

The 2PI-loop expansion revisited

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in collaboration with

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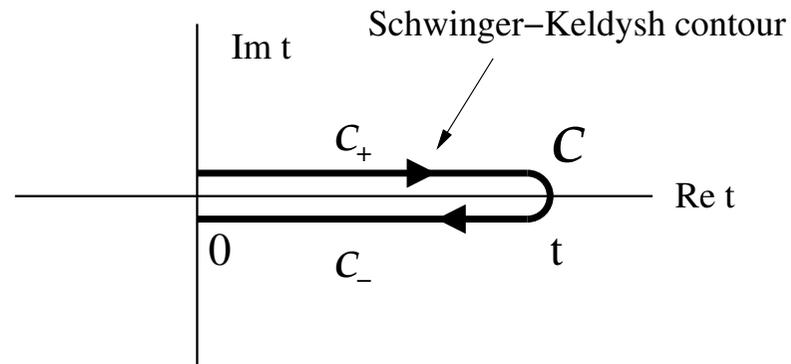
hep-ph/0503287 → Phys.Rev.D

Quantum fields out of equilibrium

Out of equilibrium, the density matrix and operator expectation values **evolve in time**, ($\mathcal{Z} = \text{Tr}[\rho] = 1$)

$$\langle \mathcal{O}(t) \rangle_\rho = \text{Tr}[\rho(t) \mathcal{O}] = \text{Tr}[\rho(0) e^{-iHt} \mathcal{O} e^{iHt}].$$

We can write this as a path-integral, as an evolution along a **path** in complex t -space,



$$\mathcal{Z}[\rho, \mathcal{C}, J_i, K] = \int \mathcal{D}\phi_+ \mathcal{D}\phi_- \langle \phi_+ | \rho(0) | \phi_- \rangle$$

$$\times \int \mathcal{D}\phi \exp \left\{ i \left(S_{\mathcal{C}}[\phi] + J\phi + \frac{1}{2}\phi K\phi \right) \right\},$$

$$S_{\mathcal{C}}[\phi] + J\phi + \frac{1}{2}\phi K\phi = \int_{\mathcal{C}} dt \int d^3x [\mathcal{L}(\phi) + J_{\pm}\phi + \frac{1}{2}\phi K\phi].$$

The currents J_{\pm} and K live on upper (+)/lower (-) branches, and $\phi(\mathbf{x}, 0_+) = \phi_+, \phi(\mathbf{x}, 0_-) = \phi_-$. The **initial condition** in $\langle \phi_+ | \rho(0) | \phi_- \rangle$ can be included in the currents.

Approximate quantum dynamics

Φ -derivable approximations defined in terms of truncated loop expansions of the **2PI effective action** provide a framework for approximating the **quantum dynamics** (in and) out of equilibrium.

Equations of motion are derived in terms of the **1-** and **2-point functions** (the **mean field** $\bar{\phi}$ and the **propagator** G). The derived equations of motion have to be **solved numerically**, by discretizing the field on a **lattice**.

[Berges&Cox:2000, Aarts&Berges:2001, Berges:2002, Aarts et al.:2002, Cassing et al.:2003, Berges et al.:2004], thermalization for scalar fields in 1+1, 2+1 and 3+1 dimensions, [Berges et al.:2003] fermions and scalars in 3+1 dimensions. [Berges&Serreau:2003, Arrizabalaga et al.(AT):2004] for studies of preheating. For gauge fields, see [Arrizabalaga&Smit:2003, Berges:2004]. [van Hees&Knoll:2002, Blaizot et al.:2003, Berges et al.:2005], renormalization].

2PI effective action

$$Z[J, K] = \int d\phi e^{iS[\phi] + iJ\phi + \frac{i}{2}\phi K\phi} = e^{iW[J, K]},$$

$$\bar{\phi}(x) = \frac{\delta W[J, K]}{\delta J(x)}, \quad G(x, y) = 2 \frac{\delta W[J, K]}{\delta K(x, y)} - \bar{\phi}(x)\bar{\phi}(y).$$

Define:

$$\Gamma^{2PI}[\bar{\phi}, G] = W[J] - J\bar{\phi} - \frac{1}{2}K(G + \bar{\phi}\bar{\phi}).$$

Then:

$$\frac{\delta \Gamma^{2PI}[\bar{\phi}, G]}{\delta \bar{\phi}(x)} = -J(x) - \int d^4y K(x, y)\bar{\phi}(y),$$
$$\frac{\delta \Gamma^{2PI}[\bar{\phi}, G]}{\delta G(x, y)} = -\frac{1}{2}K(x, y).$$

