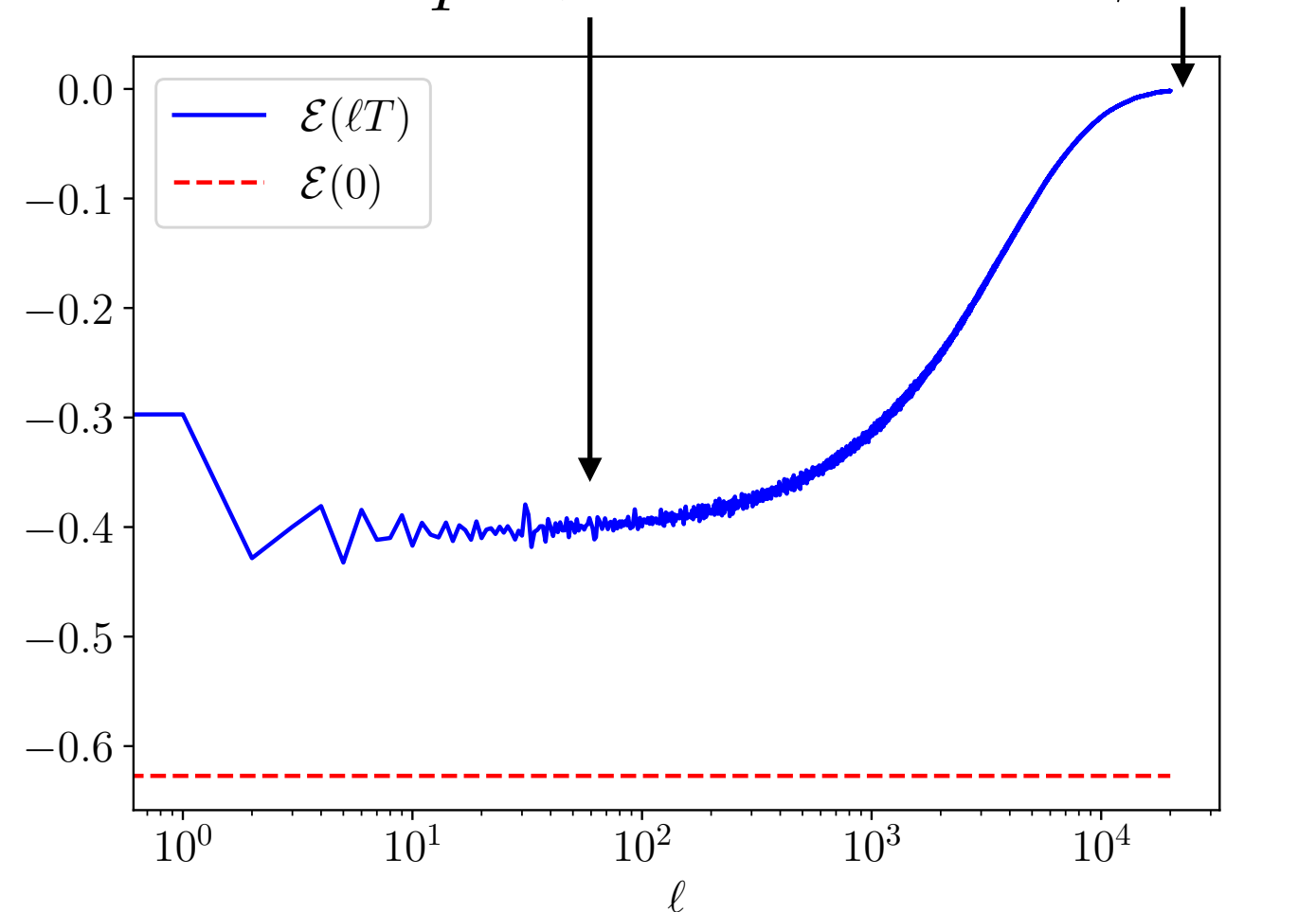
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➔ associate a temperature using ETH w.r.t. $H_F^{[n]}$: $\beta^{\text{prethermal}}$



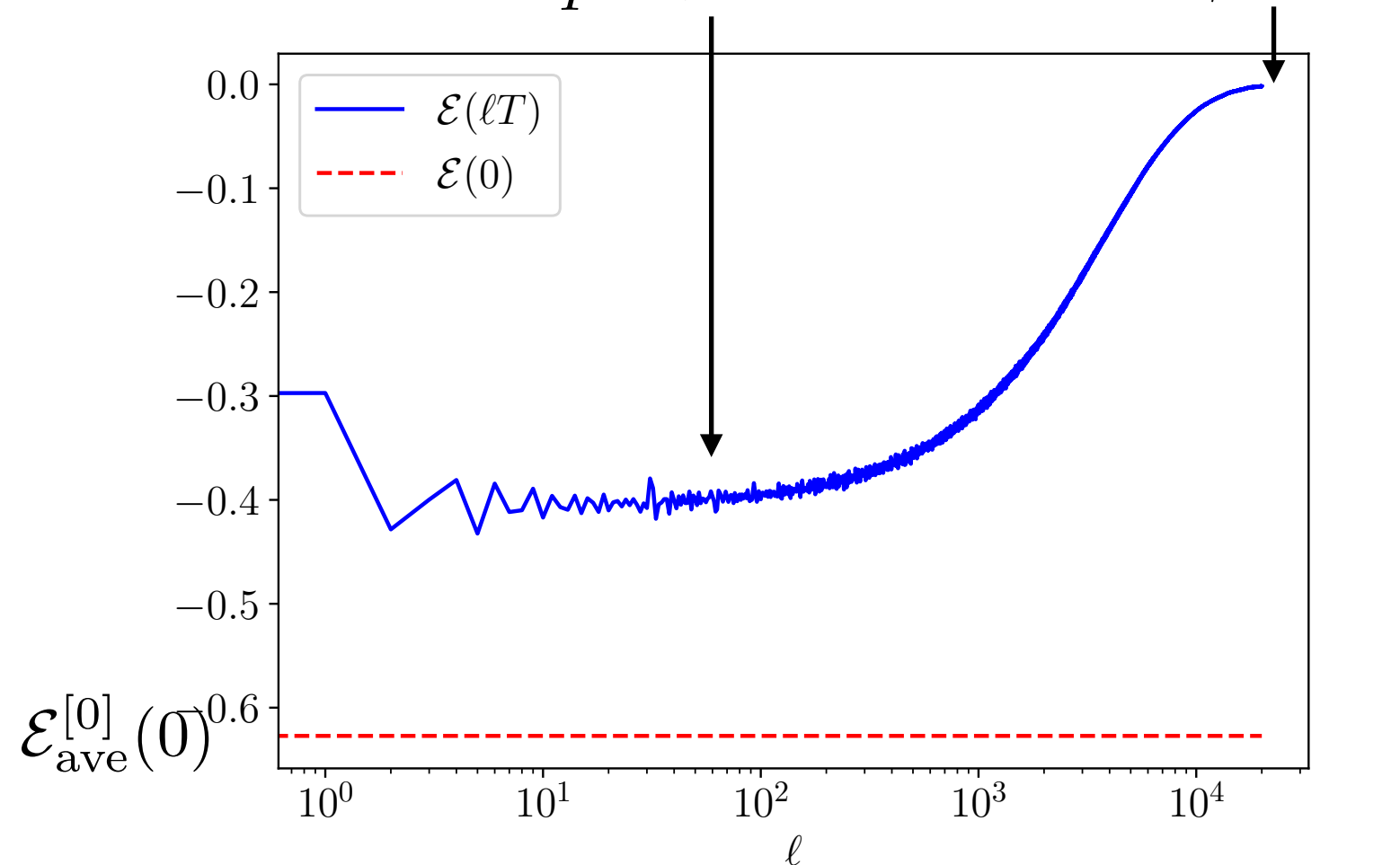
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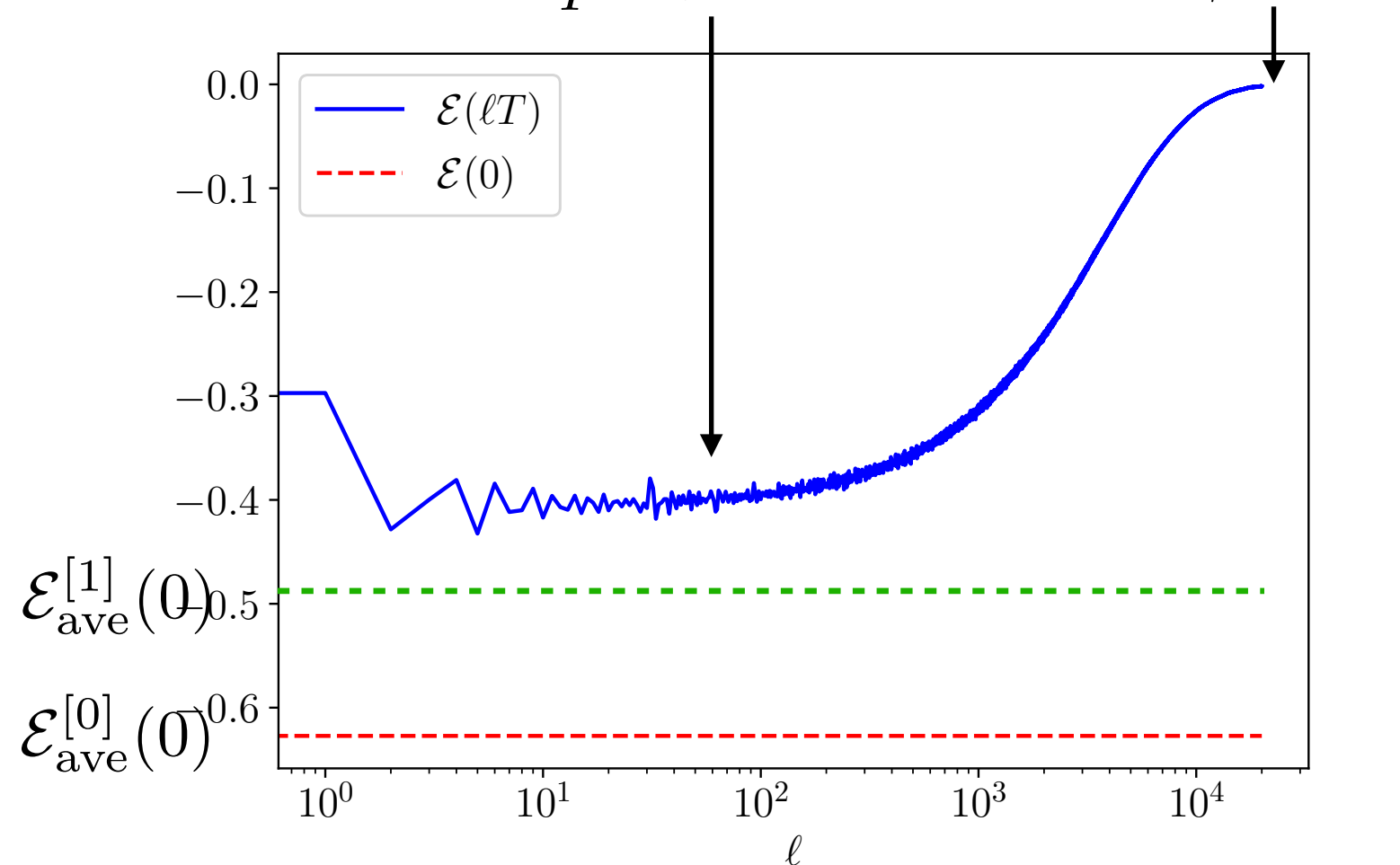
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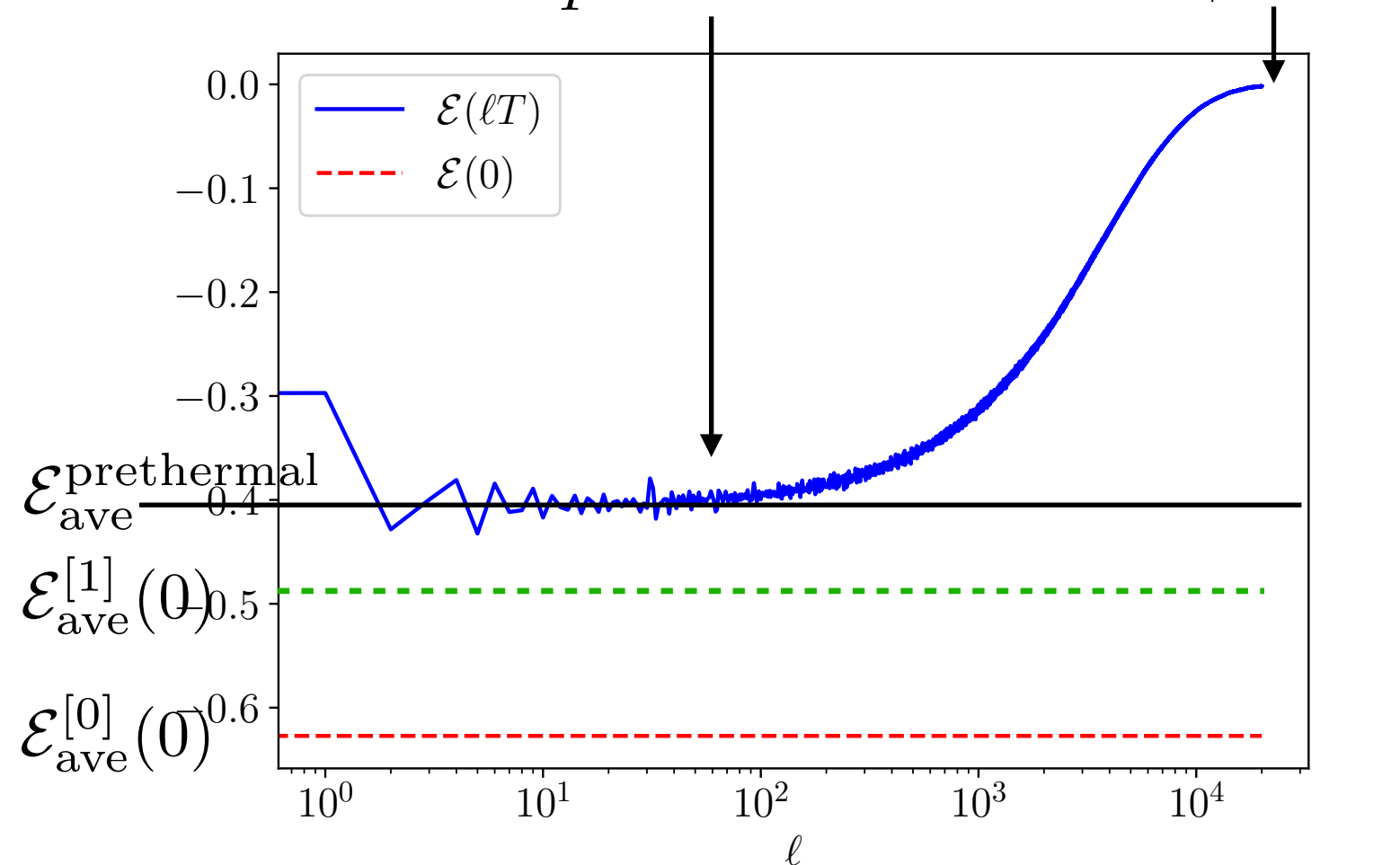
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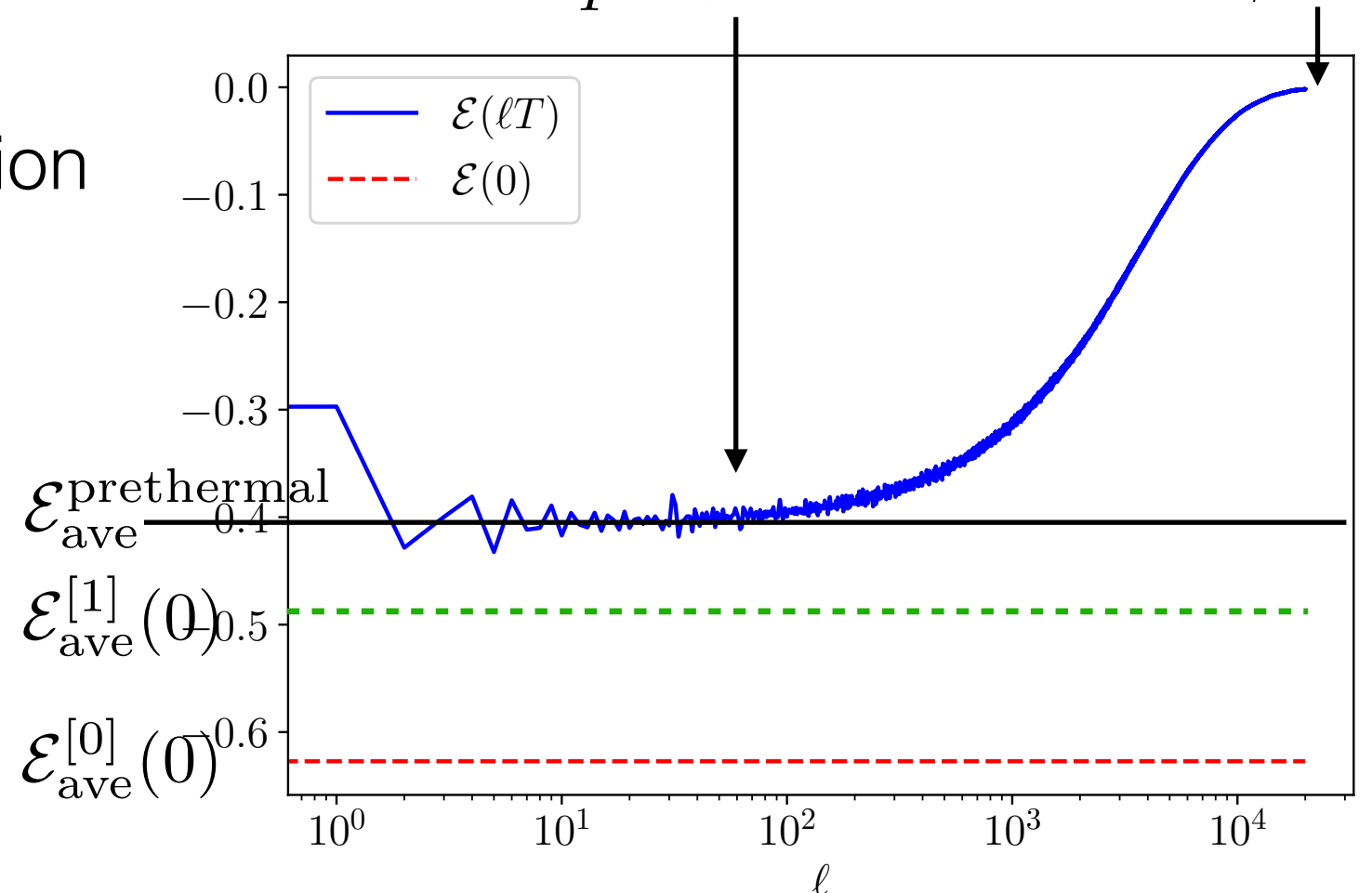
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→ four stages of thermalization
(high-frequency driving)



