

Spreading of a wave packet in a disordered medium

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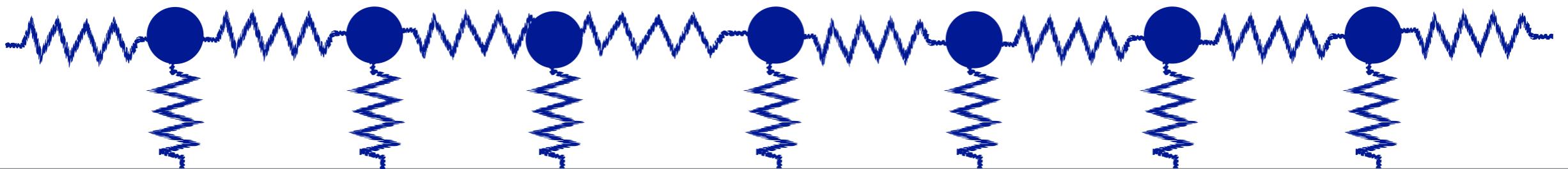
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Classical chain of oscillators

$$H(p, q) = \frac{1}{2} \sum_{x \in \mathbb{Z}} p_x^2 + \omega_x^2 q_x^2 + g(q_x - q_{x+1})^2 + \lambda q_x^4$$



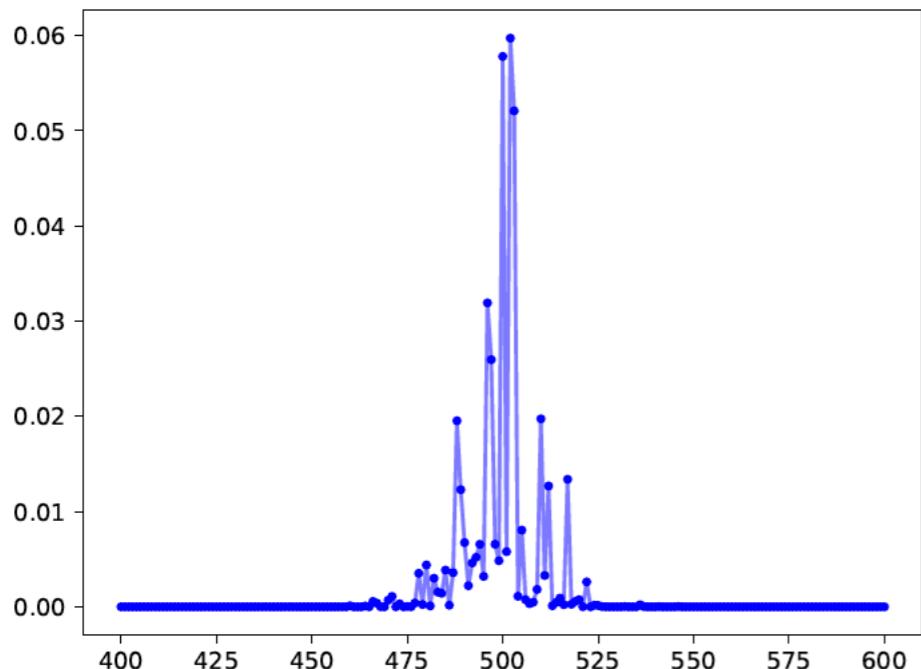
Hamiltonian dynamics:

$$\dot{q} = p, \quad \dot{p} = -\nabla_q H(p, q)$$

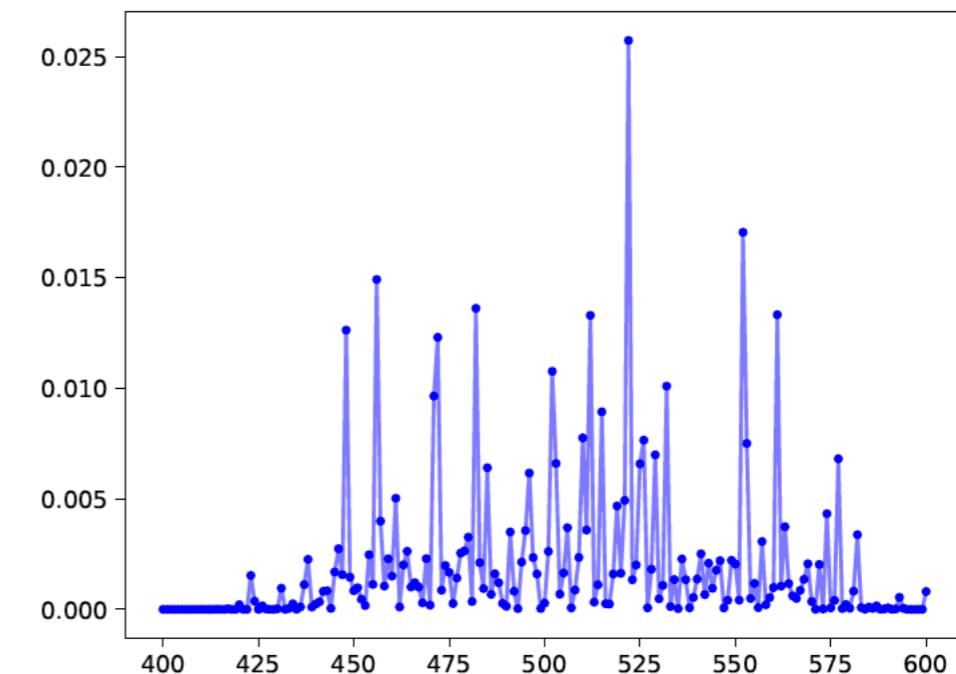
Spreading of a wave packet

The energy is conserved.