Φ -derivable approximations

The **2PI** effective action can be written as [Cornwall et al :1974]:

$$\Gamma^{2PI}[\bar{\phi}, G] = S[\bar{\phi}] - \frac{i}{2} \text{Tr} \ln G + \frac{i}{2} \text{Tr} G_0^{-1} G + \Phi[\bar{\phi}, G],$$

with

$$iG_0^{-1}(x, y) = \frac{\delta^2 S[\bar{\phi}]}{\delta\bar{\phi}(x)\delta\bar{\phi}(y)} \delta(x - y).$$

Φ is the sum of all **2PI skeleton diagrams** with **no external lines**, **full propagators** and **bare vertices**. A choice of truncation of this series $\Phi_{\text{tr}} \rightarrow \Gamma_{\text{tr}}^{2PI}$ constitutes a **Φ -derived approximation**.

Equations of motion derived from any truncation

$$\frac{\delta\Gamma_{\text{tr}}^{2PI}[\bar{\phi}, G]}{\delta G(x, y)} = 0, \quad \frac{\delta\Gamma_{\text{tr}}^{2PI}[\bar{\phi}, G]}{\delta\bar{\phi}(x)} = 0$$

conserve an energy functional.

Equations of motion

The equations of motion for a **homogeneous system** :

$$\begin{aligned}\partial_t^2 F(t, t', \mathbf{x}) &= \partial_i^2 F(t, t', \mathbf{x}) - M_{\text{eff}}^2(t) F(t, t', \mathbf{x}) \\ &+ \int_0^t dt'' d\mathbf{y} \Sigma_\rho(t, t'', \mathbf{y}) F(t'', t', \mathbf{x} - \mathbf{y}) \\ &- \int_0^{t'} dt'' d\mathbf{y} \Sigma_F(t, t'', \mathbf{y}) \rho(t'', t', \mathbf{x} - \mathbf{y}), \\ \partial_t^2 \rho(t, t', \mathbf{x}) &= \partial_i^2 \rho(t, t', \mathbf{x}) - M_{\text{eff}}^2(t) \rho(t, t', \mathbf{x}) \\ &+ \int_{t'}^t dt'' d\mathbf{y} \Sigma_F(t, t'', \mathbf{y}) \rho(t'', t', \mathbf{x} - \mathbf{y}), \\ \partial_t^2 \bar{\phi}(t) &= - M_{\text{eff}}^2(t) \bar{\phi}(t) - \int_0^t dt'' d\mathbf{y} \Sigma_\phi(t, t'', \mathbf{y}) \bar{\phi}(t'').\end{aligned}$$

with $F = \langle \{ \phi, \phi \} \rangle$, $\rho = i \langle [\phi, \phi] \rangle$. These are the **Kadanoff-Baym equations** (if non-relativistic) [**Kadanoff&Baym:1962**].

ϕ^4 -model

We study a single **real scalar field** in a ϕ^4 potential:

$$S = - \int dt \int d^3x \left[\frac{\partial_\mu \phi \partial^\mu \phi}{2} + \frac{m^2}{2} \phi^2 + \frac{\lambda}{24} \phi^4 \right].$$

In the **symmetric phase** $m^2 > 0$, $\langle \phi \rangle \simeq 0$ and in the **broken phase** $m^2 < 0$, $\langle \phi \rangle \simeq \sqrt{6|m^2|/\lambda}$.

The functional Φ

The functional Φ in a **loop expansion** is

$$\begin{aligned} \Phi[\bar{\phi}, G] = & \frac{1}{4!} \text{diagram}_1 + \frac{1}{2} \text{diagram}_2 + \frac{1}{8} \text{diagram}_3 \\ & + \frac{1}{12} \text{diagram}_4 + \frac{1}{48} \text{diagram}_5 + \dots \end{aligned}$$

Motivation

- **Thermalization** for scalar fields in $3+1d$.
- Most 2PI studies compare $O(\lambda)$ (**Hartree**) to $O(\lambda^2)$ (**Basketball**) or **LO-1/N** to **NLO-1/N**. $\langle\phi\rangle = 0$.
- Hartree/LO-1/N: **Leading** contribution to $\text{Re}\Sigma$.
Basketball/NLO-1/N: **Next-to-Leading** contribution to $\text{Re}\Sigma$.
- Basketball/NLO-1/N: **Leading** contribution to $\text{Im}\Sigma$. **No results** exist at **NNLO**. Numerically **Intractable**(?).
- The $\text{Im}\Sigma$ is related to **decay/damping rates, equilibration times**.
- $\langle\phi\rangle \neq 0$ induces **three point vertex**, and an **additional diagram** at $O(\lambda)$. Perturbatively, $\text{Im}\Sigma = 0$ (on shell). **2PI-resummed**?
- Three point vertex present in **non-abelian gauge theories**.