Understanding the pre-thermal plateau

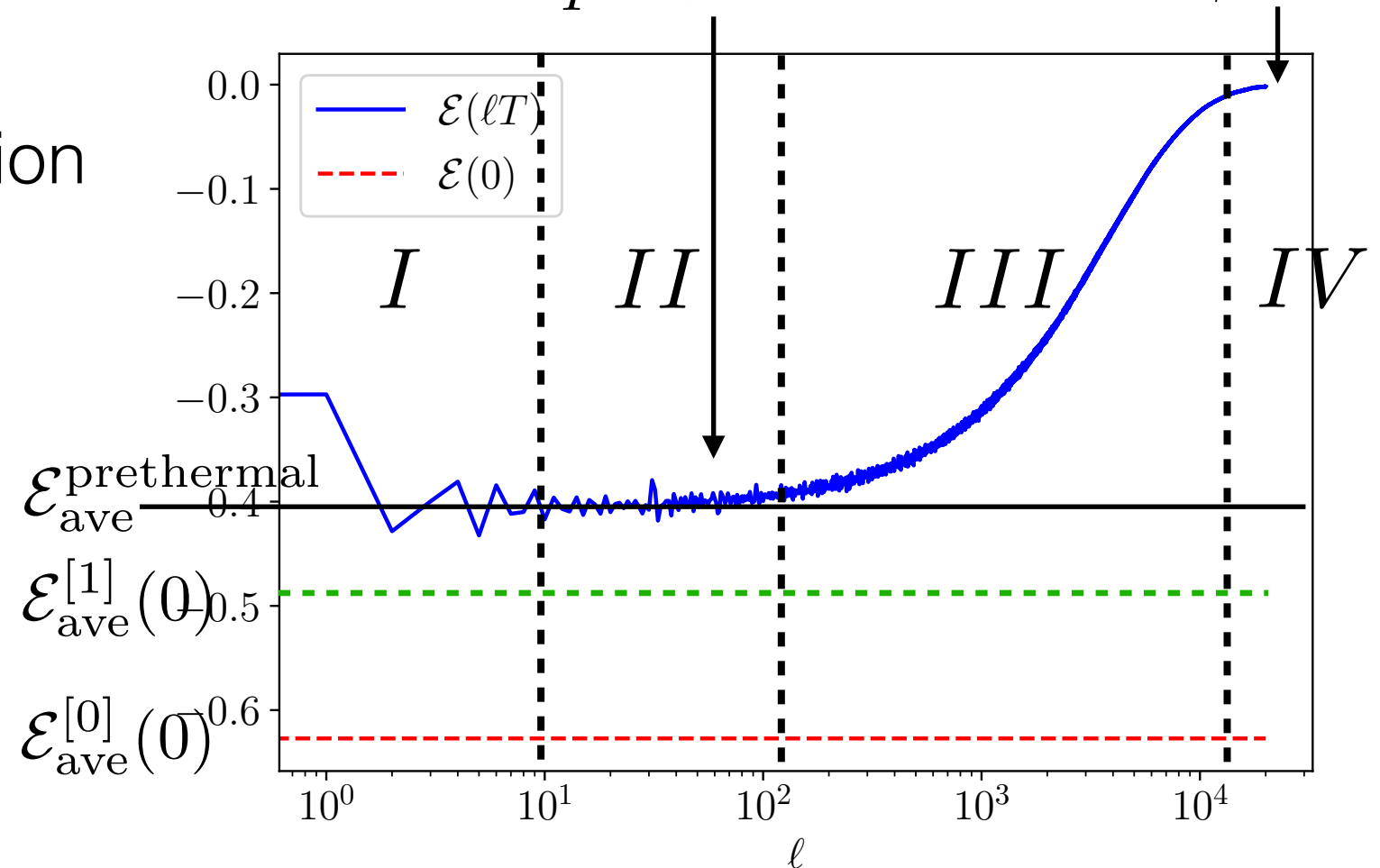
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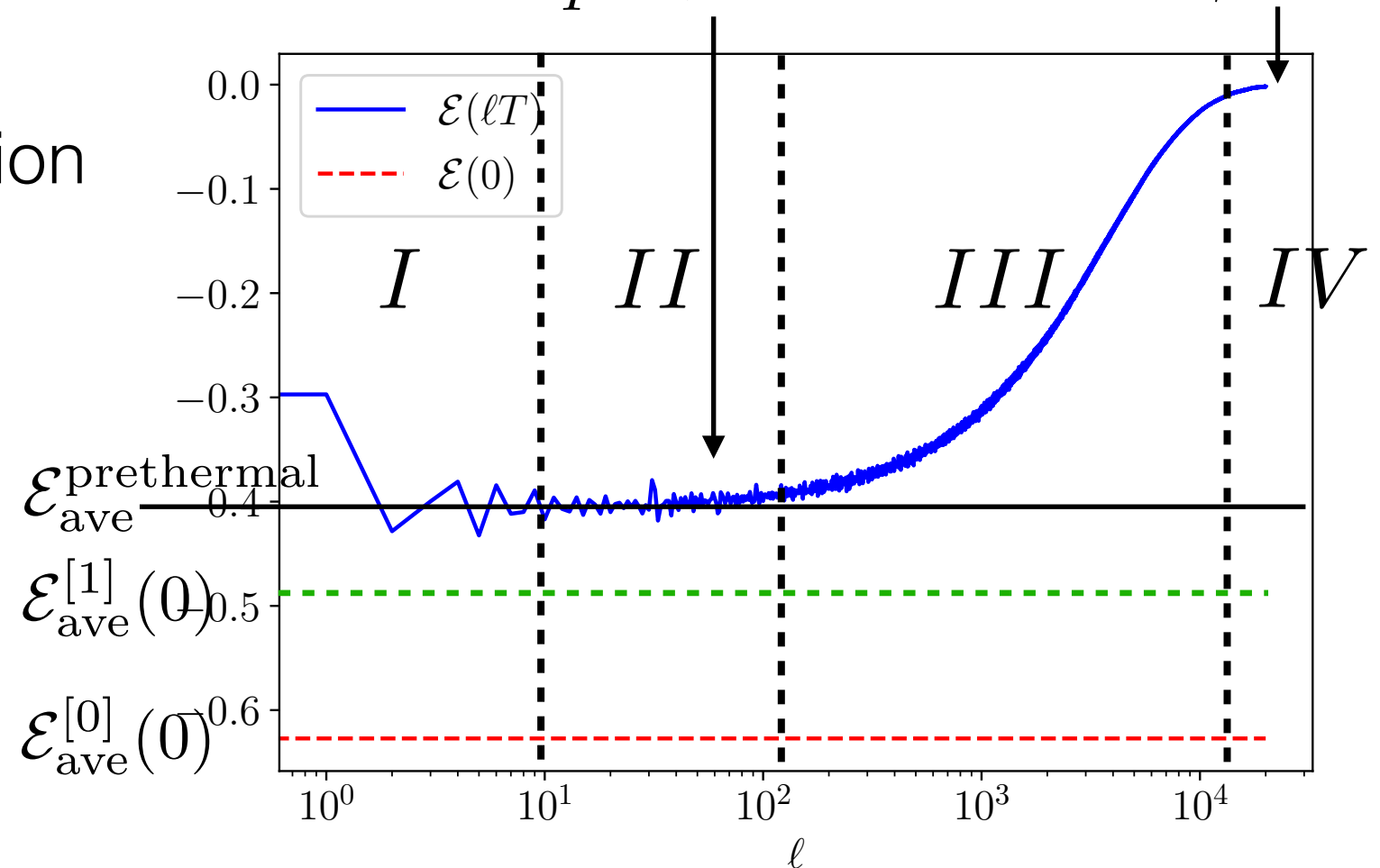
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I+II: quench to $H_F^{[n]}$

I+...+IV: quench to H_F



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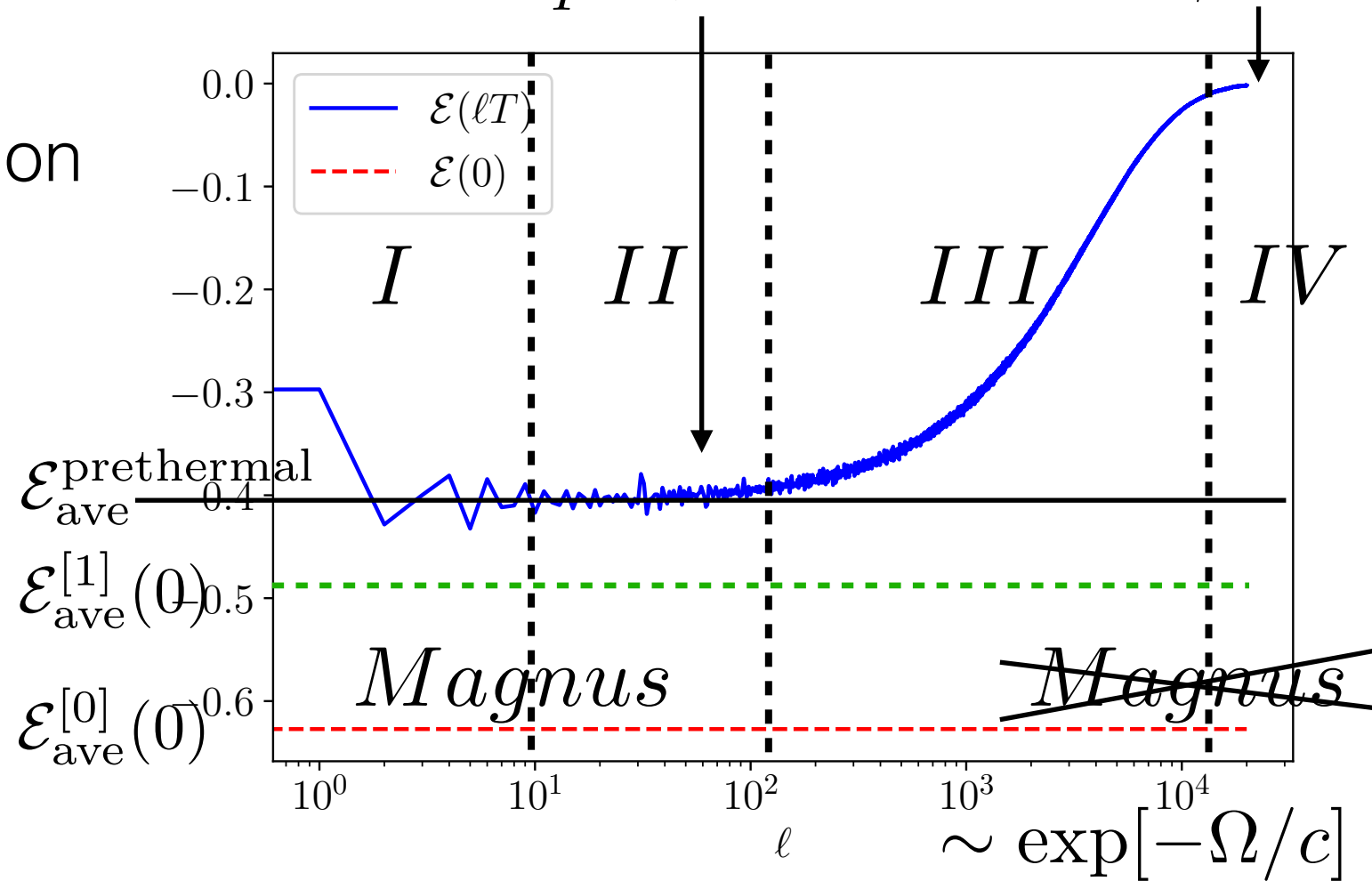
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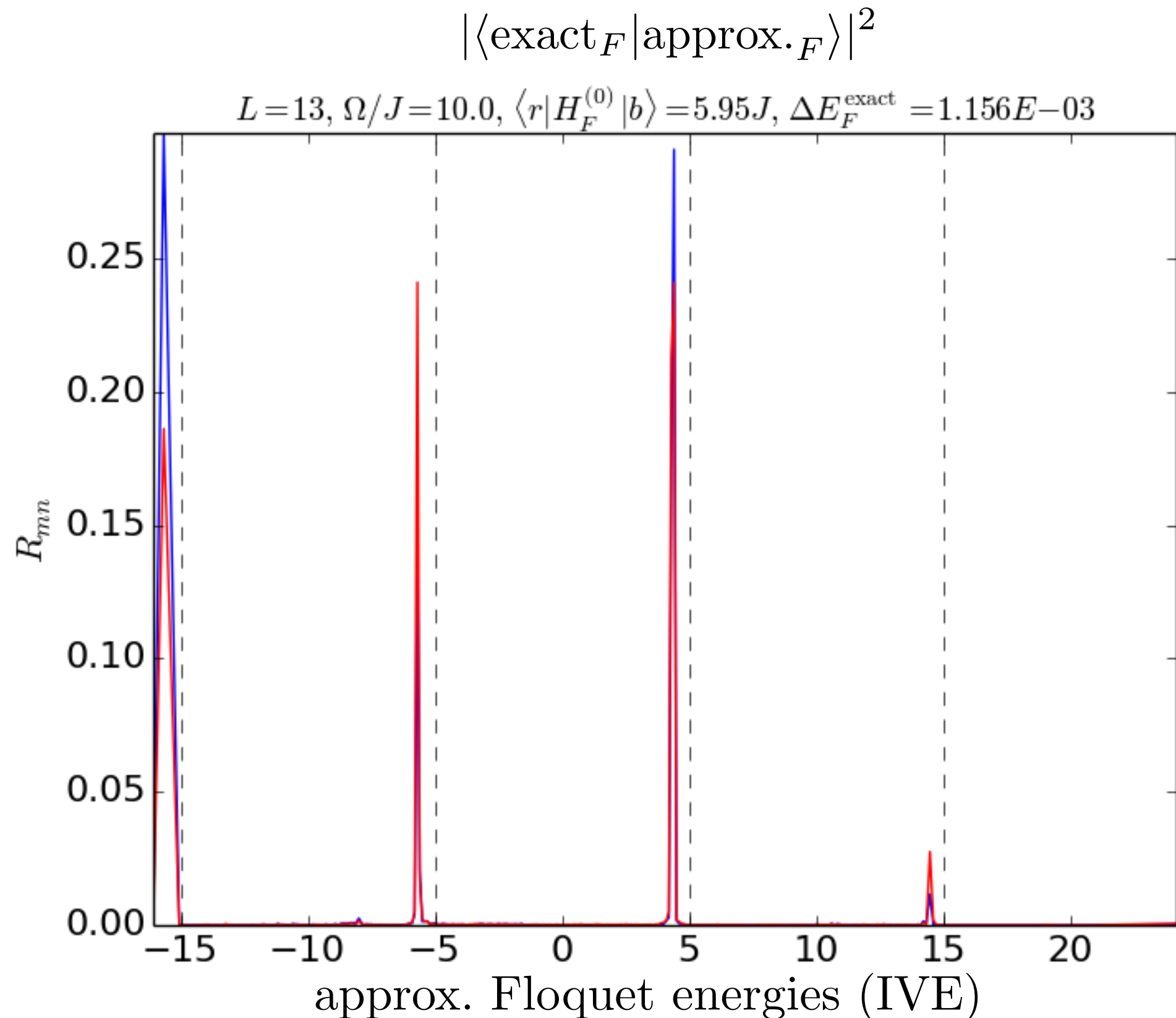
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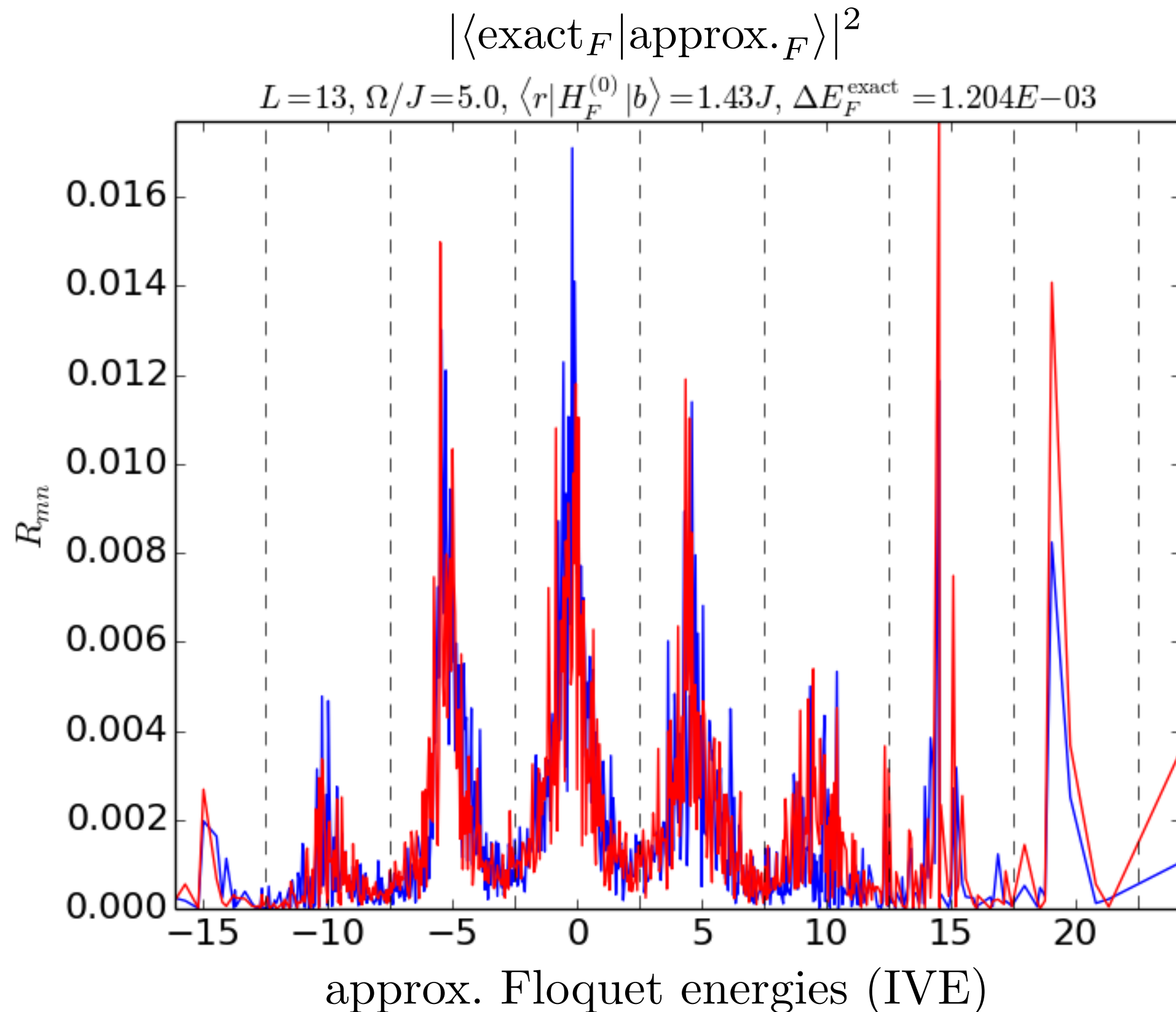
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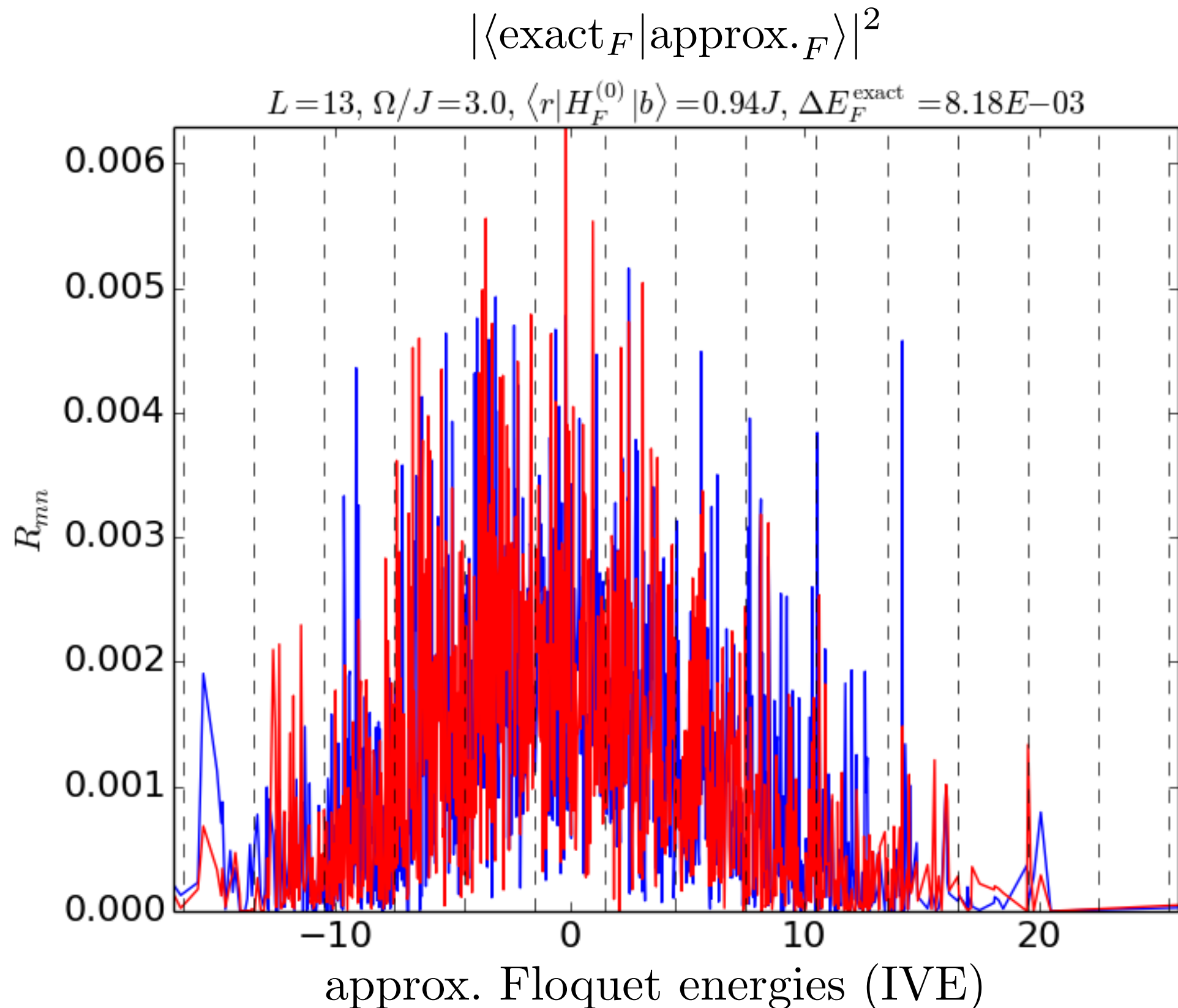
Stage III: the onset of heating: many-body resonances (**not** captured by inverse-frequency expansion)



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➔ only 100% safe method to study thermalization is (still) ED

but see also [Ye et al., arXiv:1902.01859](#)

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➔ finite system size limits the number of resonance peaks

➔ discrete spectrum: finite-size gaps close to GS

➔ appropriate order of limits: $\lim_{\Omega \rightarrow \infty} \lim_{L \rightarrow \infty}$ hard to achieve

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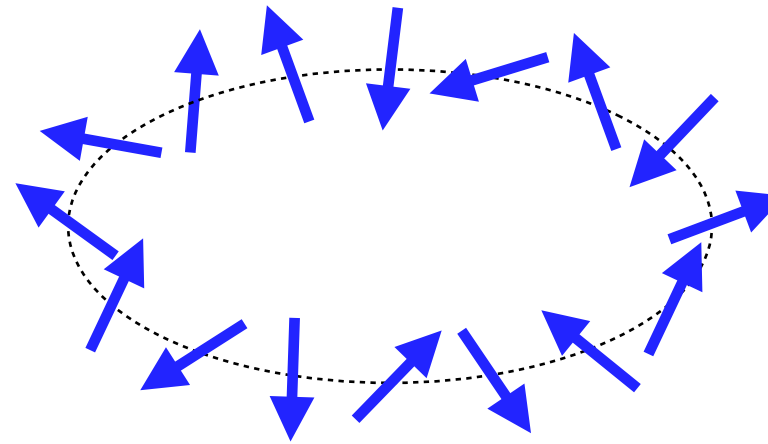
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can we avoid these issues?

which feature(s) can we sacrifice?