Finite quantity of energy in the system (zero temperature)



initial packet



later on

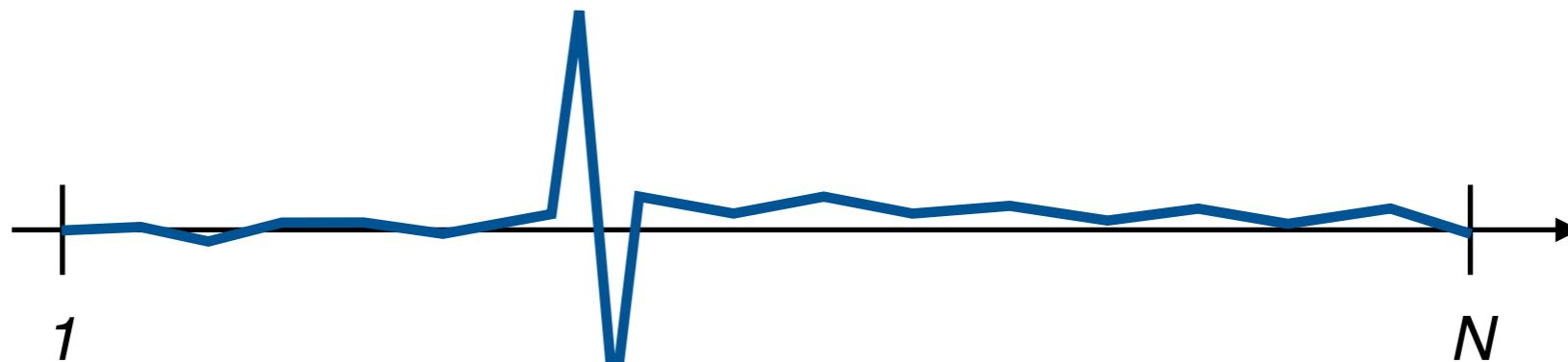
$$H = \sum_x H_x$$

Harmonic case ($\lambda=0$)

Linear equations of motion (Anderson localization):

$$\ddot{q} = -(V - g\Delta)q$$

If $(\omega_x)_{x \in \mathbb{Z}}$ i.i.d. , the eigenmodes are localized :

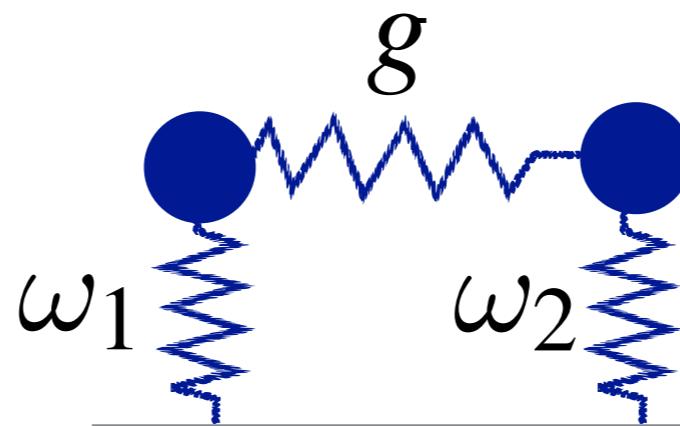


a mode for a chain of size N

$$\mathbb{E} \left(\sum_{\psi} |\psi(x)\psi(y)| \right) \leq C e^{-c|x-y|}, \quad \forall x, y$$

Intuition for localization

Oscillators at different frequencies “don’t talk to each other”,
i.e. they are not in resonance

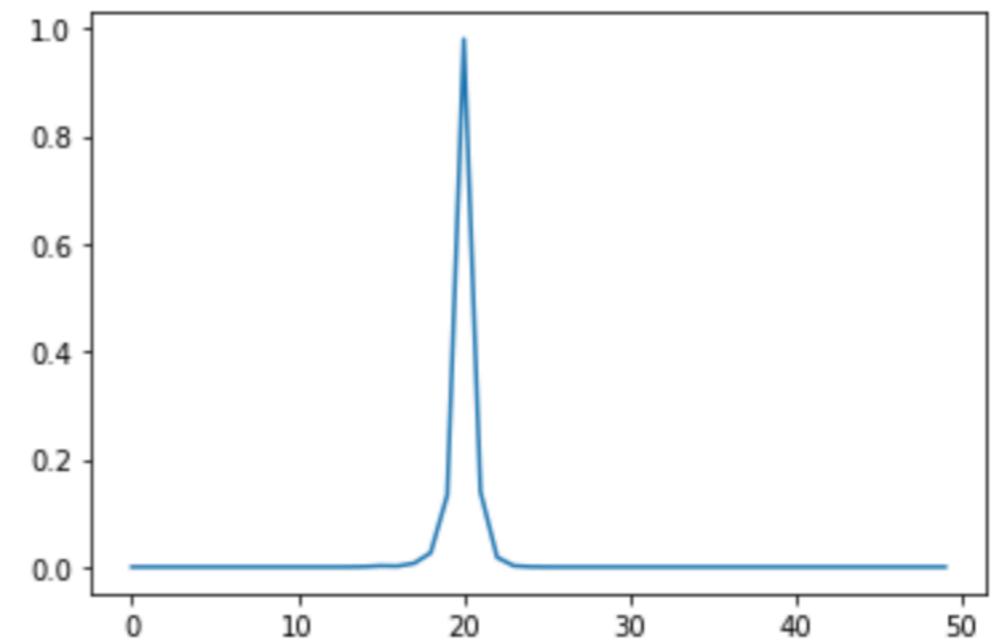
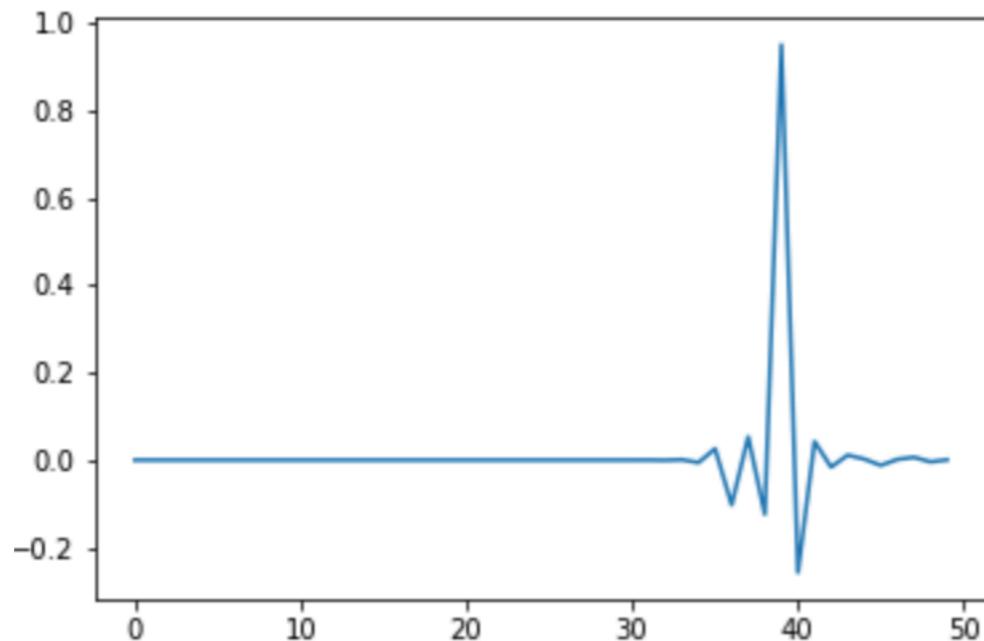