Initial conditions

We apply two **initialization schemes**, Top-hat and Thermal, determined by a choice for the **distribution function** n_k and the **dispersion relation** ω_k , through:

$$\langle \phi_k \phi_{-k} \rangle = \frac{n_k + 1/2}{\omega_k}, \quad \langle \pi_k \pi_{-k} \rangle = (n_k + 1/2)\omega_k,$$

- **Top-hat:**

$$n_k = 0; \quad n_k = c > 0, \quad k_{\min} < k < k_{\max},$$

for some $c > 0$.

- **Thermal:**

$$n_k = (e^{\omega_k/T} - 1)^{-1}.$$

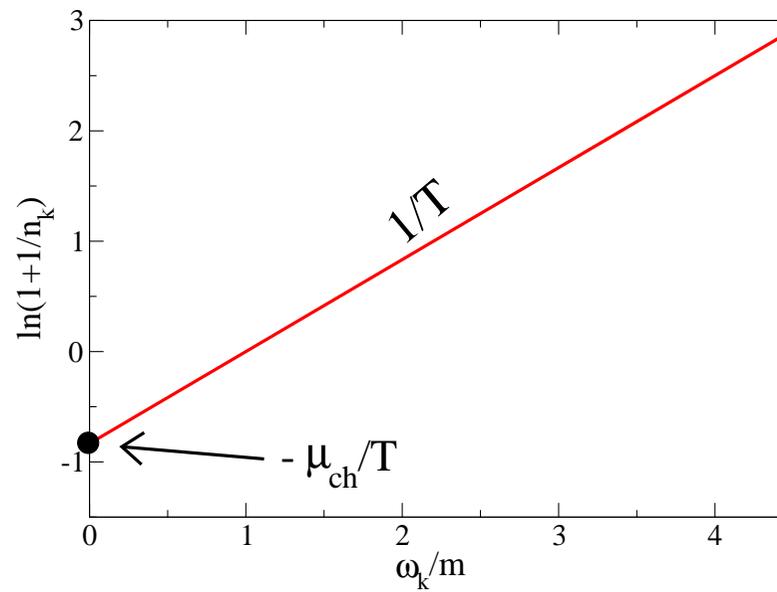
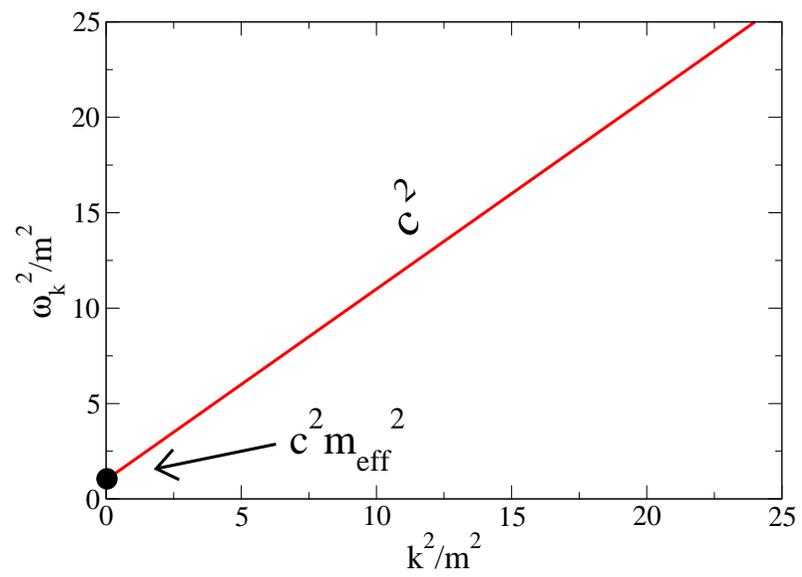
In both cases, $\omega_k^2 = k^2 + m^2$.