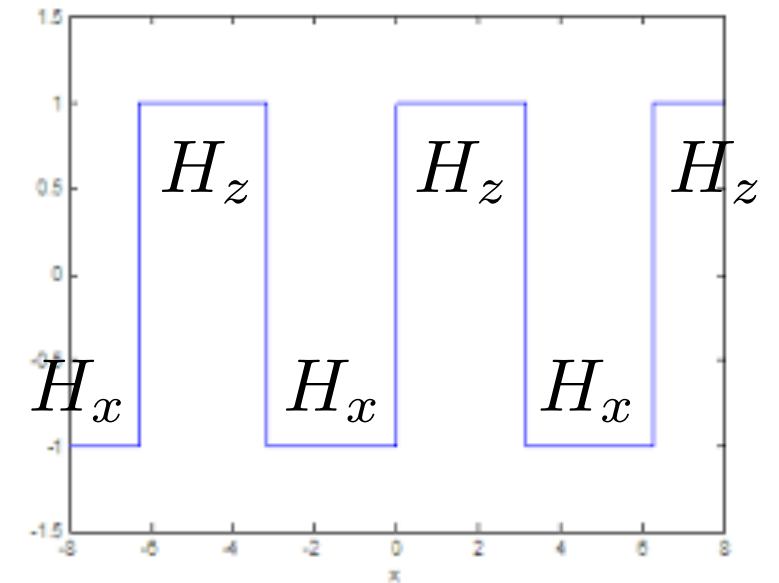
Classical Floquet systems

→ driven coupled rotors



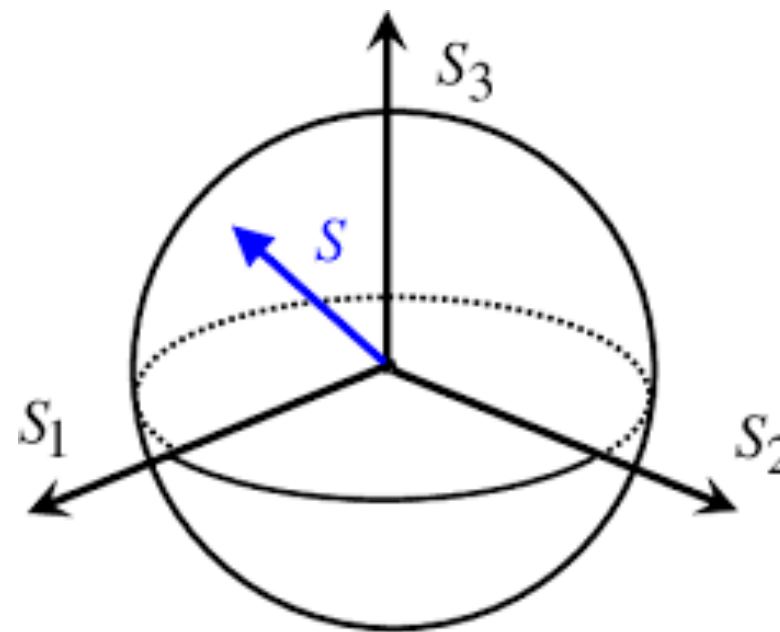
Owen Howell

$$H(t) = \begin{cases} \sum_{j=1}^L J S_j^z S_{j+1}^z + h S_j^z & \text{for } t \in [0, T/2] \bmod T \\ \sum_{j=1}^L g S_j^x & \text{for } t \in [T/2, T] \bmod T \end{cases}$$



$$\{S_i^\alpha, S_j^\beta\} = \delta_{ij} \epsilon^{\alpha\beta\gamma} S_j^\gamma$$

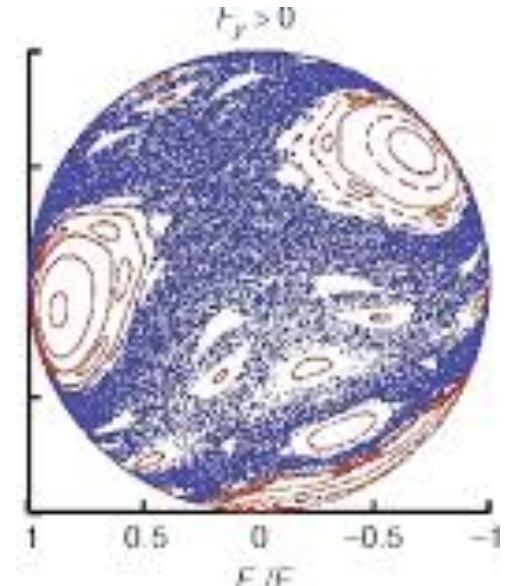
$$|\vec{S}| = 1$$



A crucial feature of the step drive

→ want *exponentially* long times in a **chaotic** system!

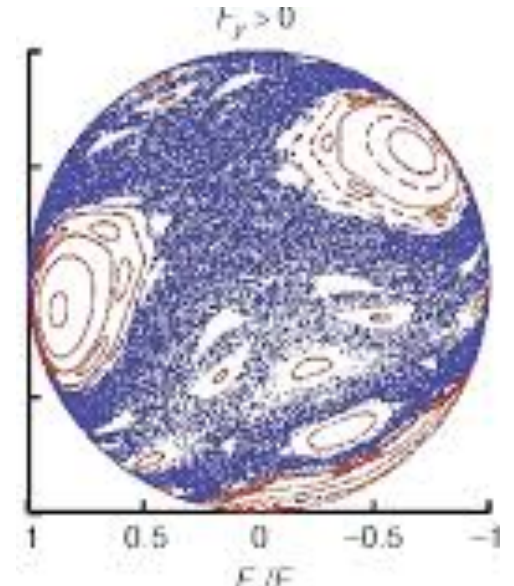
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→ solve Hamilton's EOM **exactly**: $\vec{S}_j(\ell T) = [\tau_2 \circ \tau_1]^\ell (\vec{S}_j(0))$

$$\tau_1(\vec{S}_j) = \begin{bmatrix} S_j^x \cos(\kappa_j T/2) - S_j^y \sin(\kappa_j T/2) \\ S_j^x \sin(\kappa_j T/2) + S_j^y \cos(\kappa_j T/2) \\ S_j^z \end{bmatrix} \quad \tau_2(\vec{S}_j) = \begin{bmatrix} S_j^x \\ S_j^y \cos(gT/2) - S_j^z \sin(gT/2) \\ S_j^y \sin(gT/2) + S_j^z \cos(gT/2) \end{bmatrix}$$

$$\kappa_j = J(S_{j-1}^z + S_{j+1}^z) + h$$

Ensemble pre-thermalization

→ time-averaged energy

$$H_{\text{ave}} = \frac{1}{2} \sum_{j=1}^N J S_j^z S_{j+1}^z + h S_j^z + g S_j^x$$

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→ measure energy absorption and its fluctuations

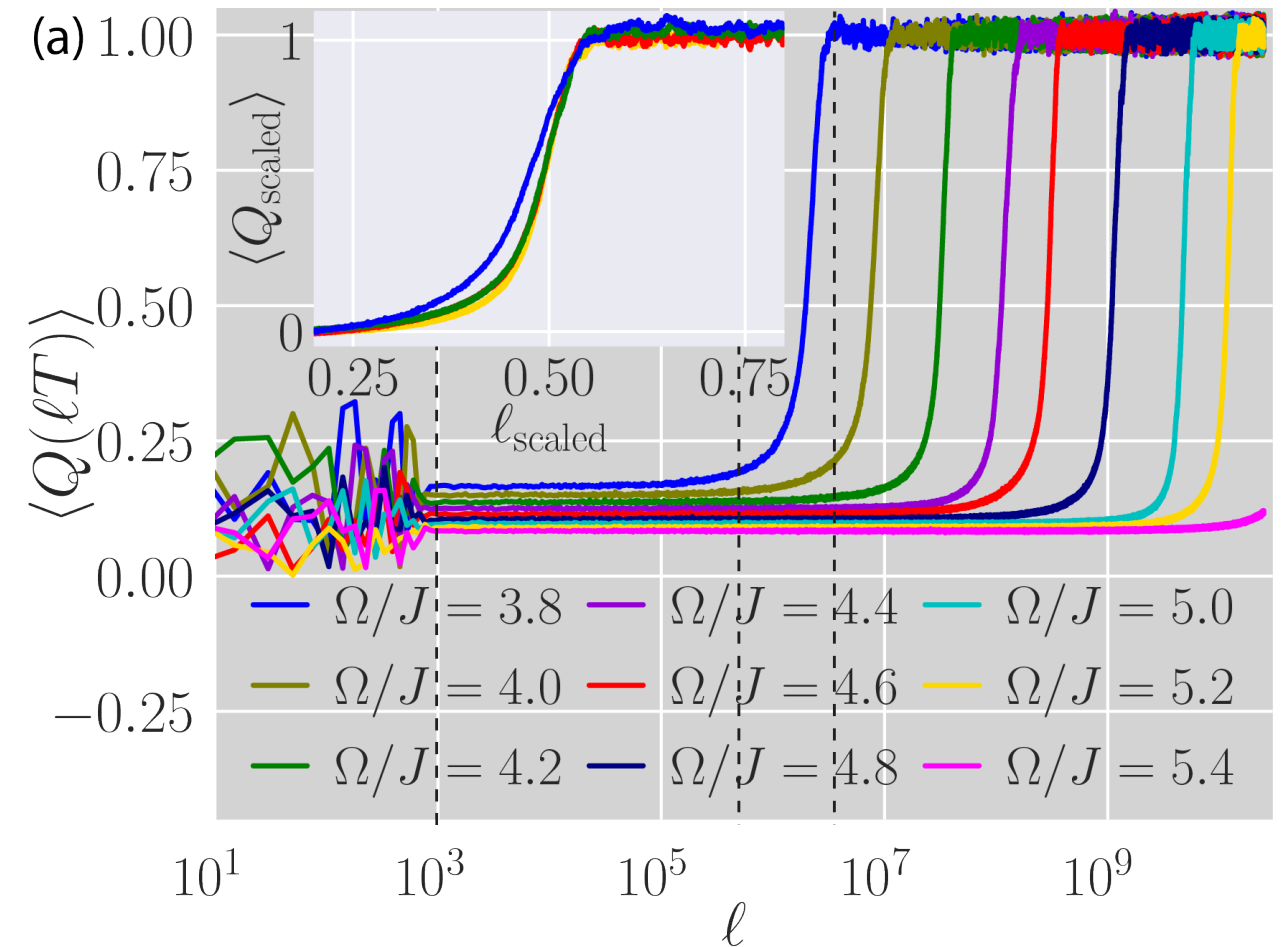
$$\langle Q(\ell T) \rangle = \frac{\langle H_{\text{ave}}[\{\vec{S}_j(\ell T)\}] \rangle - E_{\text{GS}}}{\langle H_{\text{ave}} \rangle_{\beta=0} - E_{\text{GS}}} \in [0, 1],$$

$$\langle \delta Q(\ell T) \rangle = \sqrt{\frac{\langle H_{\text{ave}}^2[\{\vec{S}_j(\ell T)\}] \rangle - \langle H_{\text{ave}}[\{\vec{S}_j(\ell T)\}] \rangle^2}{\langle H_{\text{ave}}^2 \rangle_{\beta=0} - \langle H_{\text{ave}} \rangle_{\beta=0}^2}}$$

Classical Floquet pre-thermalization

$L = 100$

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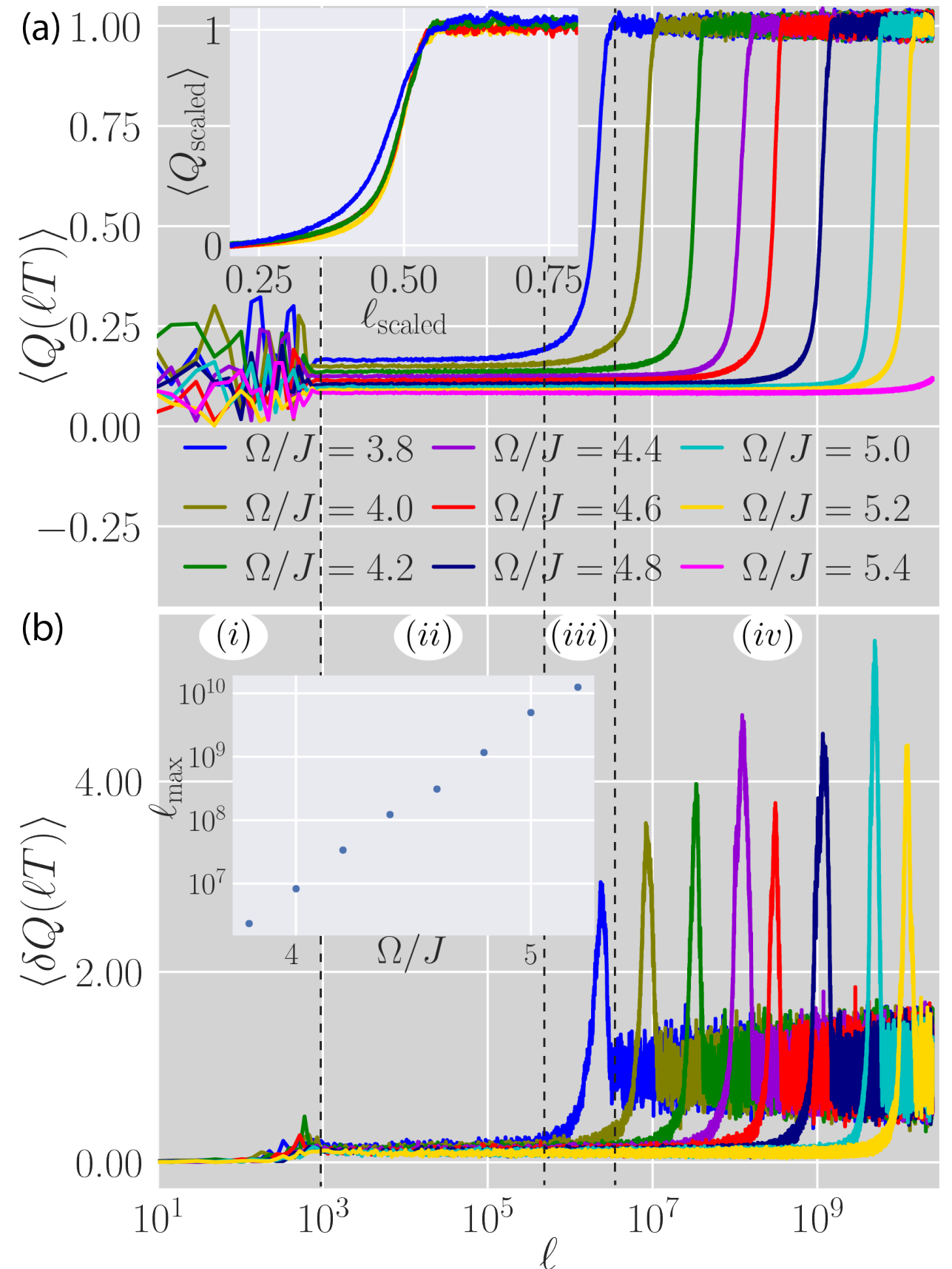


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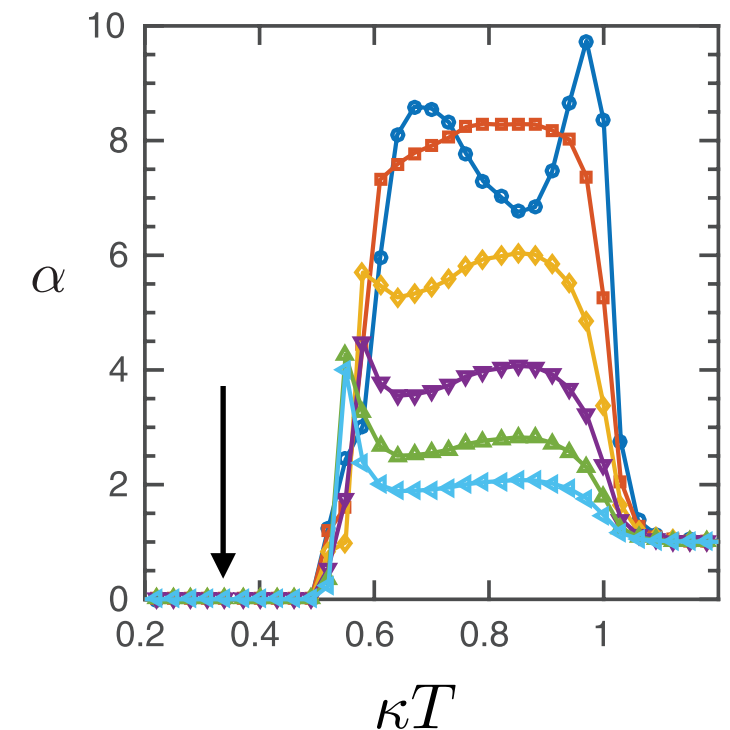
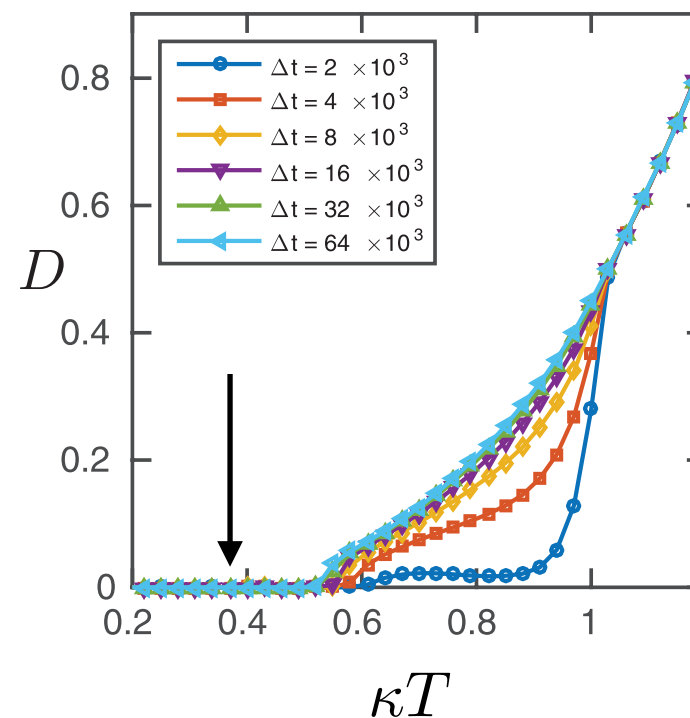
Related phenomenon: subdiffusion

→ coupled kicked pendula:

$$H(t) = \sum_{j=1}^L \left[\frac{p_j^2}{2} - \kappa \cos(\phi_j - \phi_{j+1}) \sum_{m=-\infty}^{+\infty} \delta(t - mT) \right]$$

$$D = \frac{\langle p^2(t_f) \rangle - \langle p^2(t_i) \rangle}{t_f - t_i} \propto t^{\alpha-1}$$