$$-(V - g\Delta) = \begin{pmatrix} -\omega_1 & g \\ g & -\omega_2 \end{pmatrix} \quad g \ll |\omega_1 - \omega_2|$$

$$\psi_1 = (0.99, 0.14), \quad \psi_2 = (0.14, -0.99)$$

Intuition for localization

This still holds true for larger matrices:

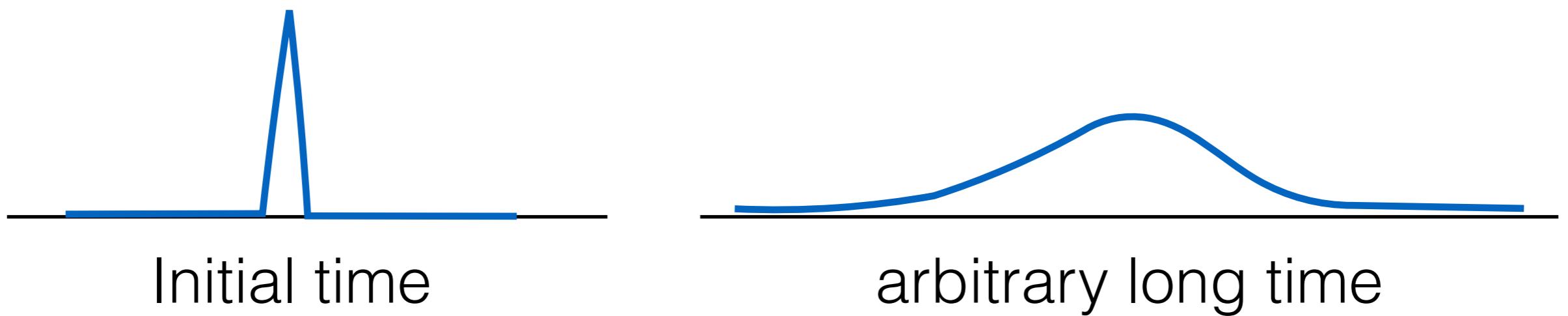


E.g.: two eigenvectors for 50 sites and $g=0.1$

Can be generalized to the full lattice, and much more...

Harmonic case ($\lambda=0$)

The packet does not spread (indefinitely):



Linearity : solution = superposition of localized modes

Anharmonic case ($\lambda \neq 0$)

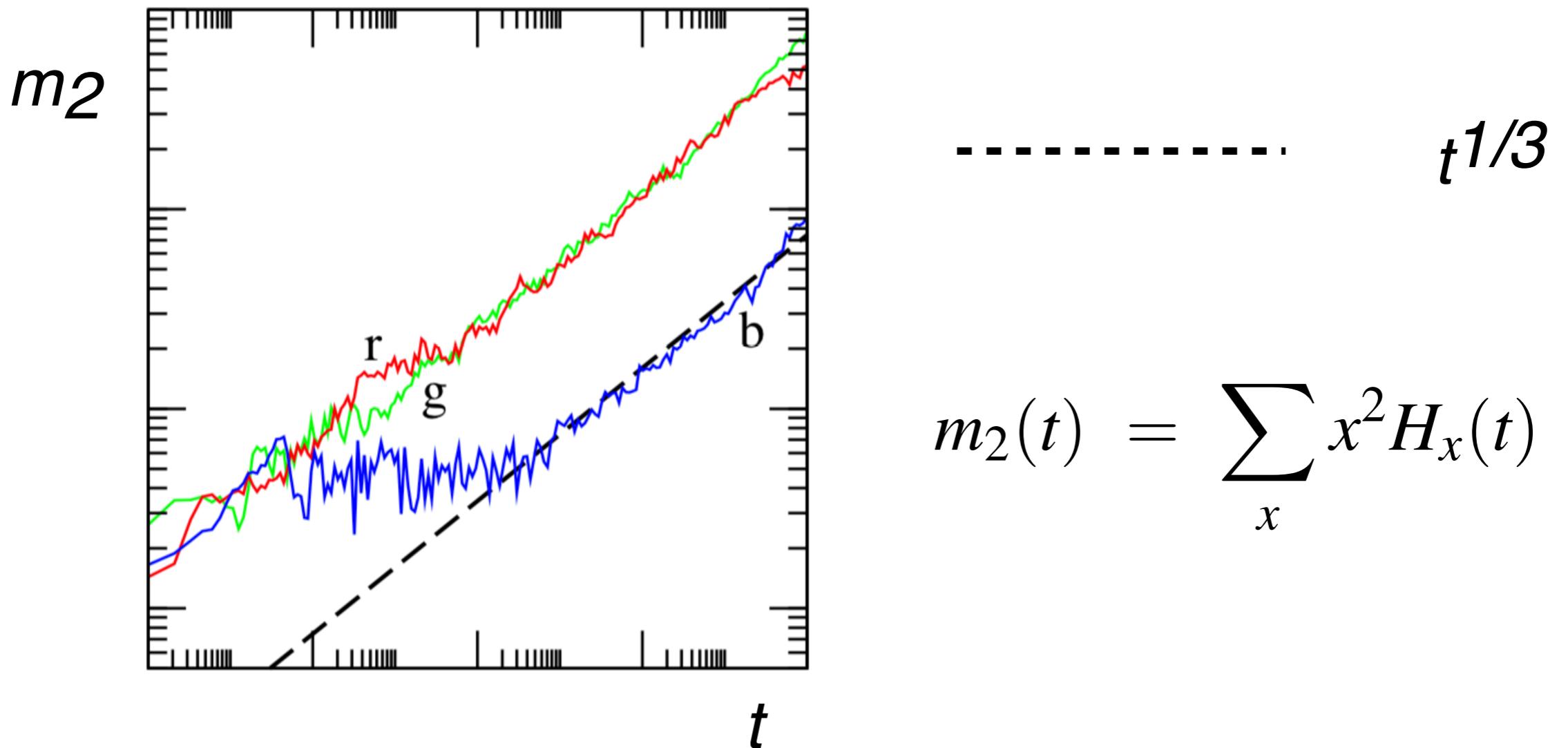
The packet does spread. At which rate ?