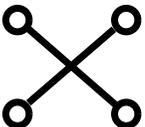
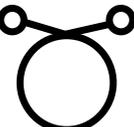
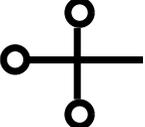
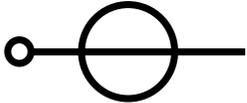
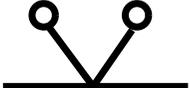
What equilibrium looks like

For small coupling we expect the equilibrium state to have a **quasi-particle** dispersion relation and a **Bose-Einstein** particle distribution. We will define (also out of equilibrium)

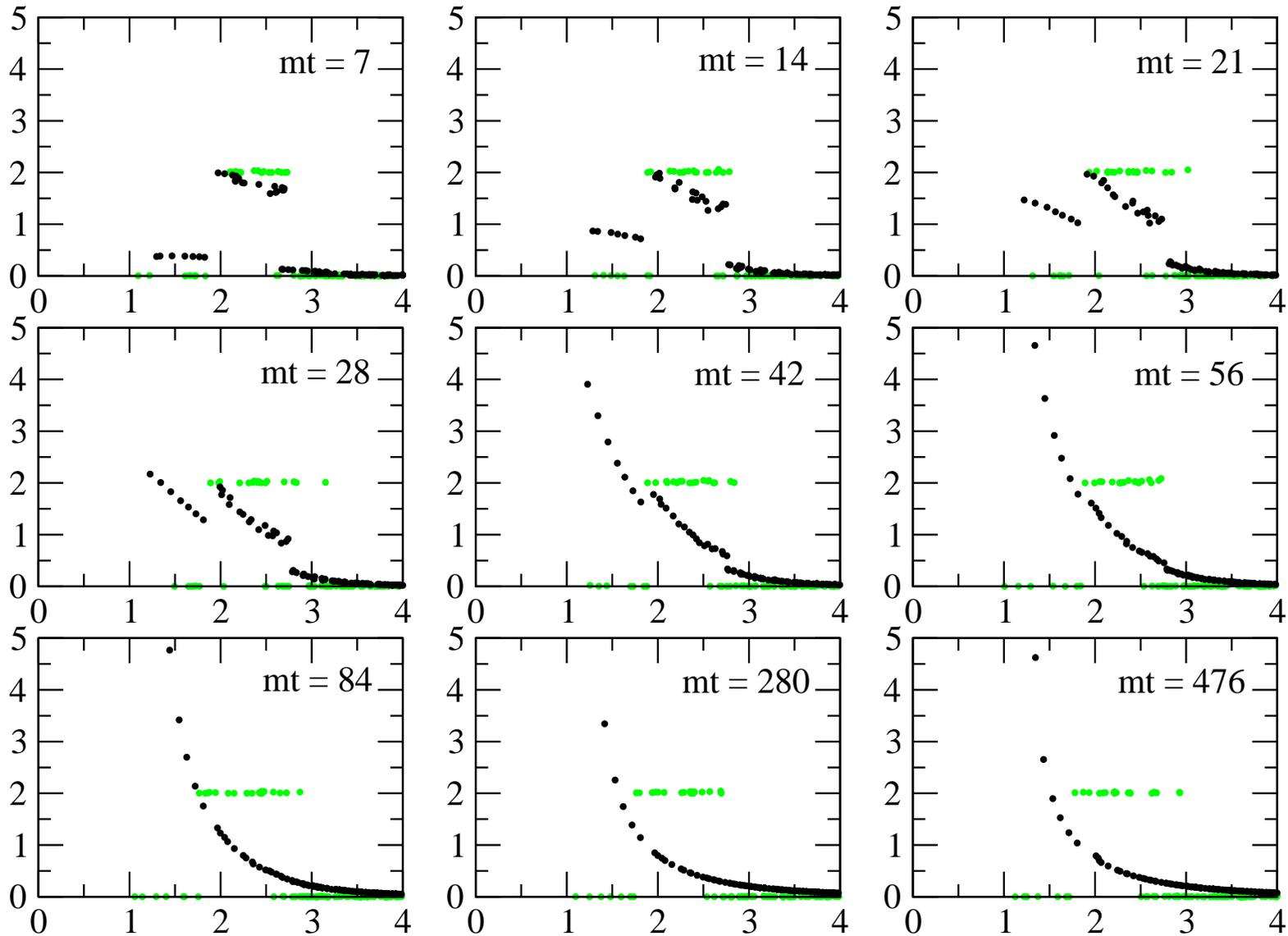
$$\begin{aligned}\omega_k &= \sqrt{\langle \pi_k \pi_{-k} \rangle / \langle \phi_k \phi_{-k} \rangle}, \\ n_k + 1/2 &= \sqrt{\langle \pi_k \pi_{-k} \rangle \langle \phi_k \phi_{-k} \rangle}.\end{aligned}$$

When plotting ω_k^2 vs. k^2 the **slope** is c^2 and the **intercept** $c^2 m_{\text{eff}}^2(T)$. In a plot of $\ln(1 + 1/n_k)$ vs. ω_k , the **slope** is T^{-1} and the **intercept** $-\mu_{\text{ch}}/T$.