diffusion coefficient



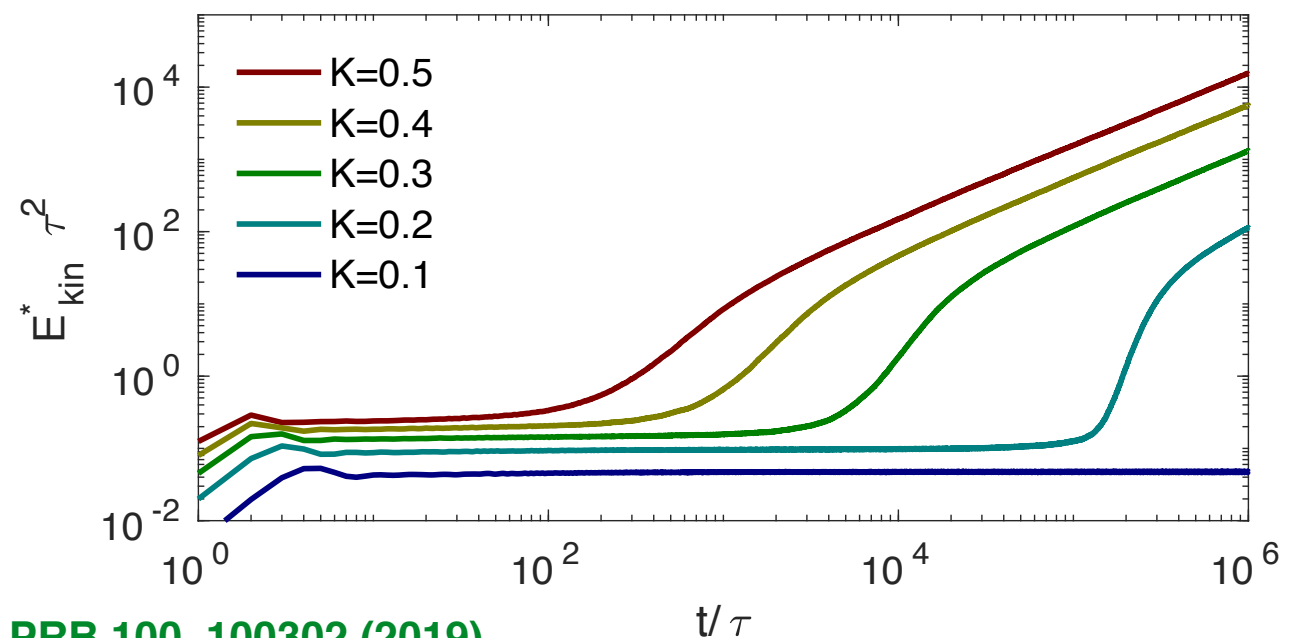
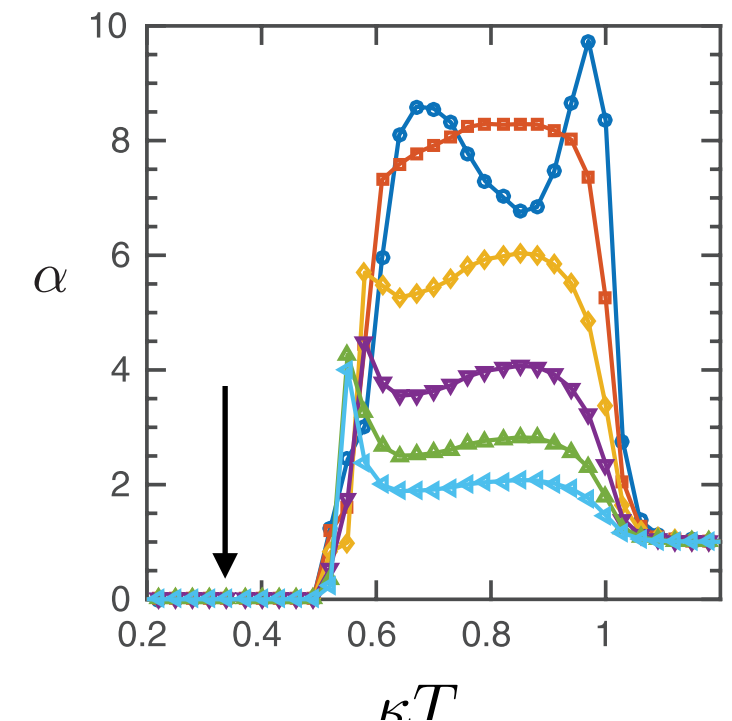
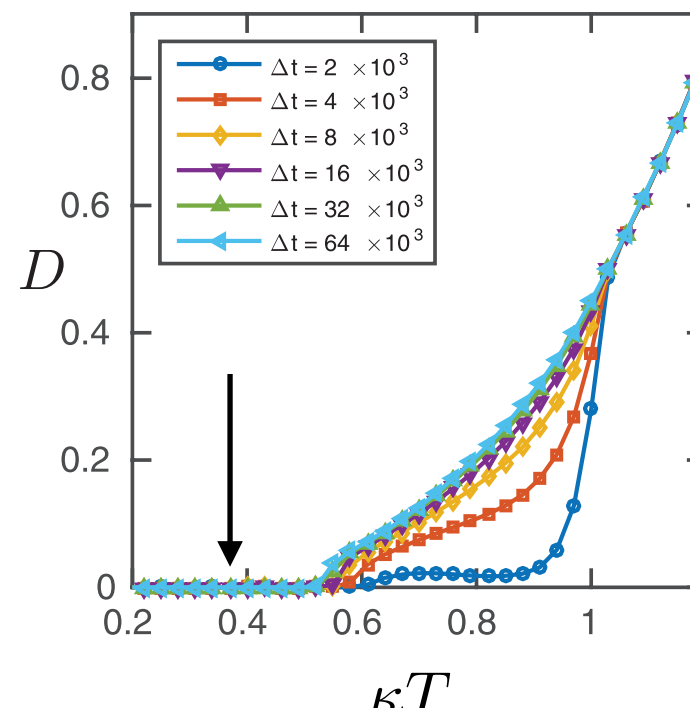
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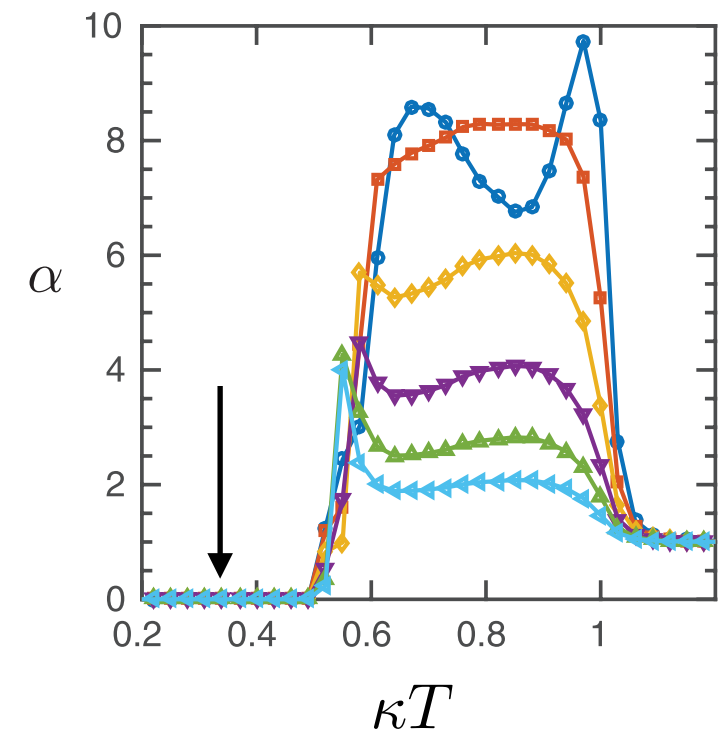
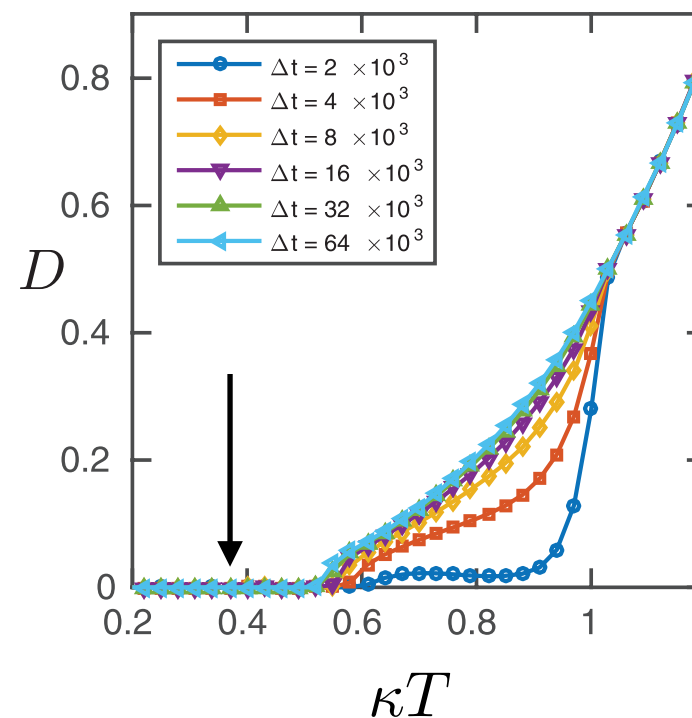
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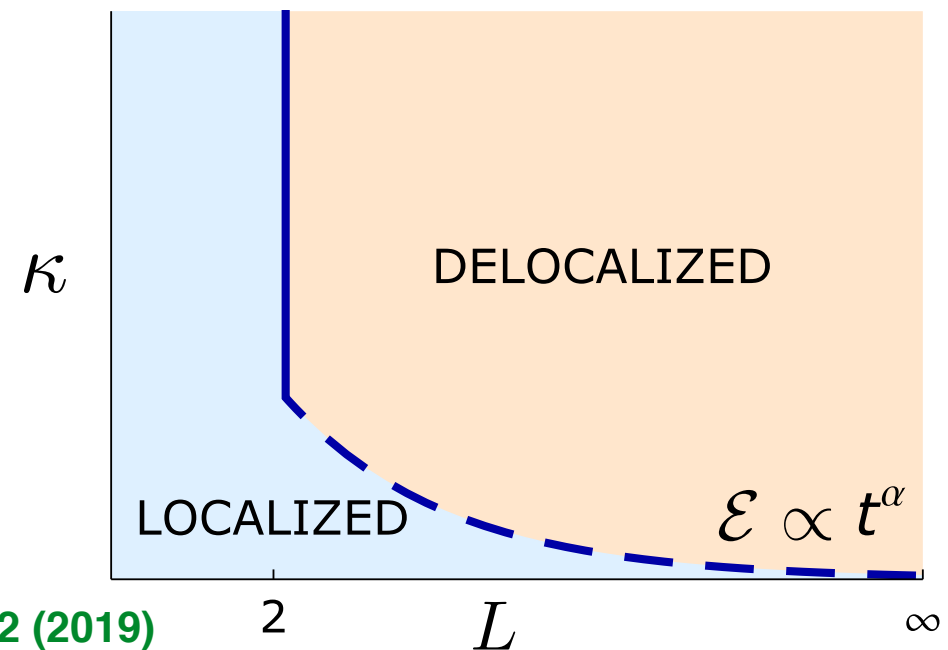
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→ schematic stability diagram



Can we use the inverse-frequency expansion?

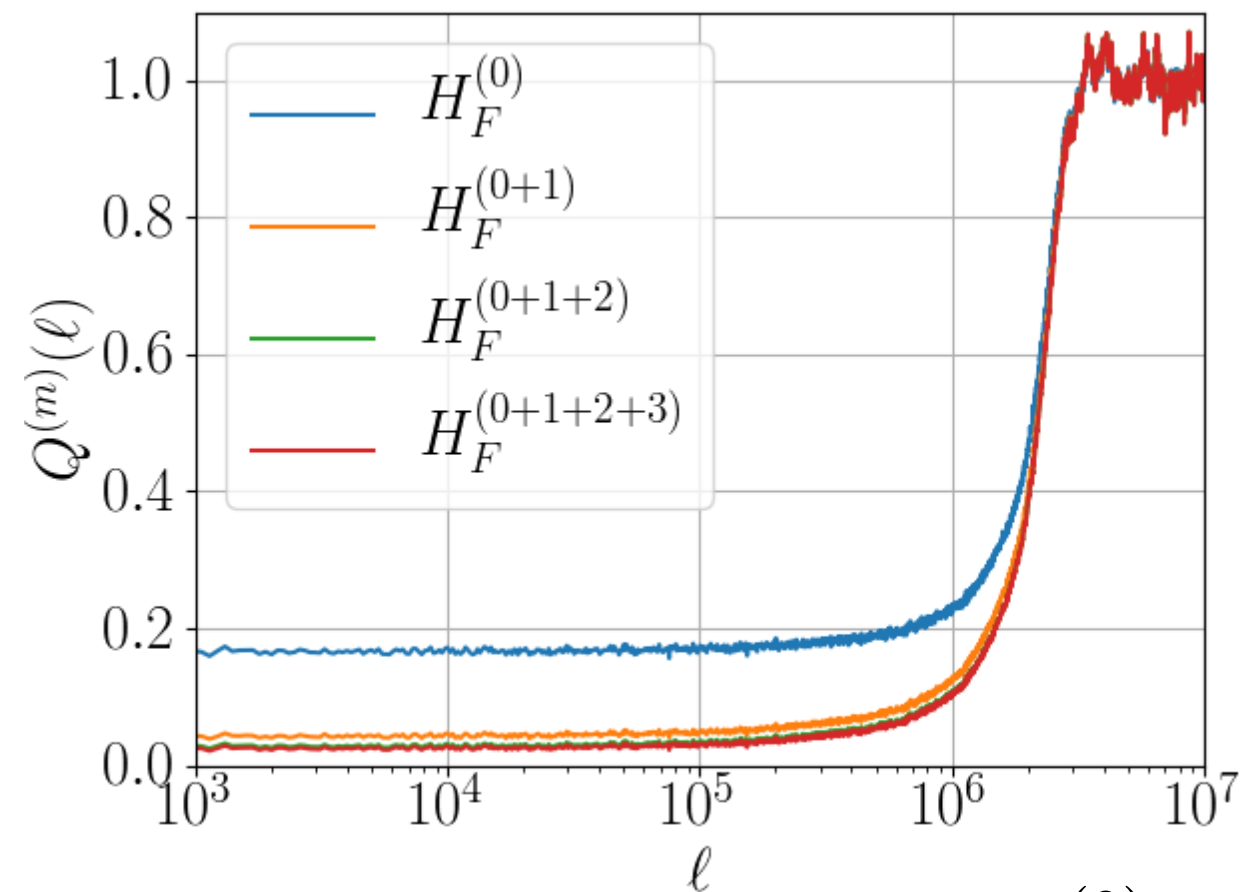
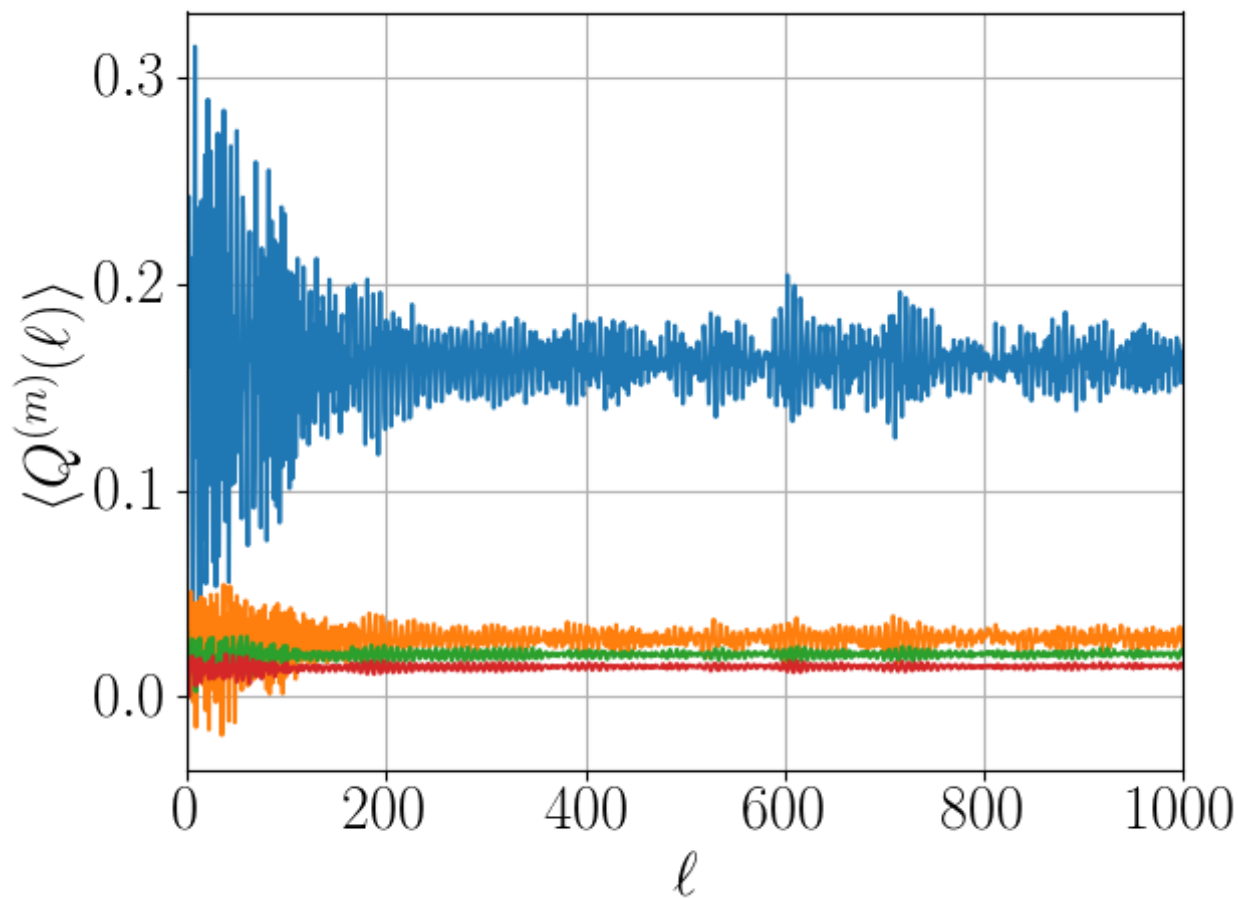
→ apply classical Magnus expansion: $-i[\cdot, \cdot] \longrightarrow \{\cdot, \cdot\}$

$$H_F^{(0)} = T^{-1} \int_0^T dt H(t)$$

$$H_F^{(1)} \sim \int_0^T dt_1 \int_0^{t_1} dt_2 [H(t_1), H(t_2)]$$

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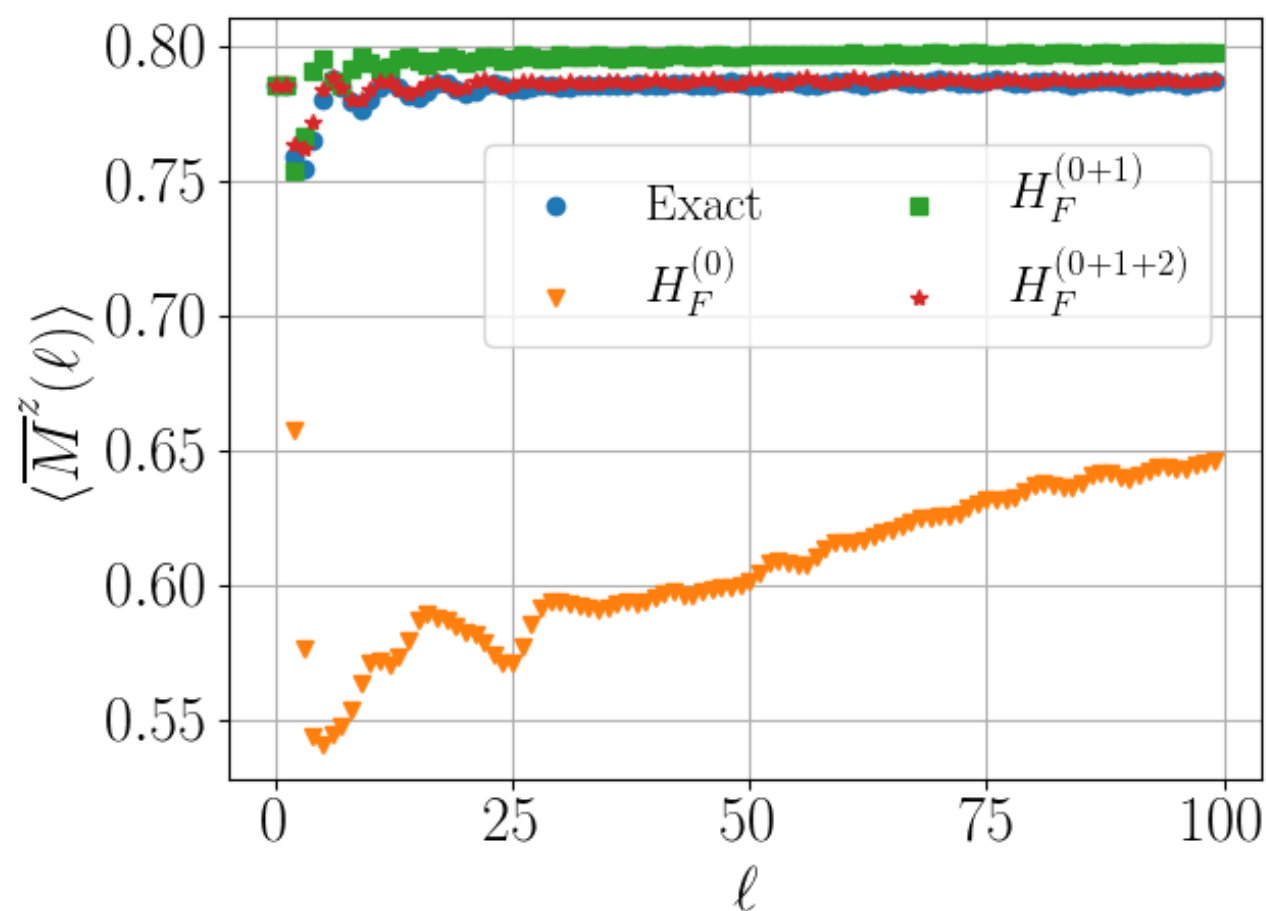
initial state : GS of $H_F^{(0)}$

$$\langle Q^{[n]}(\ell T) \rangle = \frac{\langle H_F^{[n]}[\{\vec{S}_j(\ell T)\}] \rangle - E_{\text{GS}}}{\langle H_F^{[n]} \rangle_{\beta=0} - E_{\text{GS}}} \in [0, 1],$$