- Numerical simulations,
- Analytical computations, mathematical results

The main difficulty is, that there is no regime of parameters, where analytical and numerical results agree for a long time.

From S. Fishman, Y. Krivolapov and A. Soffer (2012)

Numerics : power law



b, g, r : from low to high energy

(you may perhaps think: from small to large λ)

cf. e.g. Flach et al. (2009, 2014, 2020), and many others

Remark:

Small energy density \longleftrightarrow small effective λ

Because anharmonic interactions are given by

$$\lambda q^4 = (\lambda q^2)q^2 \simeq (\lambda E)q^2$$

In equilibrium, the effective non-linearity is indeed

$$\lambda T$$

Analytical : slower than a power law!

Theorem (W.-M. Wang and Z. Zhang, 2009)

«The wave packet stays localized for a very long time with very high probability»

$$\forall n \in \mathbb{N}, \exists \lambda_0 > 0 : \quad \lambda < \lambda_0 \quad \Rightarrow \quad \tau \geq \frac{1}{\lambda^n}$$

- τ is the 1st time that 10% of the energy exits some box around the origin
- with probability that goes quickly to 1 as $\lambda \rightarrow 0$.

Remarks:

- No proper contradiction with the numerics
- Different (more idealized) model (*atomic limit*):

$$H(p, q) = \frac{1}{2} \sum_{x \in \mathbb{Z}} p_x^2 + \omega_x^2 q_x^2 + \lambda_1 (q_x - q_{x+1})^2 + \lambda_2 q_x^4$$

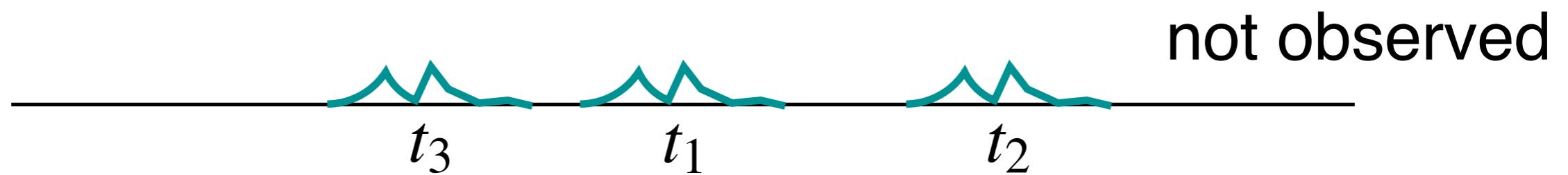
$$\lambda = \lambda_1 \vee \lambda_2$$

- Several other results of this type. E.g. improved bounds by H. Cong, Y. She and Z. Zhang (2020)

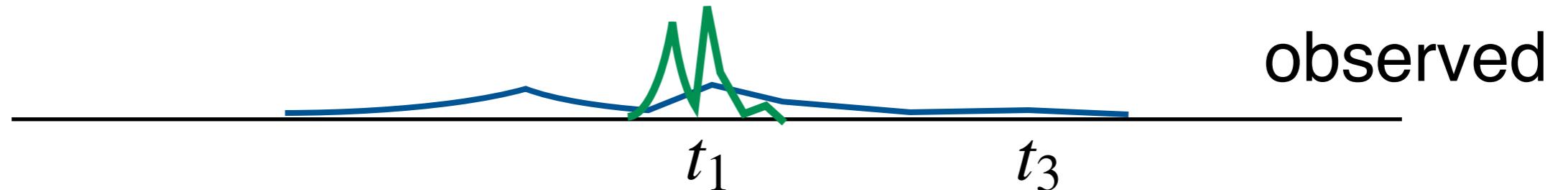
Another analytical result

Let us first contemplate two scenarios for « spreading »:

1. Wandering of a hot spot



2. Proper spreading



A theorem for yet another model

$$H(p, q) = \frac{1}{2} \sum_{x \in \mathbb{Z}} p_x^2 + \omega_x^2 q_x^2 + \lambda(q_x - q_{x+1})^2 + \lambda_x q_x^4$$

$$\lambda_x = \frac{\lambda}{(1 + |x|)^\tau}, \quad \tau > 0$$

Theorem (J. Bourgain and W.-M. Wang, 2007)

« The packet spreads slower than any power law in time. »

$$\forall n \geq 1, \exists \lambda_0 > 0 : \quad \lambda < \lambda_0 \quad \Rightarrow \quad \sum_{x \in \mathbb{Z}} x^2 H_x(t) \leq t^{1/n}$$

a.s. for all $t \geq 0$ provided that this quantity is finite at $t = 0$

Remark: If the packet properly *spreads*, the effective λ decays.

Can numerics be misleading?

The observed spreading is actually *very slow* :

$$m_2^{1/2}(t) \sim t^{1/6}$$

- When the packet spreads, the effective non-linearity (λq^4) decays.
- If the spreading is slow, you need a lot of time for the effective non-linearity to decay, and so you need a lot of time to change regime and see another power law.

Cheap but not unrealistic to think that numerics
were not run for a long enough time

Direct comparison numerics/theory

Recreate a “contradiction” numerics/math
that can be decided

We will define a quantity $I(t)$

- for the original model (technical issues),
- starting from equilibrium (simpler),
- that can be controlled by a theorem.