Ensemble-averaged observables captured by Magnus expansion

→ time-averaged magnetization

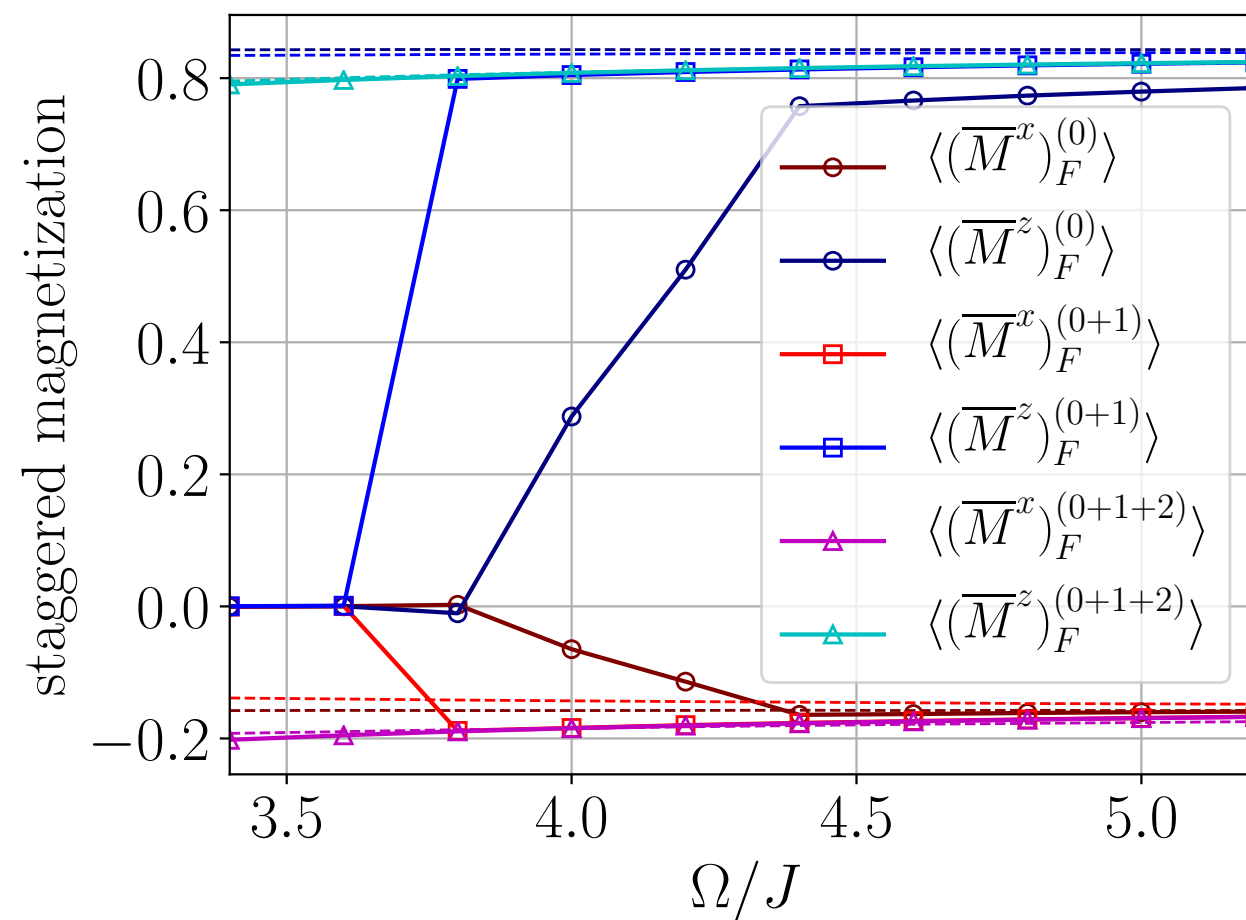
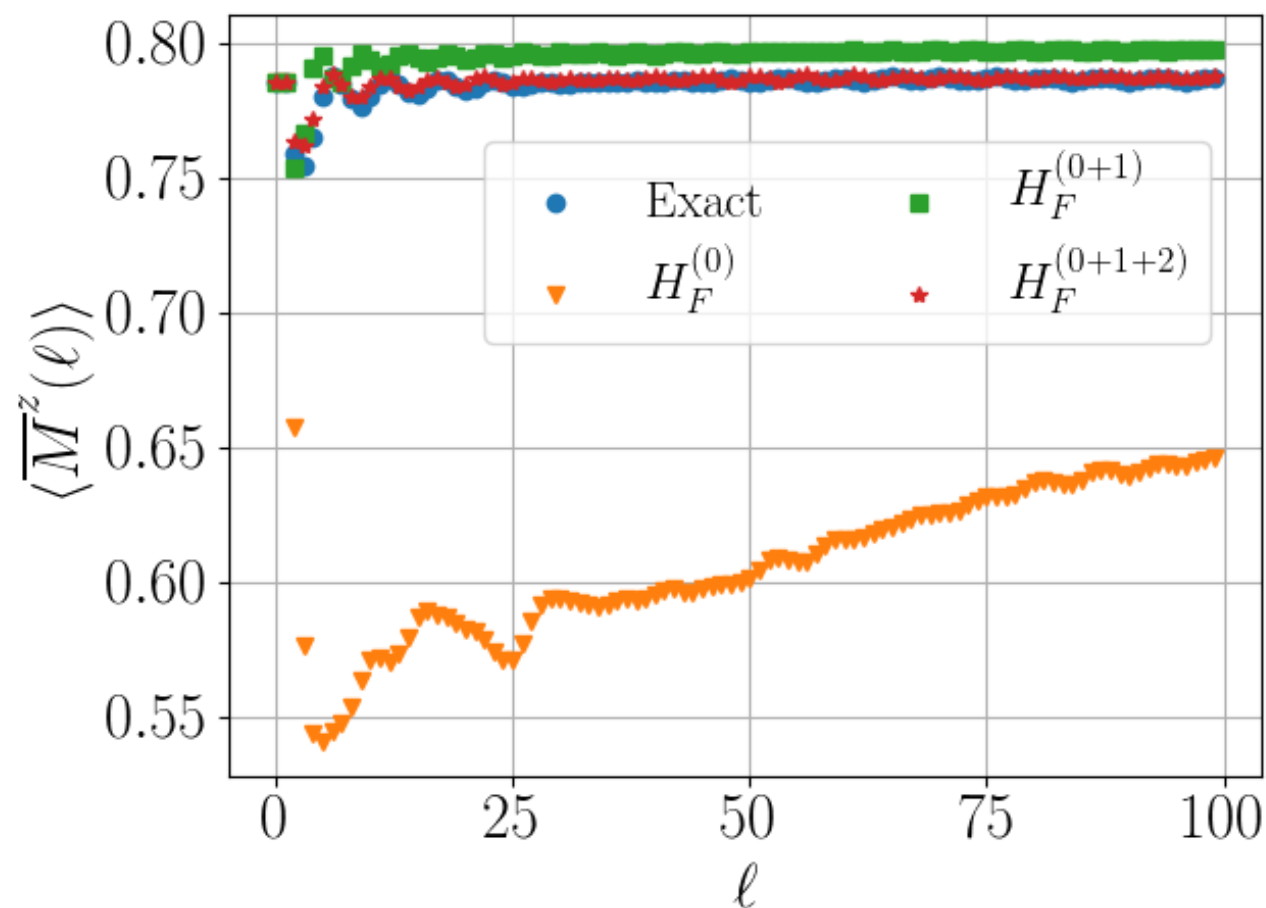
$$M^\alpha(\ell T) = \frac{1}{N} \sum_{j=1}^N (-1)^j S_j^\alpha(\ell T); \overline{M}^\alpha = \frac{1}{N_T} \sum_{\ell=N_T}^{N_T+10^3} M^\alpha(\ell T)$$



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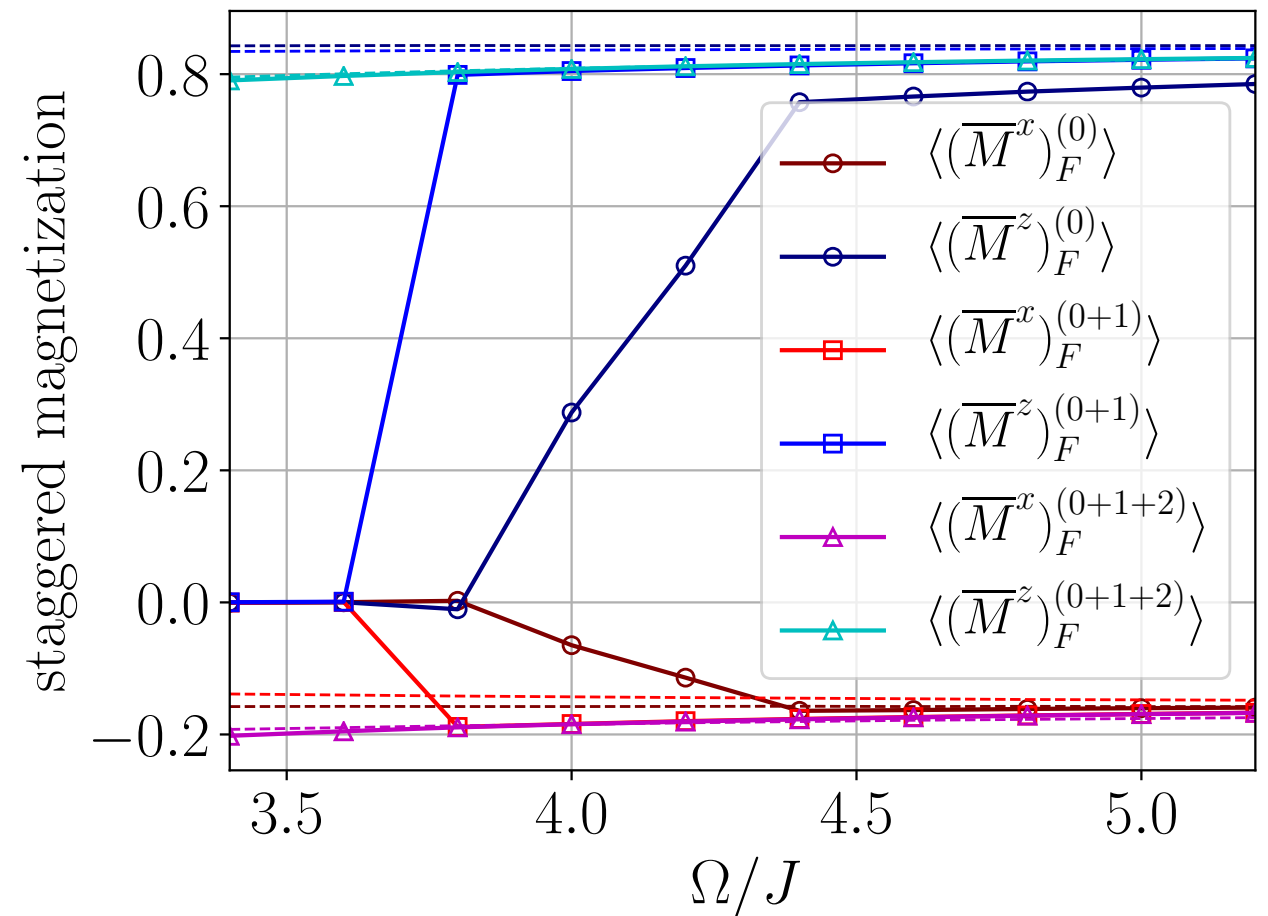
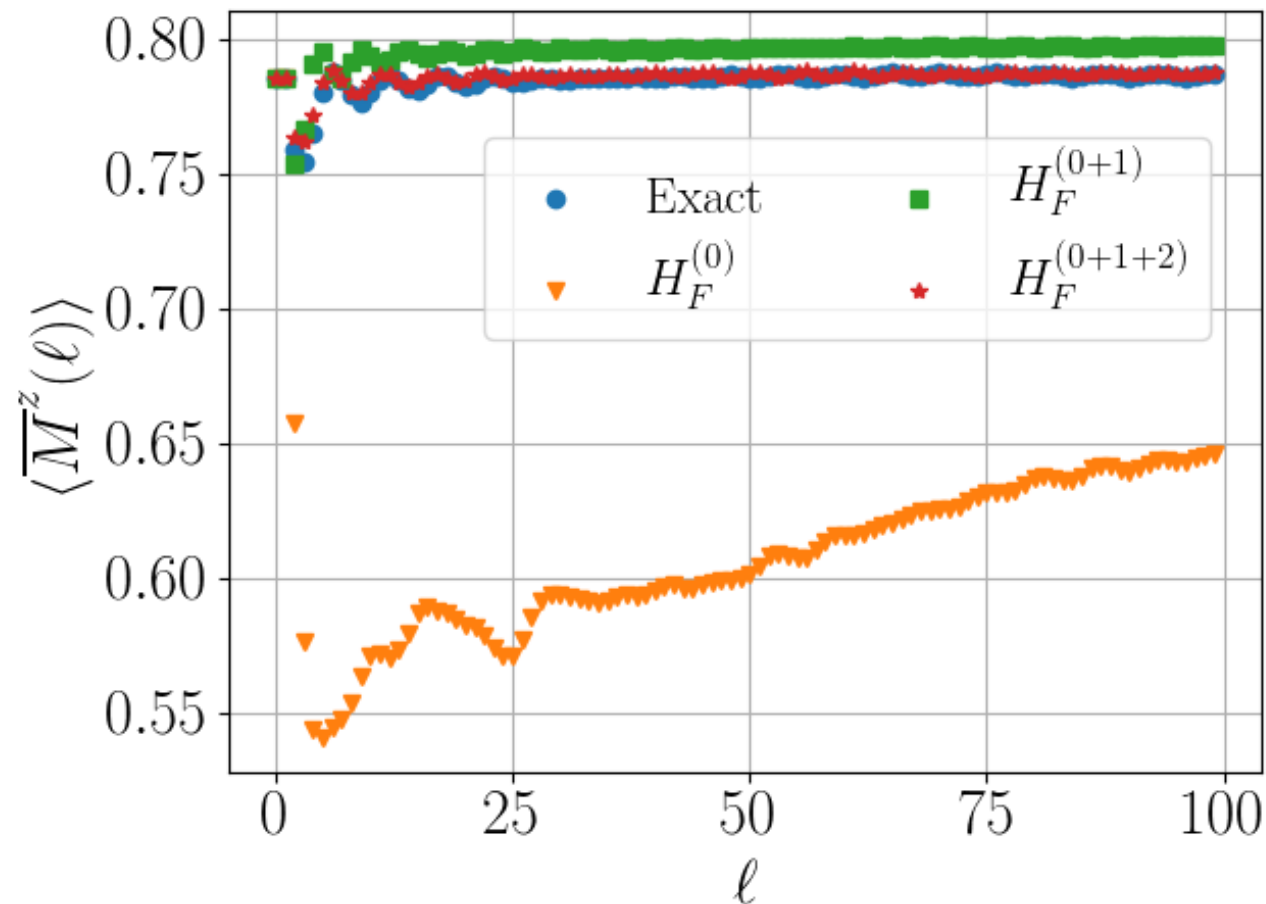
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→ classical Floquet engineering

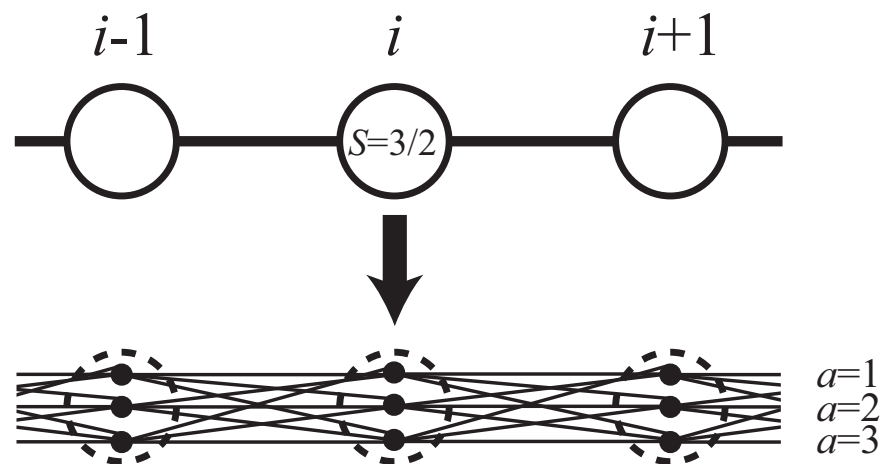
Higashikawa, Fujita, Sato, arXiv:1810.01103 (2018)

But why does Magnus work?

classical equations of motion are **nonlinear!**

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- consider proof for quantum systems:
 - holds for spin-1/2 particles
 - take large- S limit and decompose into many spin-1/2's



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$$\dot{\rho}(t) = \{\rho(t), H(t)\}$$
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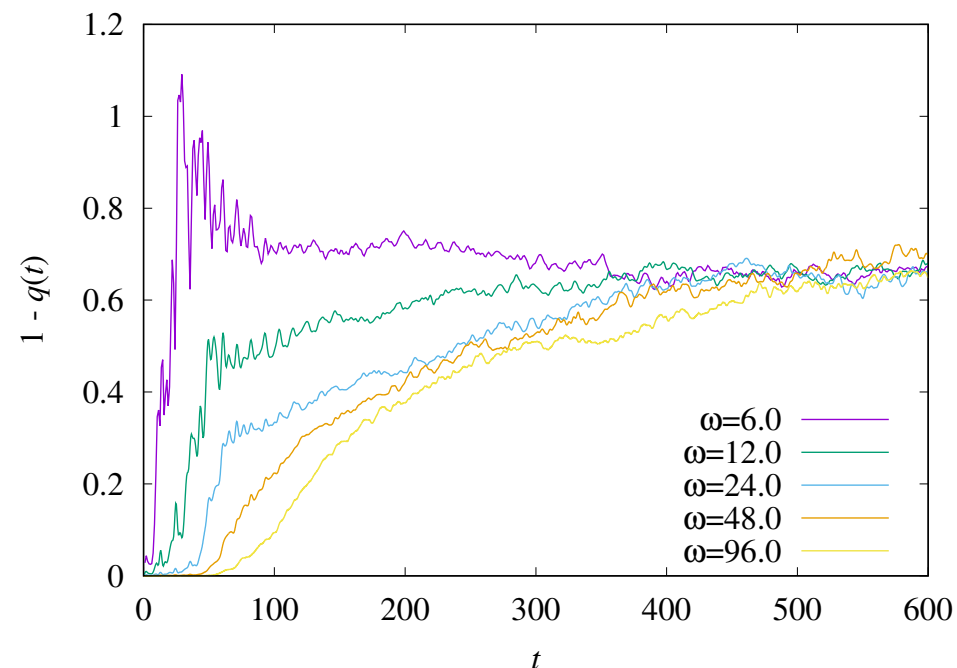
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- important **caveat**:

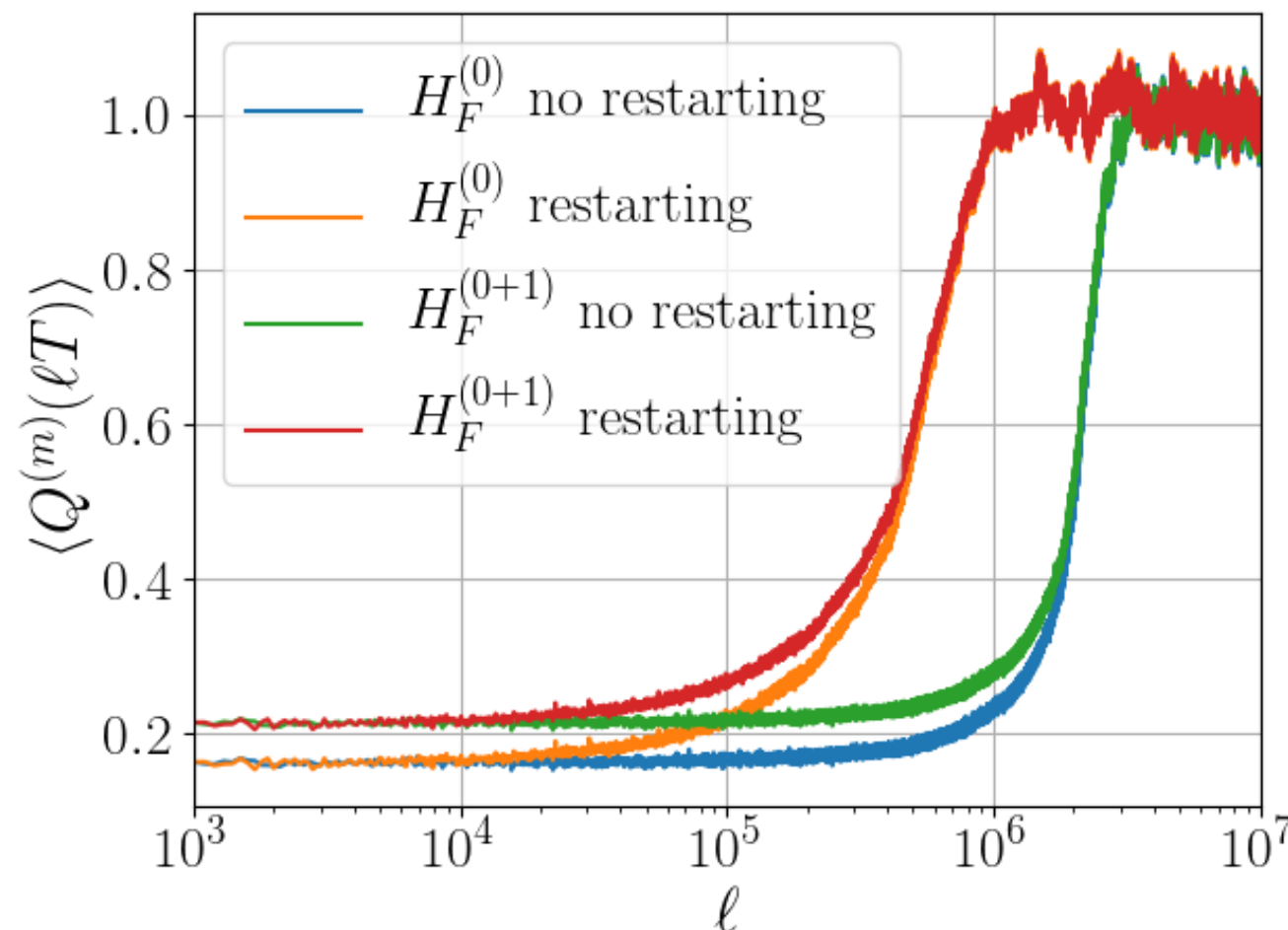
*agreement not exponentially long for **single** trajectories due to chaos*

$$q(t) = \frac{1}{L} \sum_{j=1}^L \mathbf{S}_{F,j}^{\text{exact}}(t) \cdot \mathbf{S}_{F,j}^{[0]}(t)$$



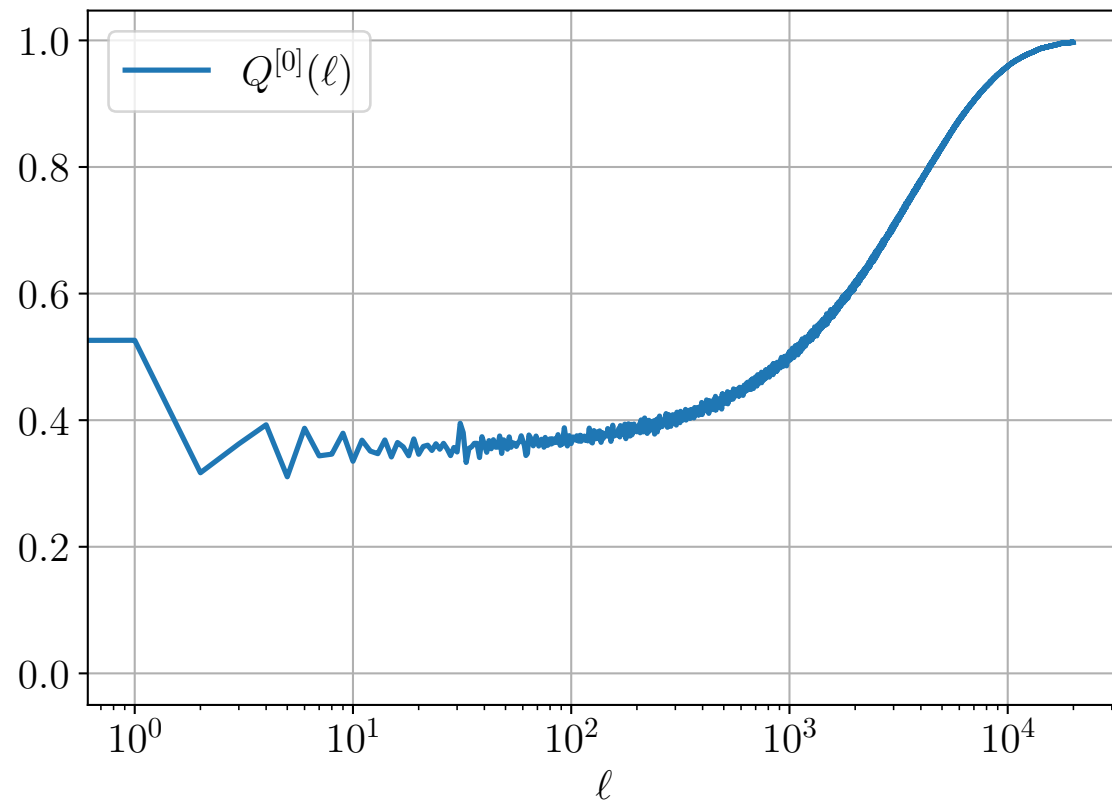
Is the prethermal plateau fully thermal?

- ➔ numerical experiment
- ➔ evolve initial ensemble into prethermal phase
- ➔ **restart ensemble:** measure energy mean and width of ensemble and draw new thermal (Gaussian) ensemble every 1000 cycles

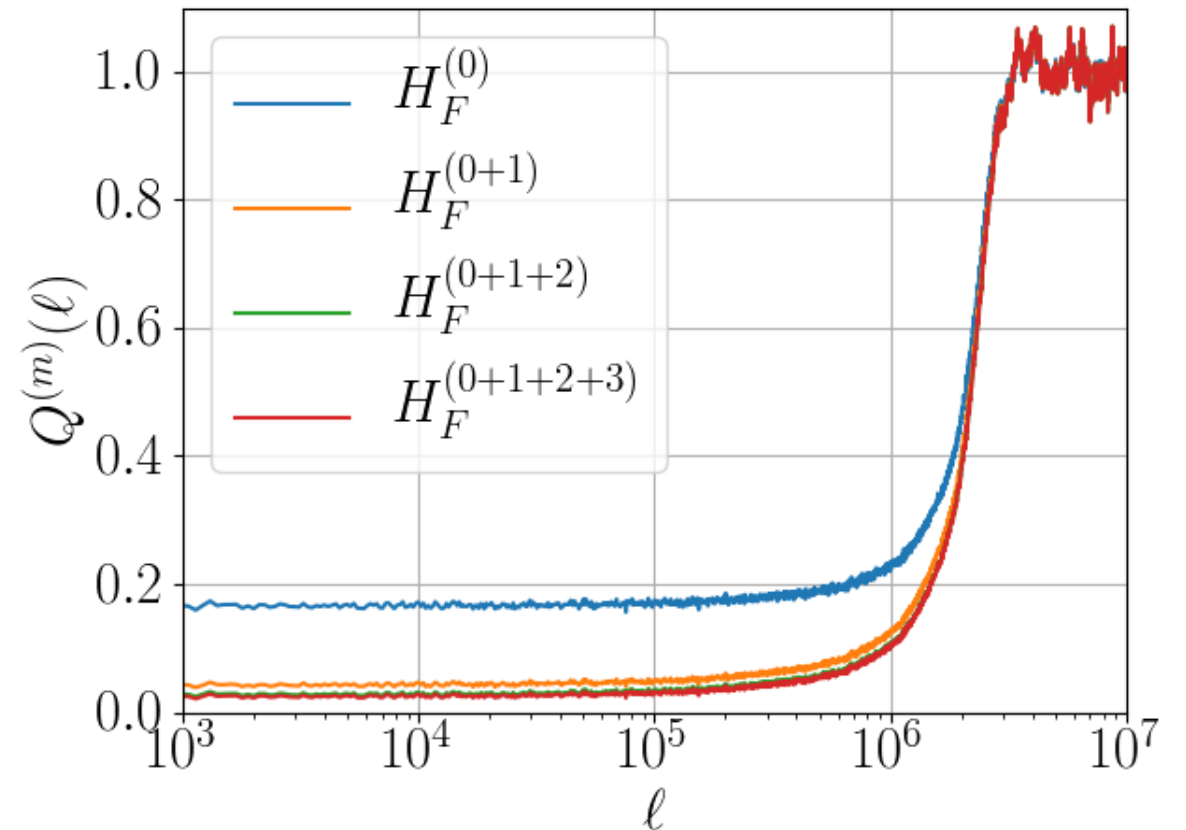


➔ what is quantum about (Floquet) prethermalization?

quantum $L = 22$



classical $L = 100$



Outlook

- ➔ what is quantum about (Floquet) prethermalization?
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Outlook

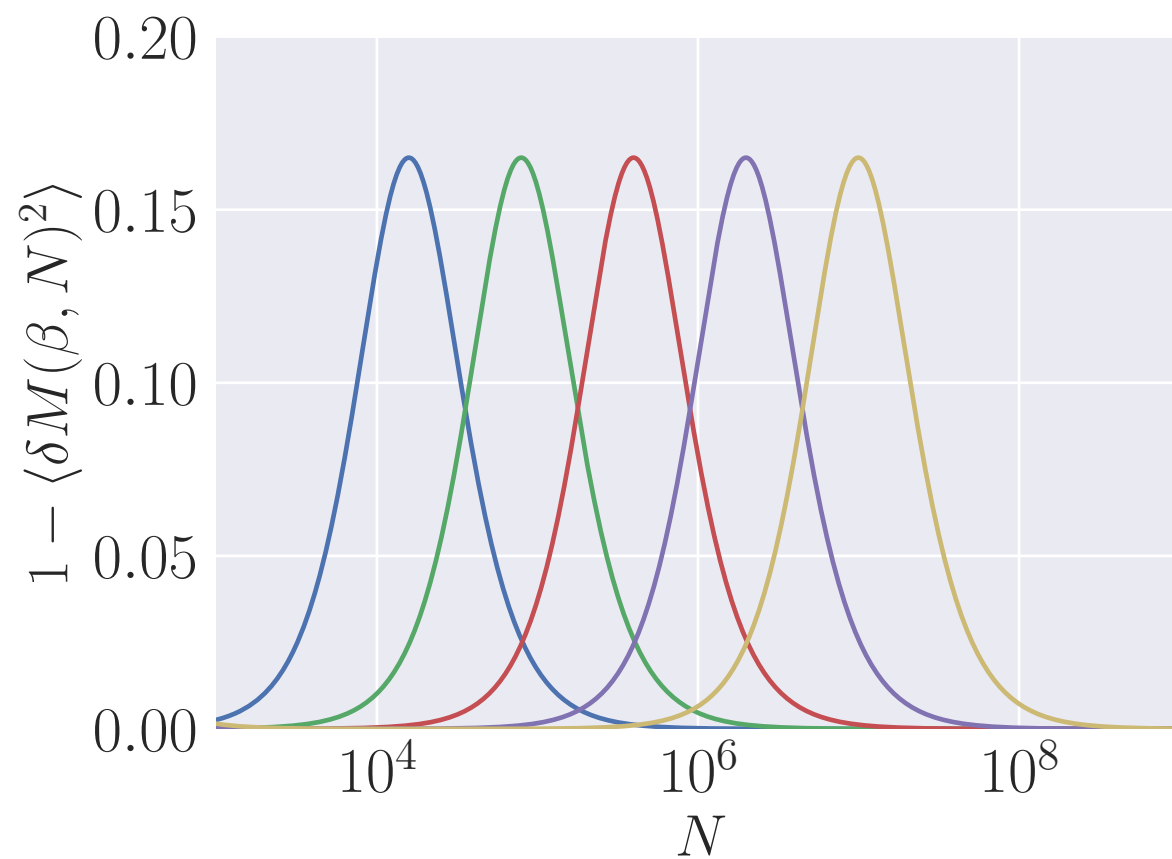
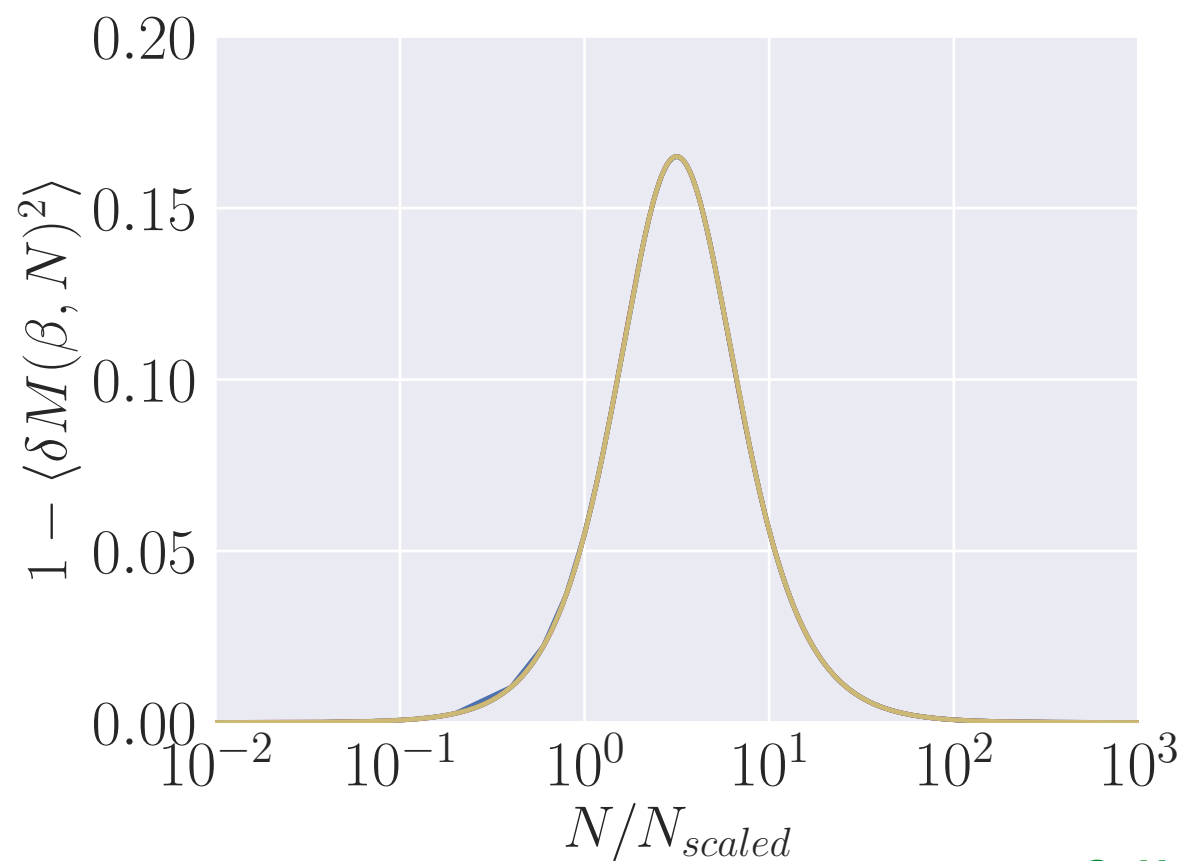
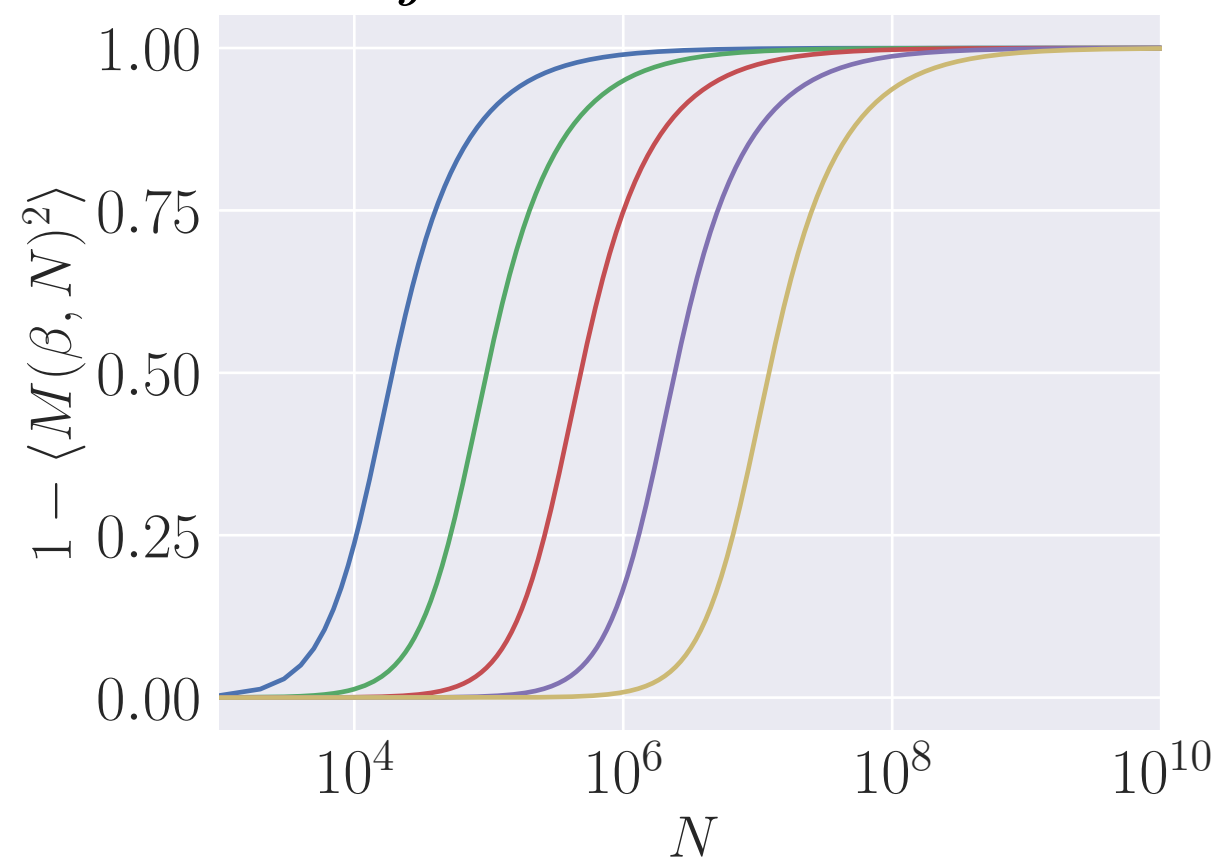
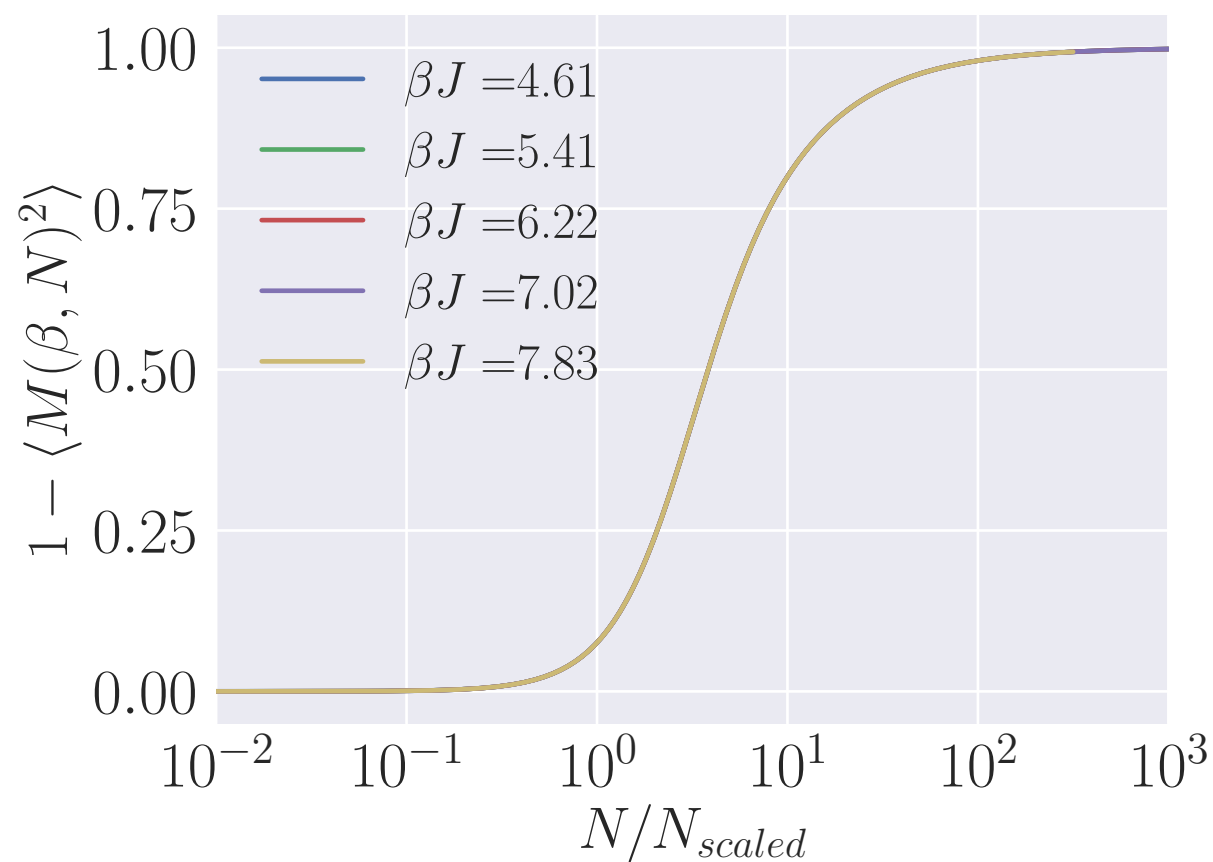
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QuSpin: [*http://weinbe58.github.io/QuSpin*](http://weinbe58.github.io/QuSpin)

python package for ED & many-body dynamics (with P. Weinberg, BU)