Roughly $I(t)$ measures the loss of memory in the system

Definition of $I(t)$

Another way to decompose the Hamiltonian at $\lambda = 0$:

$$\begin{aligned}\sum_x H_x &= \frac{1}{2} \langle p, p \rangle + \langle q, (V - g\Delta)q \rangle \\ &= \frac{1}{2} \sum_E |\langle p, E \rangle|^2 + E |\langle q, E \rangle|^2 = \sum_E H_E\end{aligned}$$

with

$$(V - g\Delta)|E\rangle = E|E\rangle$$

The energy of each mode is conserved at $\lambda = 0$:

$$\frac{dH_E}{dt} = 0 \quad \forall E$$

Definition of $I(t)$

For the coupled dynamics, $H = H_0 + \lambda H_1$, we define

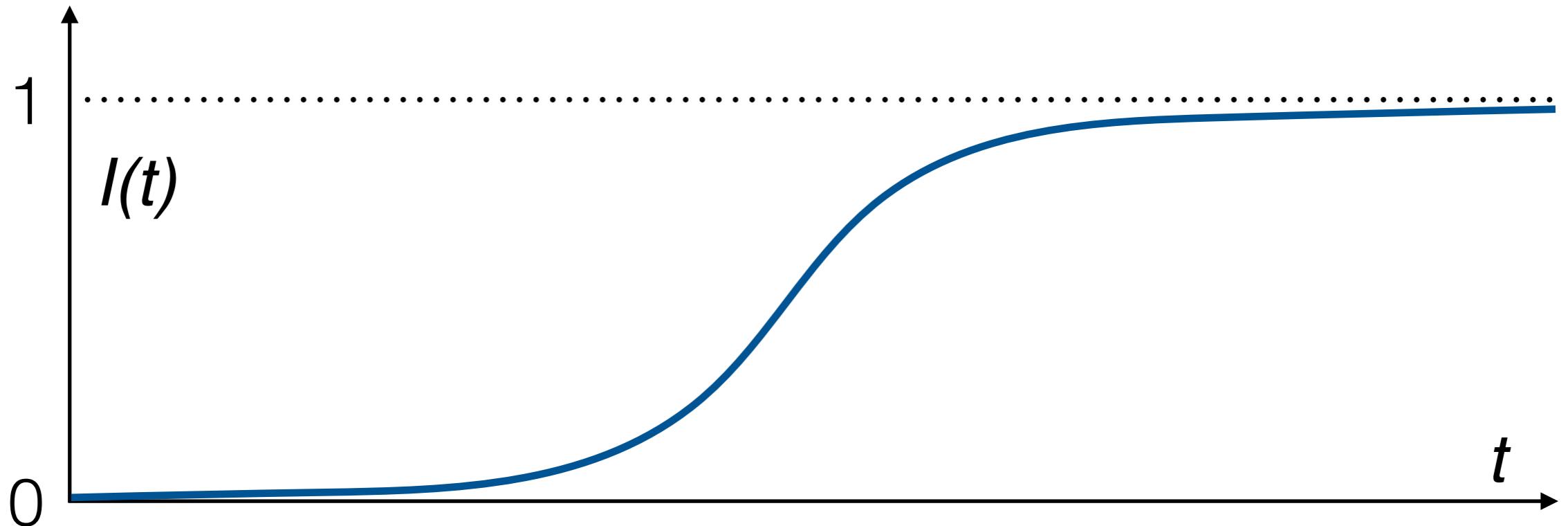
$$I(t) = \frac{1}{N} \sum_E \frac{\langle (H_E(t) - H_E(0))^2 \rangle_T}{2 \operatorname{var}(H_E)}$$

where $\langle f \rangle_T$ is the Gibbs state at temperature T :

$$\langle f \rangle_T = \frac{1}{Z} \int f(q, p) e^{-H(q, p)/T} dq dp$$

(this is an equilibrium measure for the dynamics)

Expected behavior for $I(t)$



$I(0) = 0$: by definition

$I(+\infty) = 1$: in the large N limit, $\langle H_E(t); H_E(0) \rangle_T \rightarrow 0$ as $t \rightarrow \infty$

Rigorous bound on $I(t)$

Theorem (W. De Roeck, F. H. and O. Prosnjak).

Let $n \in \mathbb{N}$. There exists a deterministic constant $C_n < +\infty$ such that for all $\lambda \geq 0$ and for all $t \geq 0$,