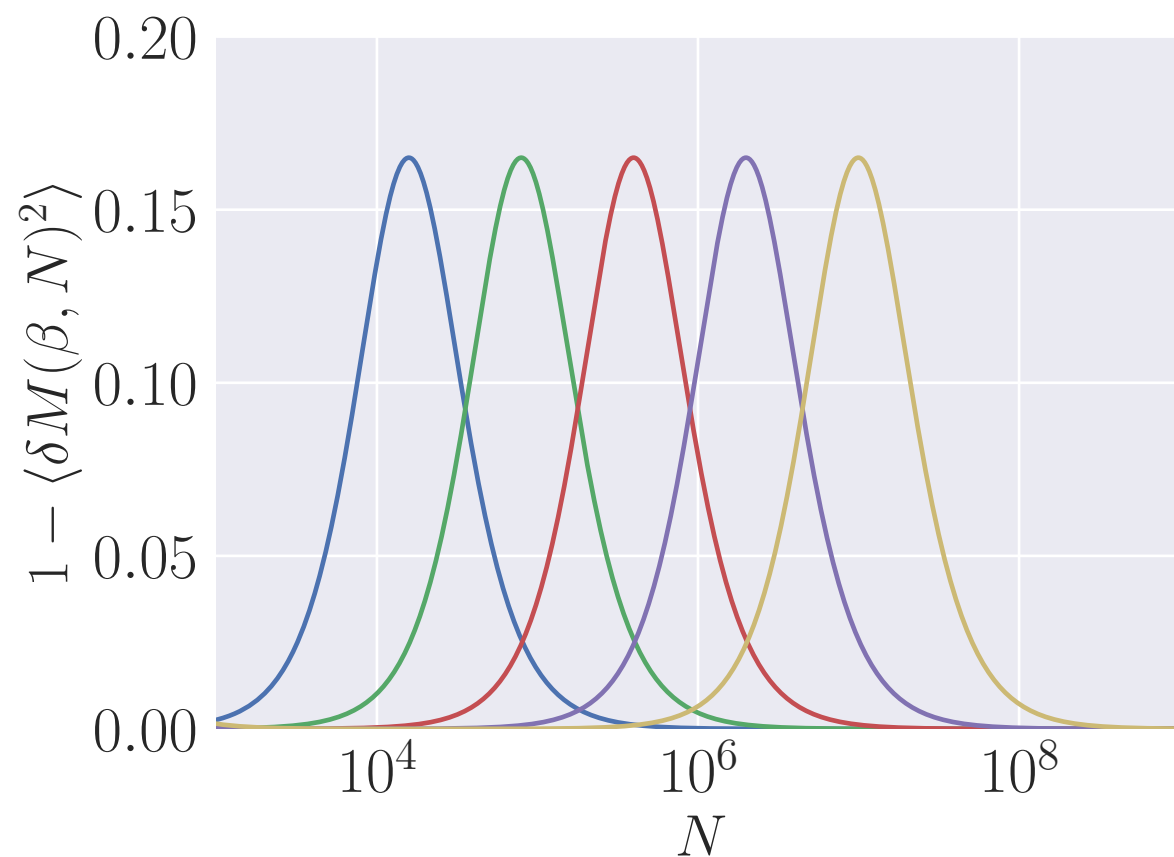
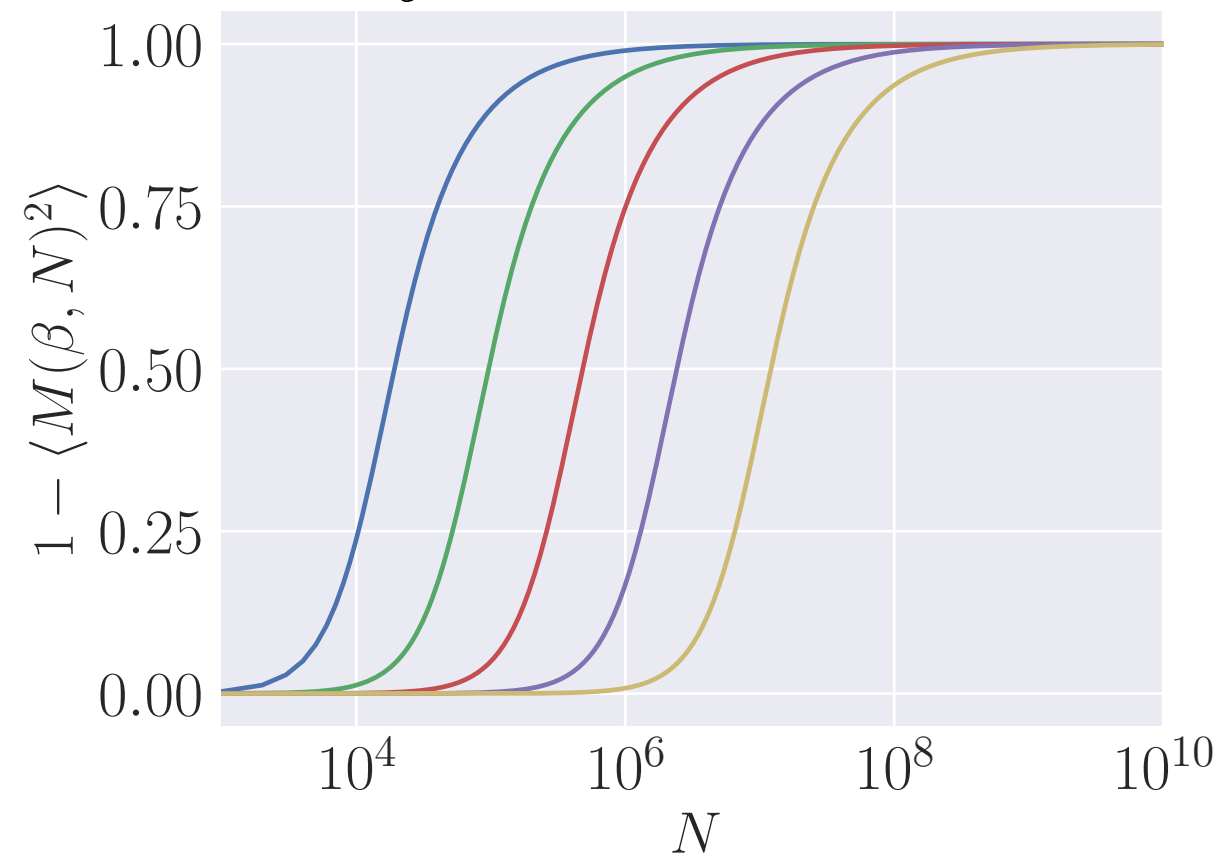
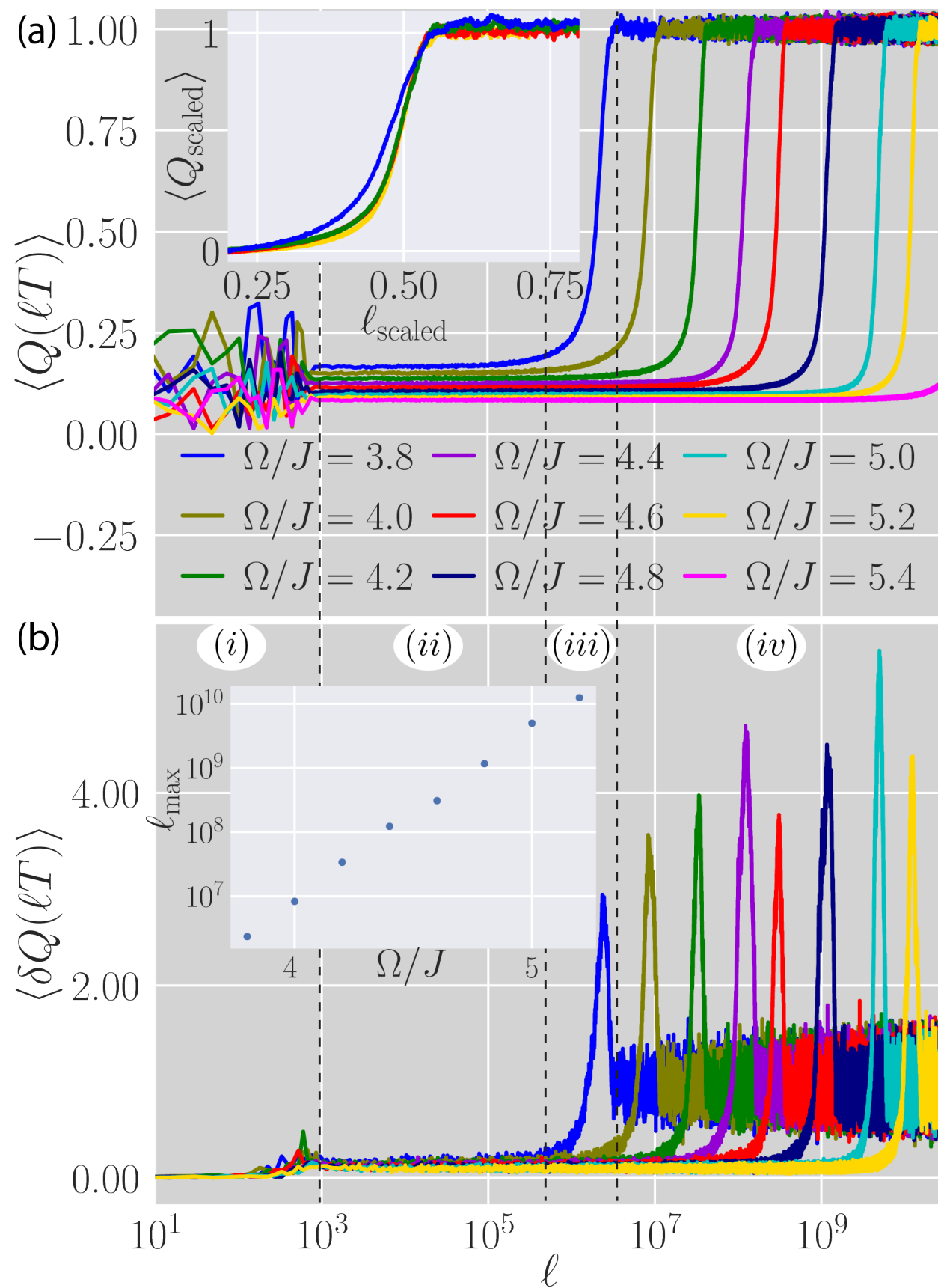
Heating transition 1d Ising:

$$H = -J \sum_j \sigma_j \sigma_{j+1}, \quad \sigma \in \{\pm 1\}$$



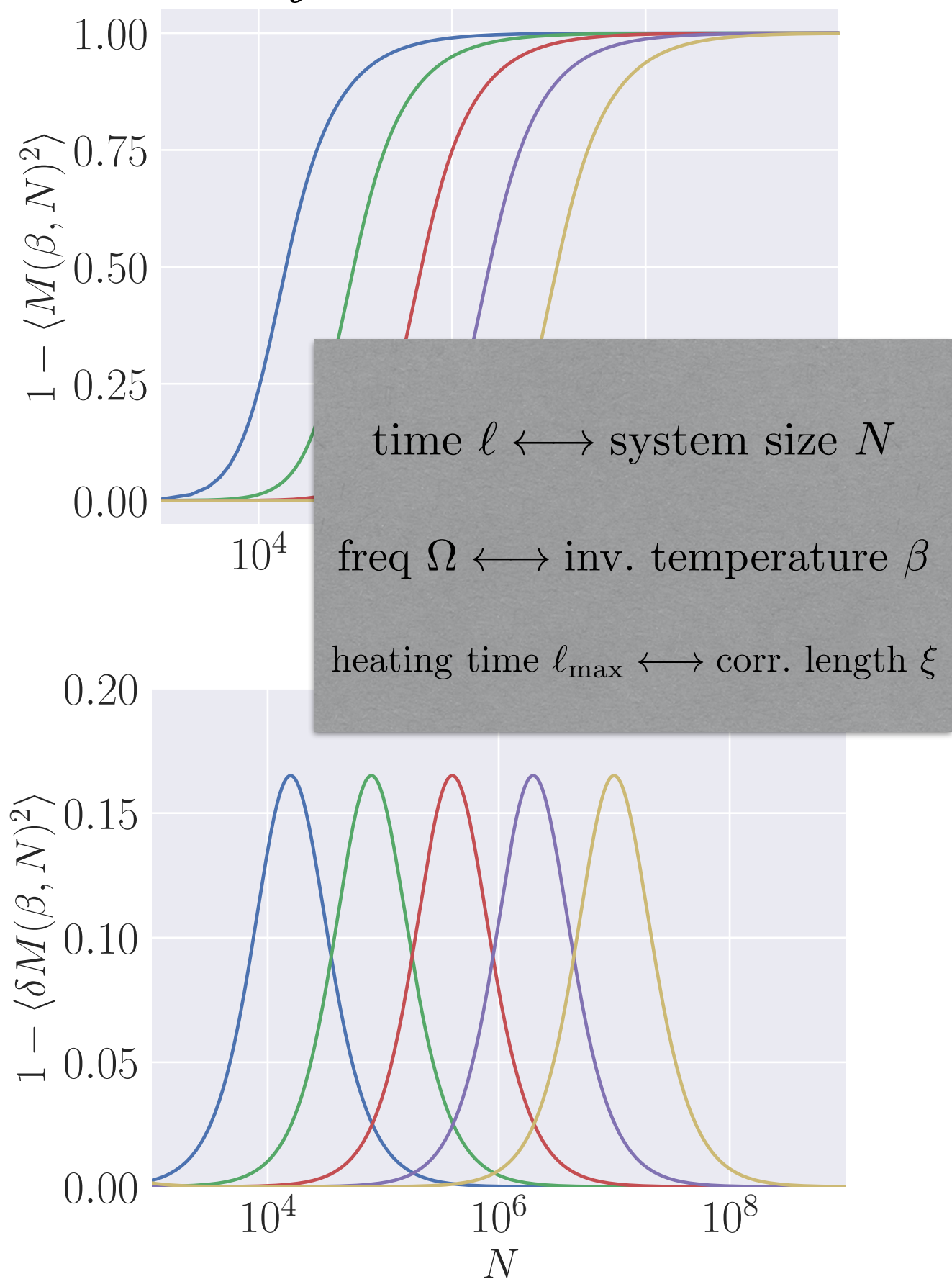
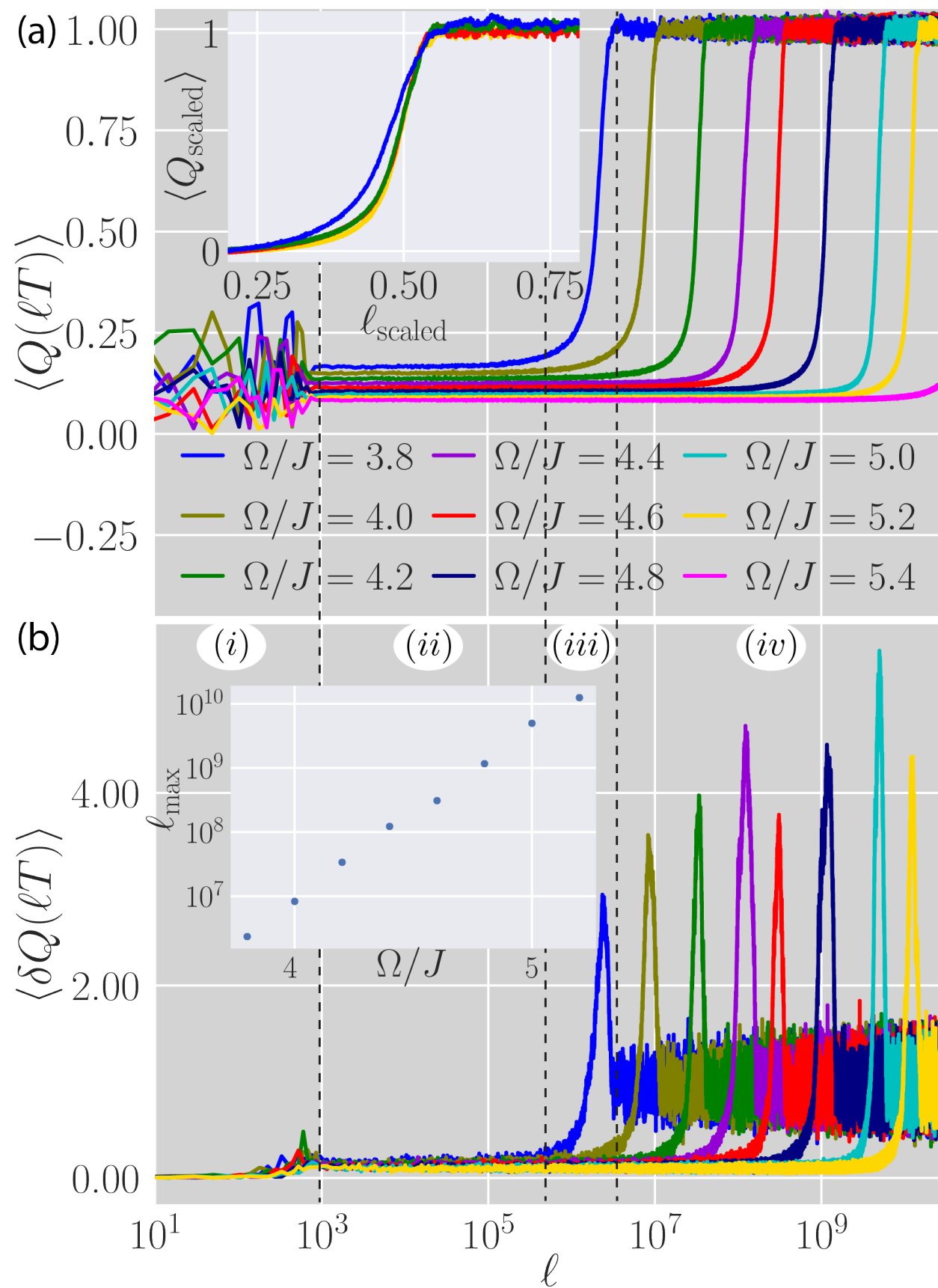
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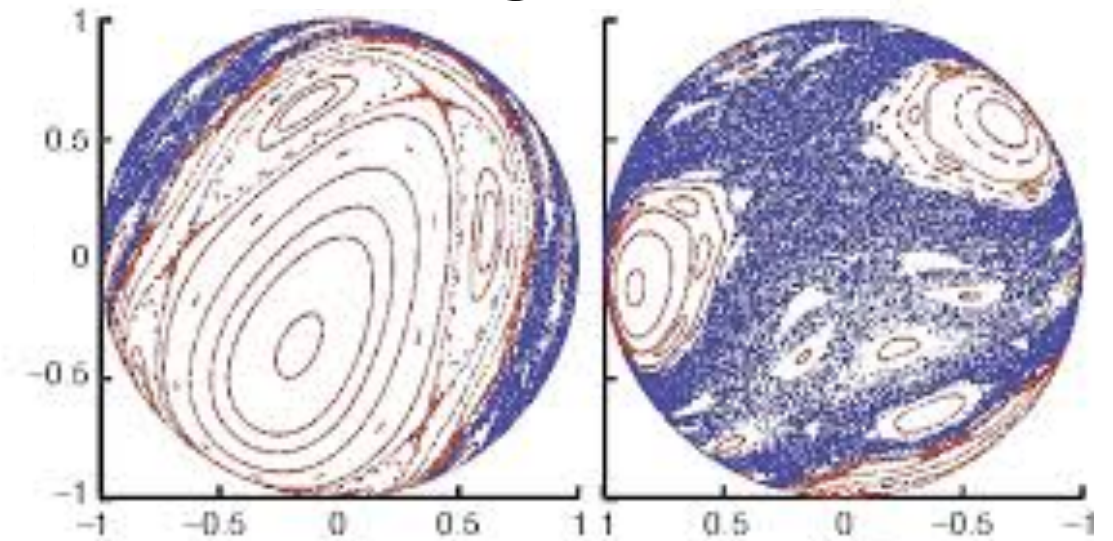


Phase space picture of heating

→ KAM theorem

$$H(I, \phi) = H_0(I) + \epsilon H_1(I, \phi)$$

→ Arnold diffusion



regular dynamics = localization
chaos = diffusion

KAM: **many** solutions to nearly integrable Hamiltonian systems persist under a perturbation for **all** time.

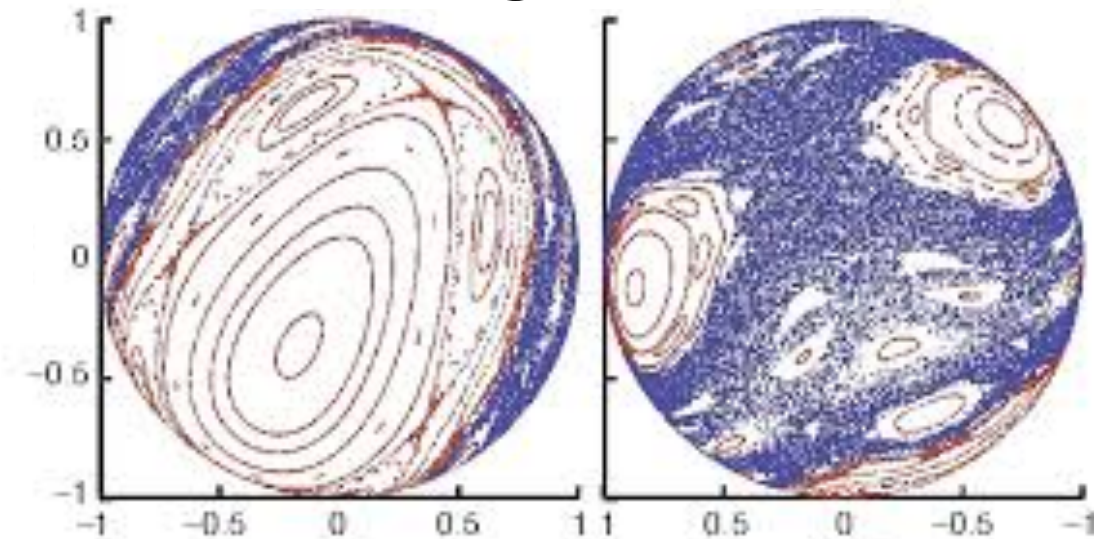
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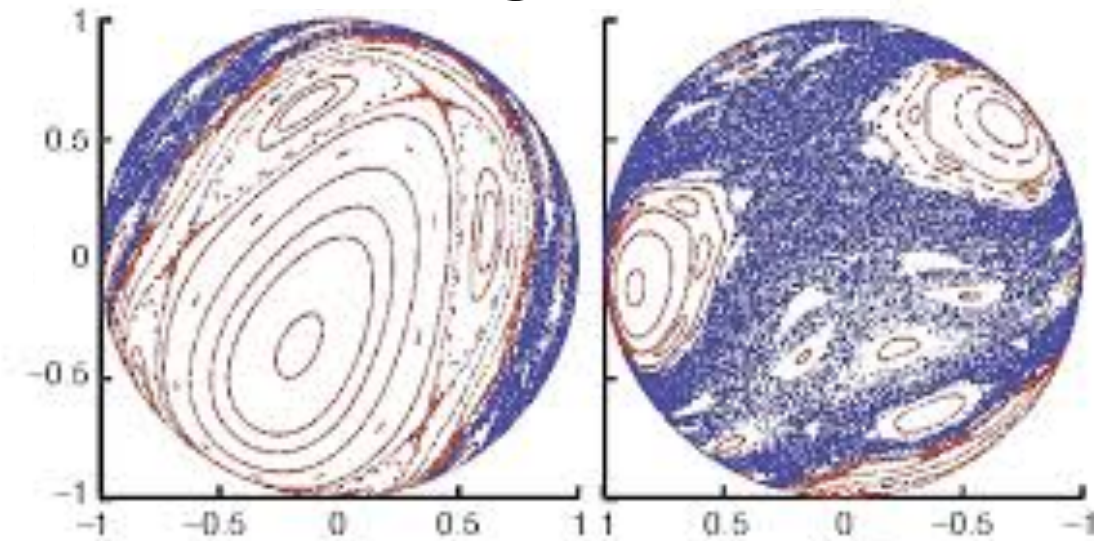
all solutions stay close to their integrable counterparts for an **exponentially long time**. These estimates restrict how quickly solutions can become unstable.