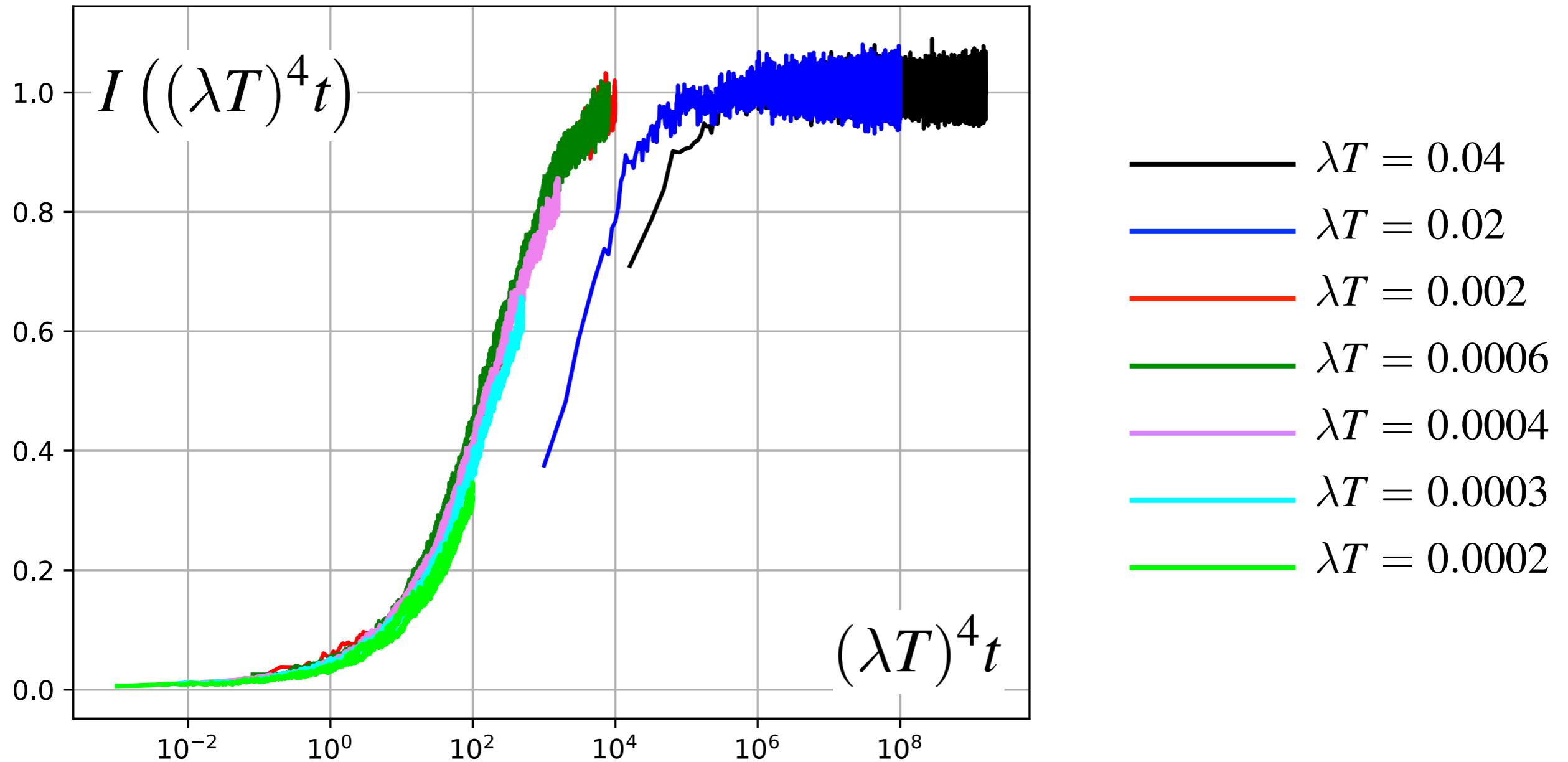
$$\limsup_{N \rightarrow \infty} I_N(t) \leq C_n(\lambda^{2-a} + (\lambda^n t)^2)$$

a.s. with $a < 2$ that can be made explicit.

Remark: We assume that the temperature T is fixed.

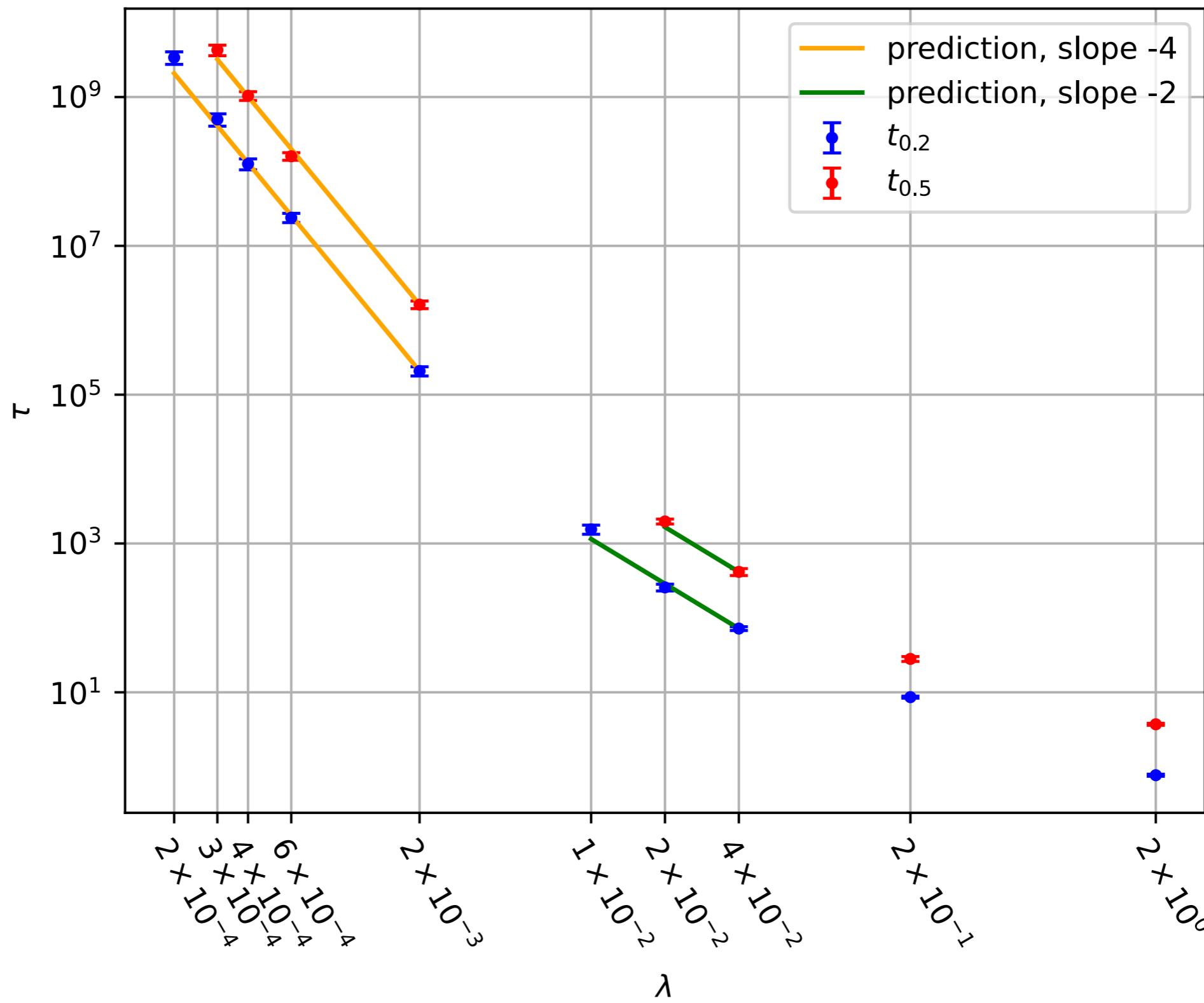
Actually $I(\lambda, T) = \bar{I}(\lambda T)$.

Numerical results for $I(t)$



It would seem that $I(t) = f((\lambda T)^4 t)$, but we know it is not!
(smallest value of λ suggests actually a deviation from this behavior)

Numerical results for $I(t)$



Power law consistent with S. Flach et al.

Back to the spreading of a wave packet. If local equilibrium holds inside the packet:

$$\partial_t E = \partial_x (D(T\lambda) \partial_x E)$$

and $T\lambda$ goes to 0 as the packet spreads

For this non-linear diffusion equation, we find

$$m_2(t) \sim t^{1/3} \quad \Rightarrow \quad D(T\lambda) \sim (T\lambda)^4$$

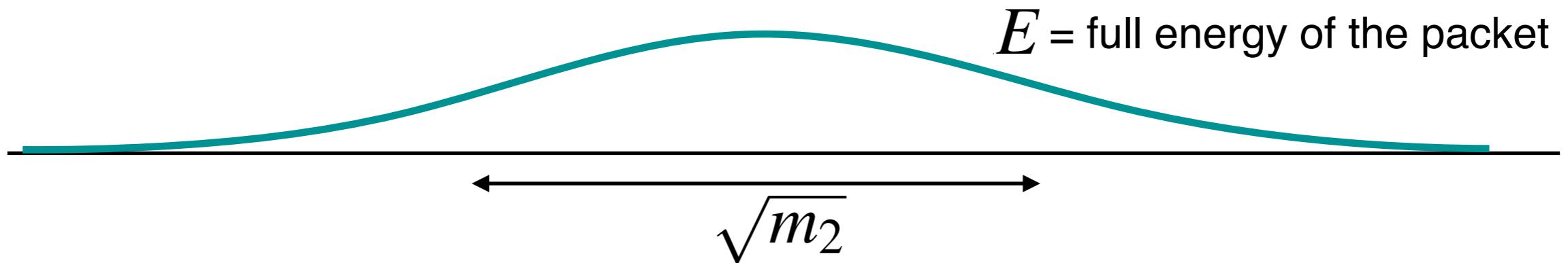
Consistent within linear fluctuating hydrodynamics:

$$I(t) = f(Dt) = f((T\lambda)^4 t)$$

Timescales consistent with S. Flach et al.

Effective temperature of the wave packet:

$$\lambda T \leftrightarrow \frac{\lambda E}{\sqrt{12m_2}}$$



Smallest effective non-linearity that is reached:

$$\lambda T = 0.0005 \text{ (Flach),}$$

$$\lambda T = 0.0002 \text{ (us)}$$

Some catch: Is the packet in equilibrium?

Applying statistical mechanics is challenging:

- finite amount of energy,
- nearly conserved local quantities

Observation: pre-thermal plateau for the *clean* chain

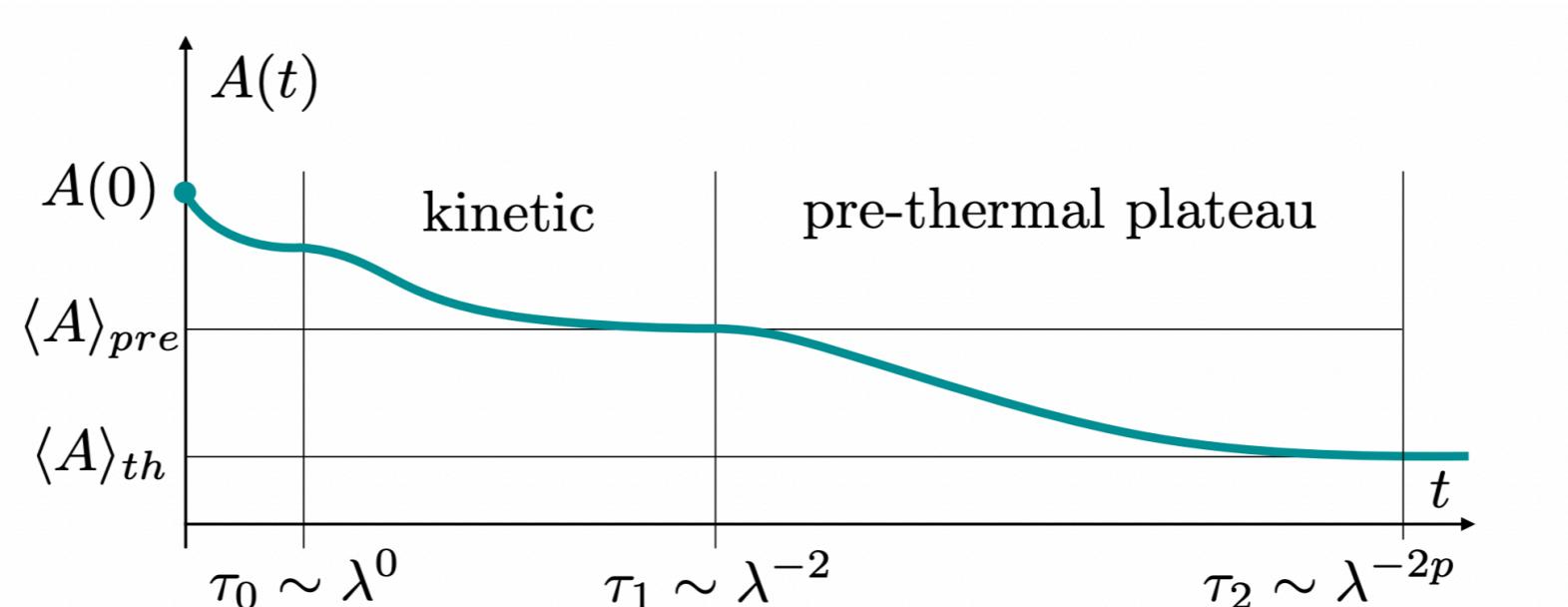
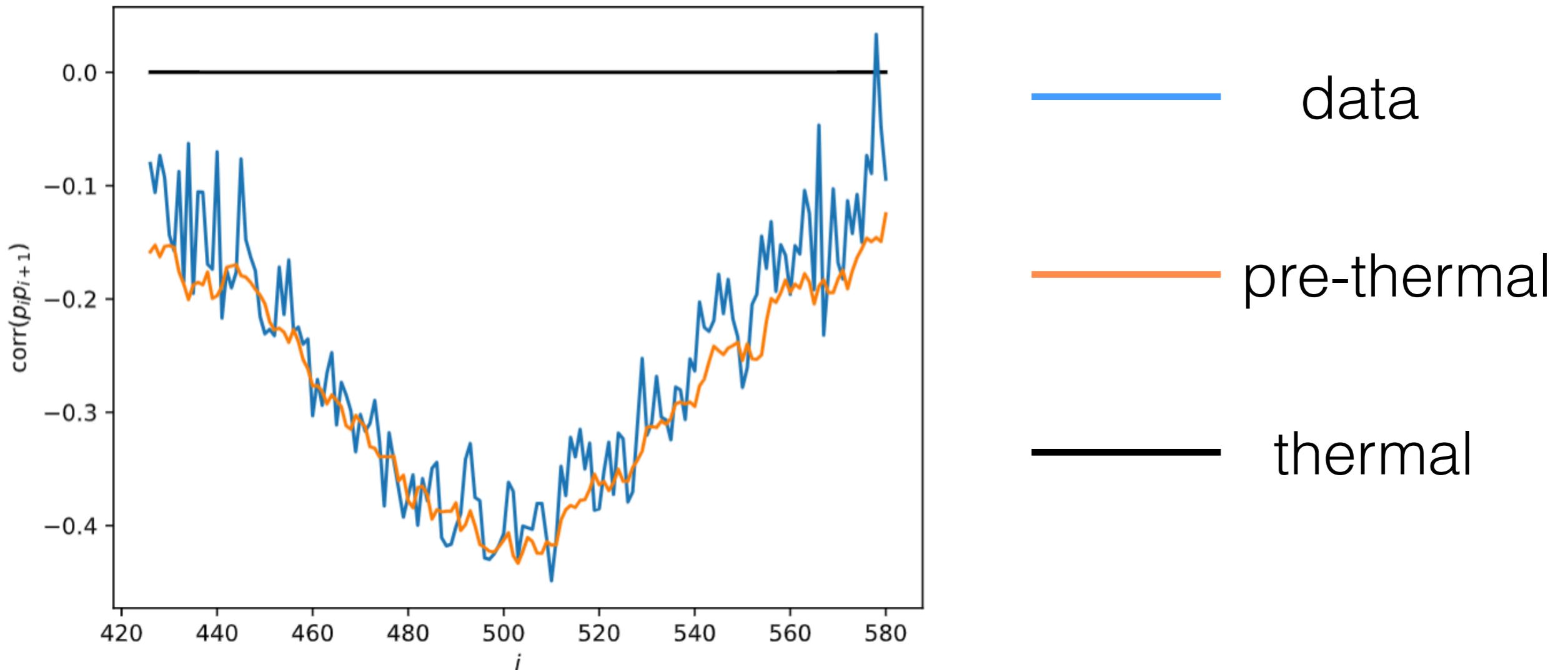