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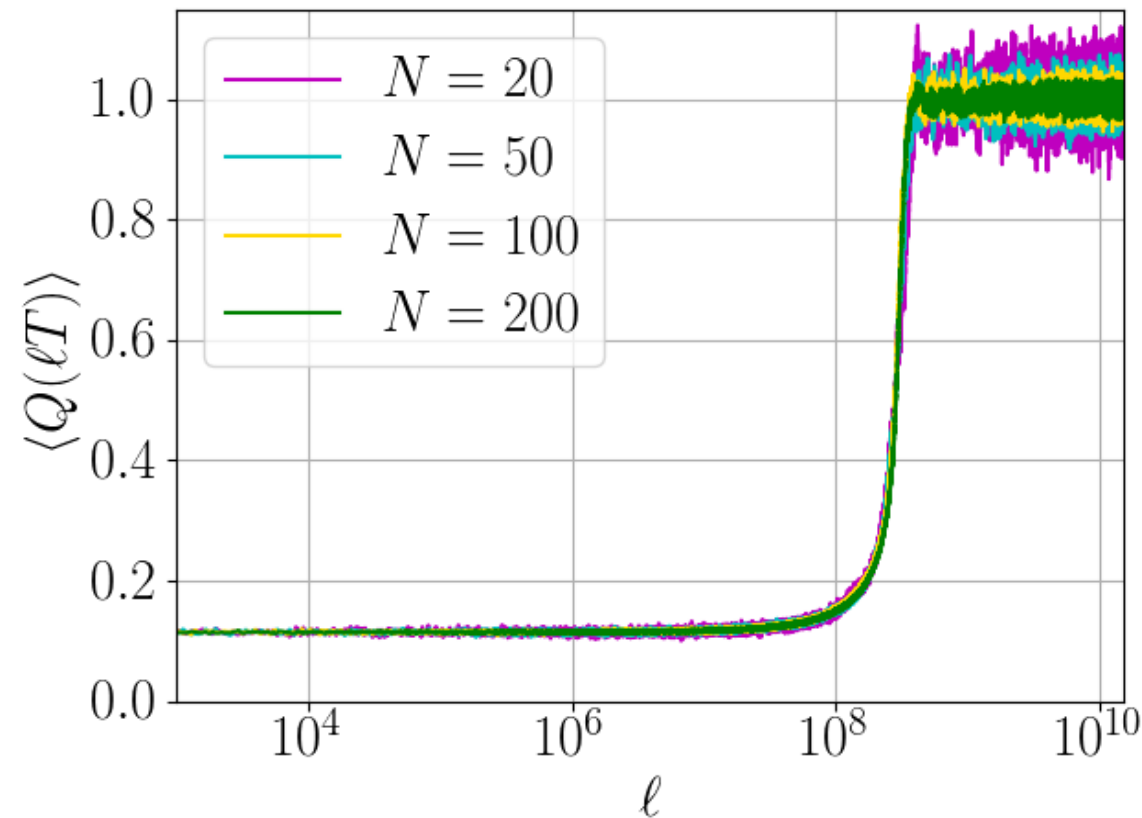
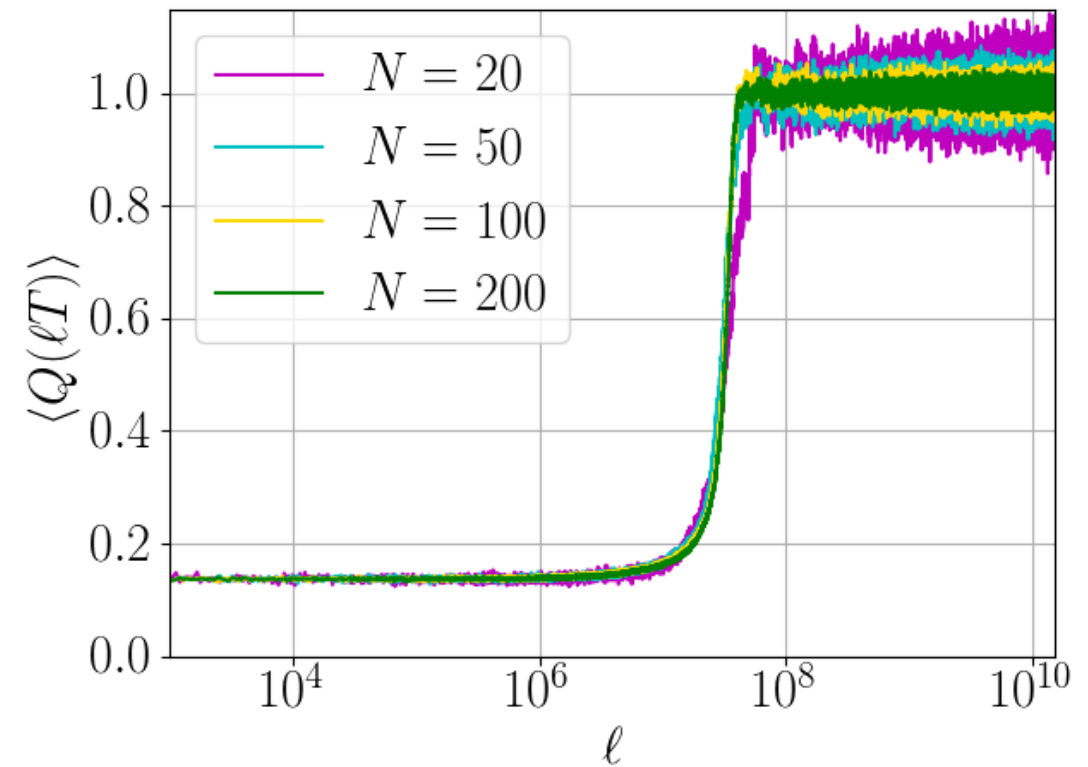
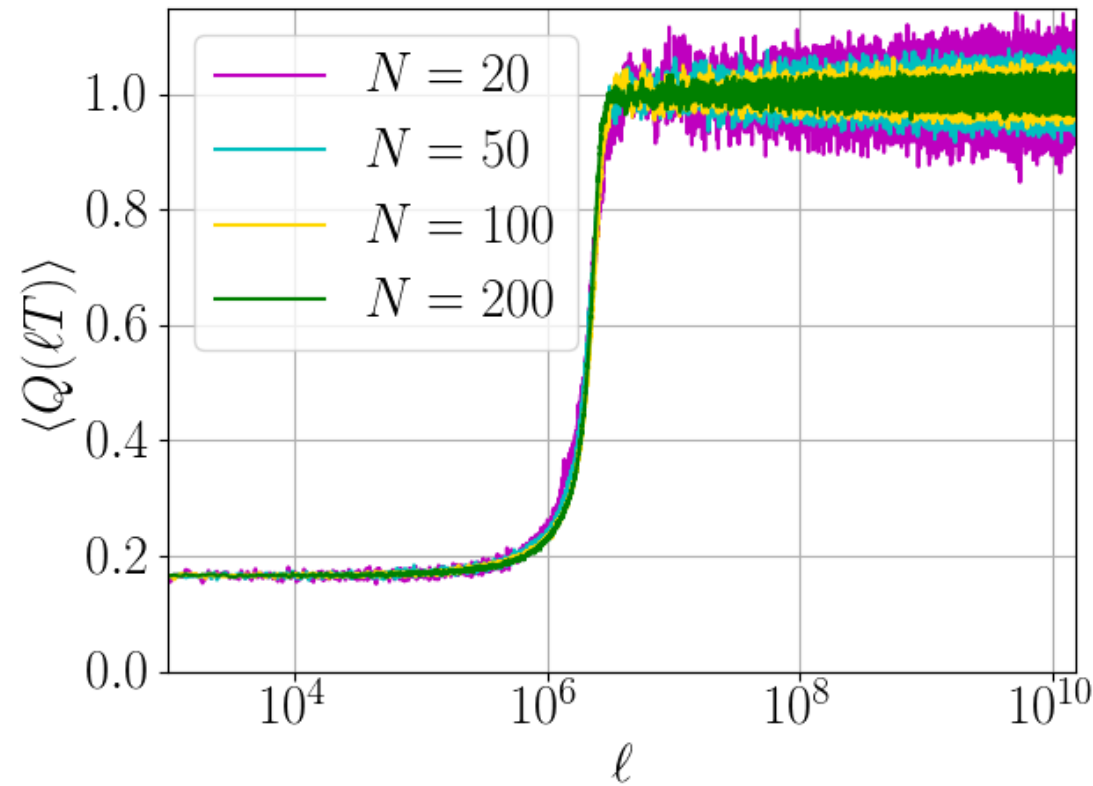
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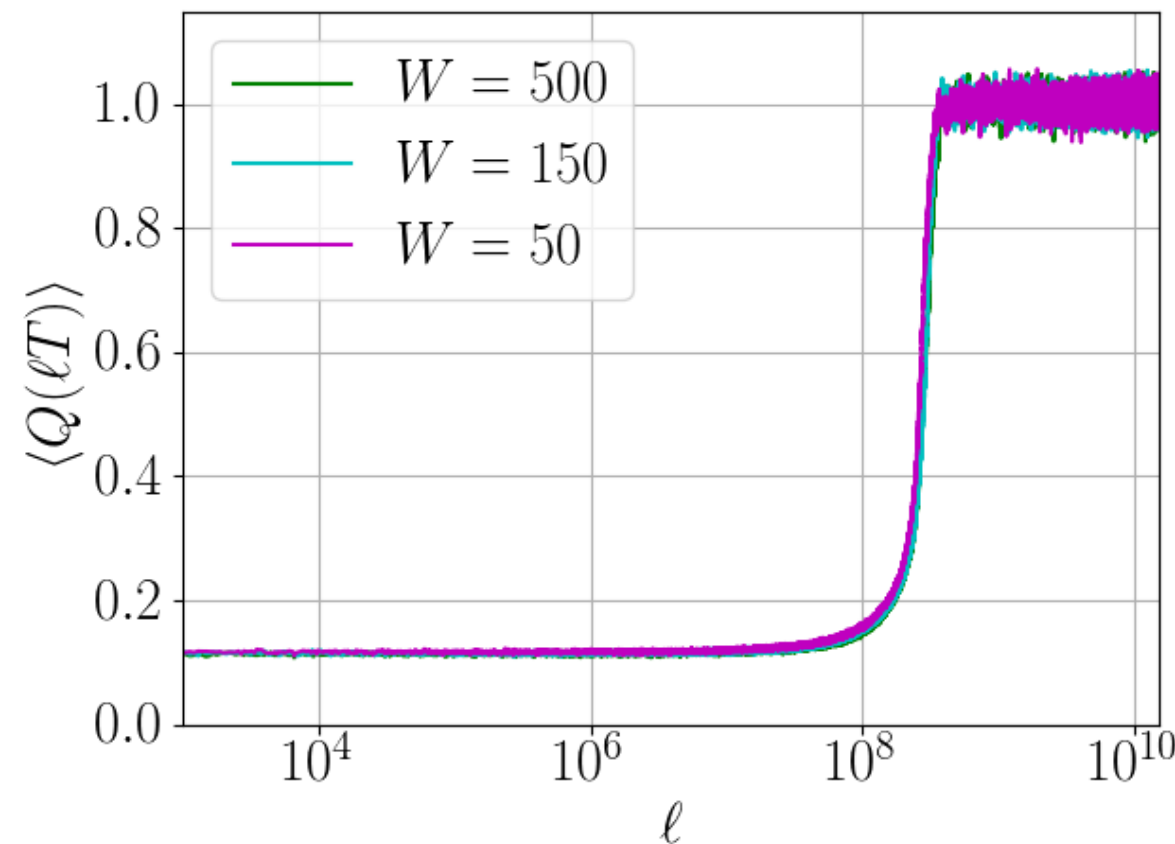
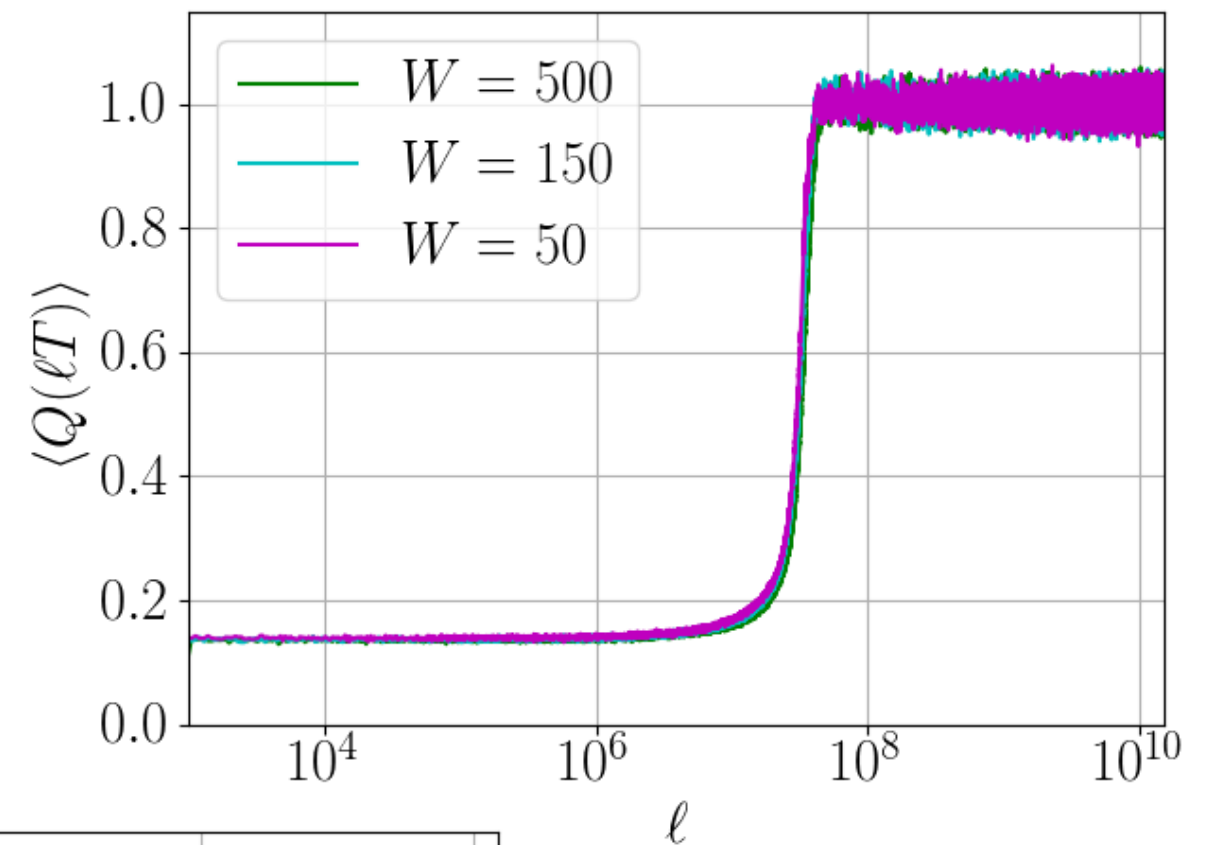
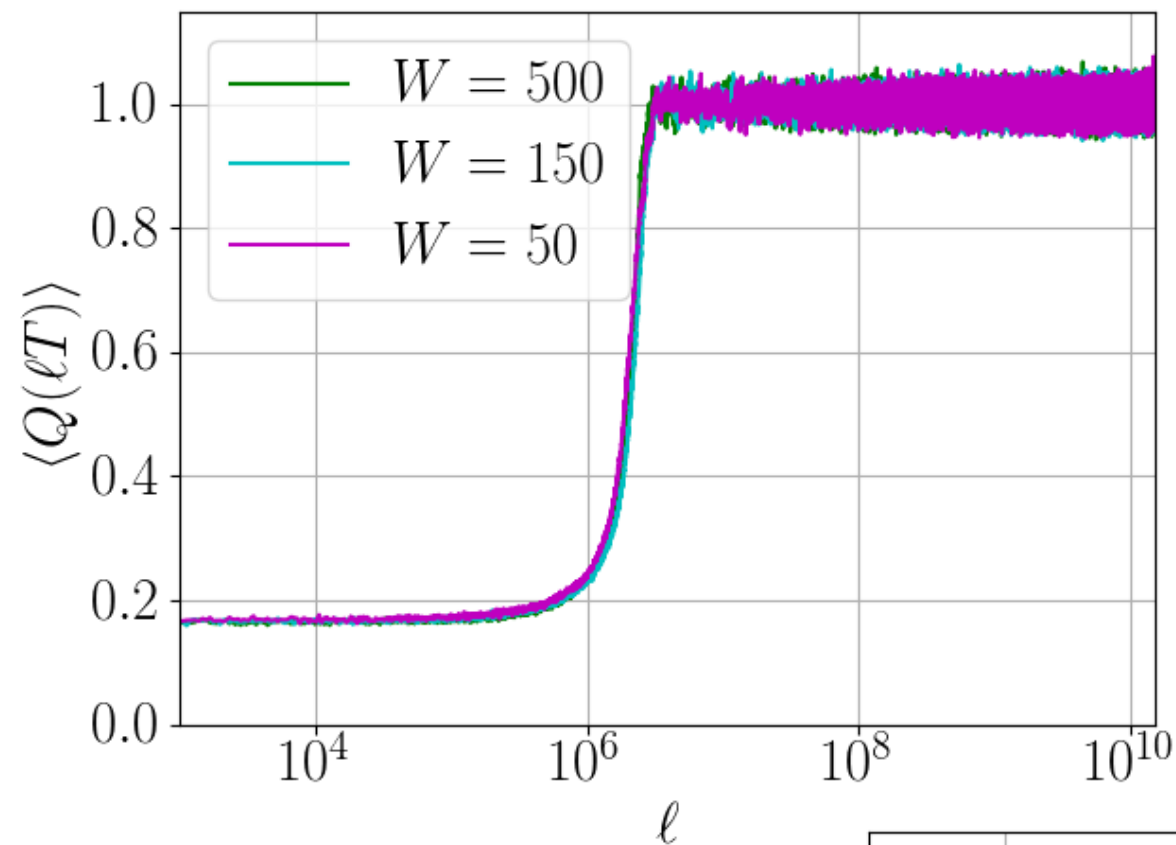
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$$|I(t) - I(0)| < \epsilon^{1/(2L)} \text{ for } |t| < \exp \left(c \left(\frac{1}{\epsilon} \right)^{1/(2L)} \right)$$

Finite size dependence

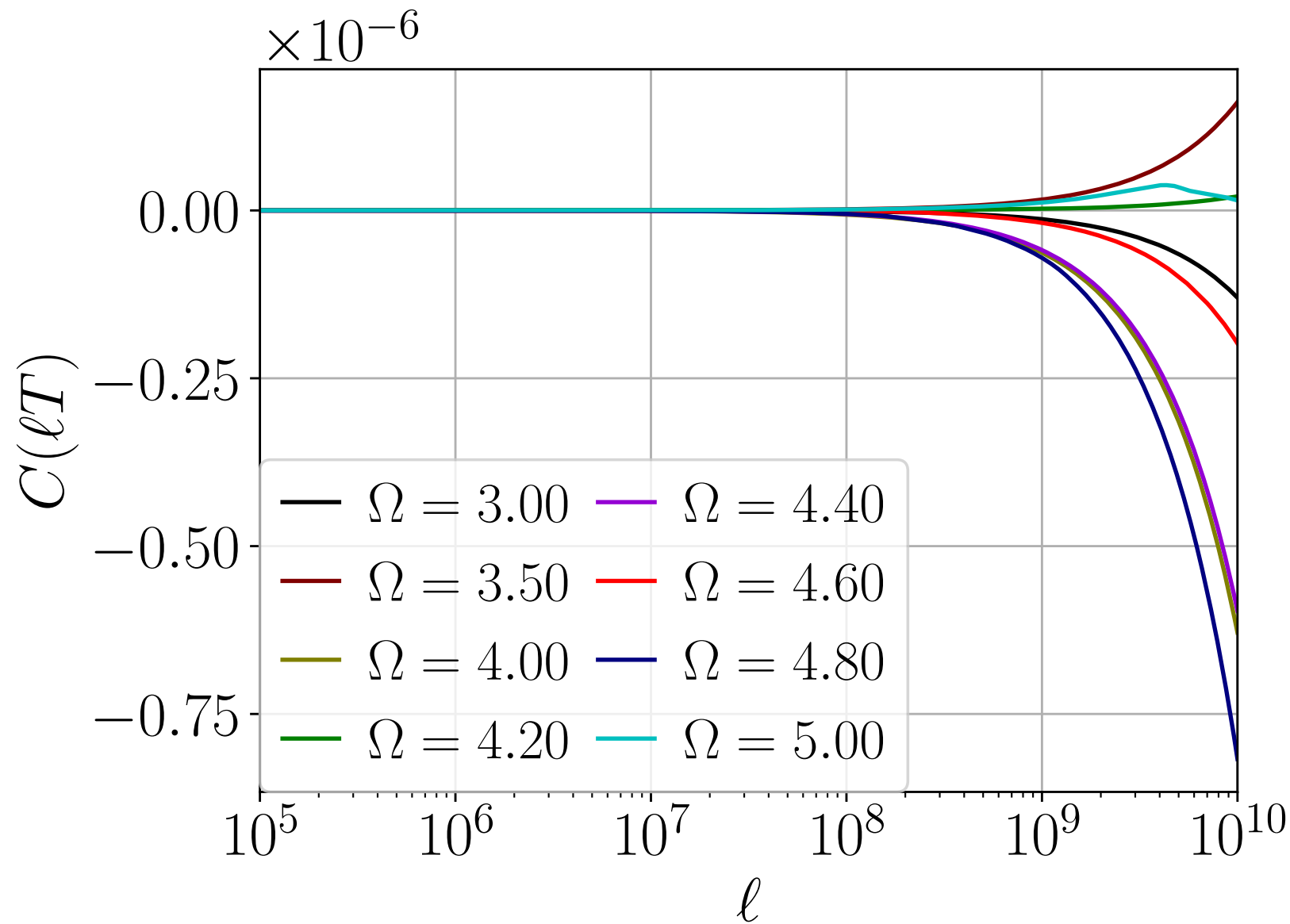


Ensemble noise dependence



$$[-\pi/W, \pi/W]$$

EOM integration error



$$|\vec{S}_j(t)| = 1$$

$$C_j(\ell T) = |\vec{S}_j(\ell T)|^2$$

$$C(\ell T) = \max_j |1 - C_j(\ell T)|$$