Figure 1. Expected time evolution of a local observable $A(t)$.

H conserved, N pseudo-conserved (number of phonons)

Pre-thermal state in the packet?

Preliminary data suggest the packet is pre-thermal:



It is probably the closest to equilibrium that we can get on these time scales

Challenges with the proof

Perturbative analysis in λ : $\forall n$, find u_n and g_n such that

$$\frac{dH_E}{dt} = \lambda\{H, u_n\} + \lambda^n g_n$$

Hence

$$H_E(t) - H_E(0) = \lambda(u_n(t) - u_n(0)) + \lambda^n \int_0^t ds g_n(s)$$



fluctuation



'dissipation'

Controlling denominators

The perturbative expansion yields *small denominators*:

$$\frac{1}{\sigma_1\nu_1 + \cdots + \sigma_m\nu_m}$$

with

$$\sigma_k = \pm 1$$

ν_k eigenfrequencies $\Leftrightarrow \nu_k^2$ eigenvalues of $H = V - g\Delta$

Heuristics: ν_1, \dots, ν_m are nearly i.i.d.

Controlling denominators

- 2 eigenvalues in a system of size L : Minami's estimate

$$P(\exists \nu_k \neq \nu_{k'} : |\nu_k^2 - \nu_{k'}^2| \leq \gamma) \leq L^2 \gamma$$

- Linear combination of > 2 eigenvalues : ?

$$\nu_{k_1}^2 + \nu_{k_2}^2 - \nu_{k_3}^2 - \nu_{k_4}^2$$

(would feature if KG would be replaced by DNLS)

- Linear combination of > 2 eigenfrequencies : New bound!

$$\nu_{k_1} + \nu_{k_2} - \nu_{k_3} - \nu_{k_4}$$

Trick to control denominators

Shift the full spectrum:

$$H \rightarrow H + \alpha \text{Id}$$

- Leaves *invariant*: $\nu_{k_1}^2 + \nu_{k_2}^2 - \nu_{k_3}^2 - \nu_{k_4}^2$
- Does *not* leave *invariant*: $\nu_{k_1} + \nu_{k_2} - \nu_{k_3} - \nu_{k_4}$

In our model, the disorder is on the diagonal:

$$H = V - g\Delta, \quad V_x = \omega_x^2 \quad \text{i.i.d.}$$

So, we can escape resonances by shifting the whole disorder

Control on denominators

This idea yields a lemma:

Lemma (WDR, FH, OP).

In a system of size L , for any $0 < \varepsilon < 1/L$,

$$P\left(\min \left| \sum_{k=1}^m \tau_k \nu_k \right| \leq \varepsilon\right) \leq C_m L^m \varepsilon^{\frac{1}{m+1}}$$

where the minimum runs over m -tuples of all different eigenvalues, and where $\tau_k \in \{-m, \dots, m\}$, $\tau_k \neq 0$ are given.

From a technical point of view, this is the key new result

Conclusion and outlook

- Our mathematical results show that the chain is asymptotically many-body localized: dissipative effects arise as a non-analytic function of λ .
- Comparison with numerical data suggest that state-of-the-art numerics do not capture correctly the asymptotic behavior of the chain.
- Mathematical results are so far limited to the dynamics in equilibrium. We are working to relax this hypothesis.