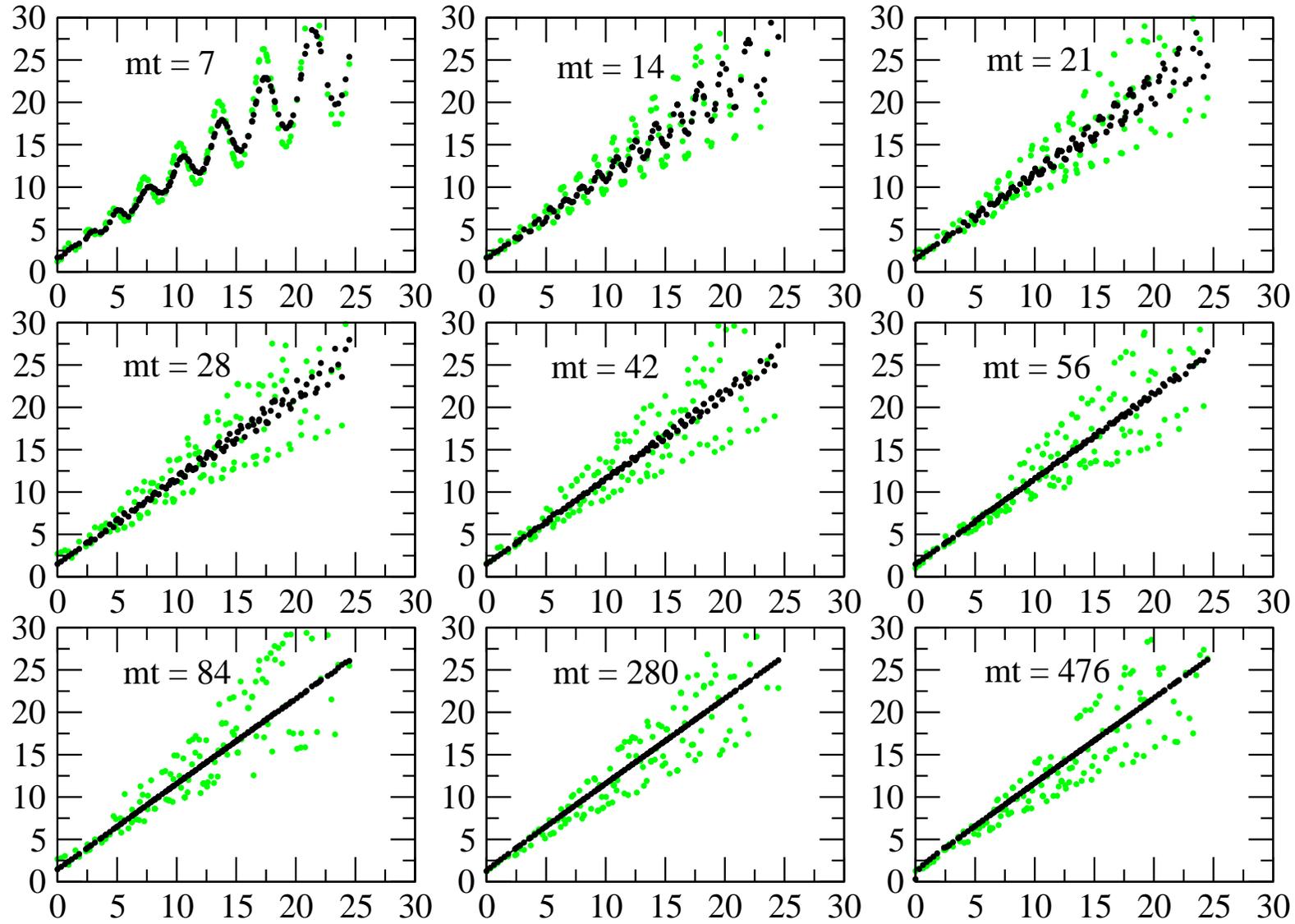


$\Phi =$					
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$\phi = \lambda^{-1/2}$	$O(\lambda^{-1})$	$O(1)$	$O(\lambda)$	$O(\lambda)$	$O(\lambda^2)$
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	$O(\delta^2, 1)$	$O(\lambda)$	$O(\lambda^2)$		
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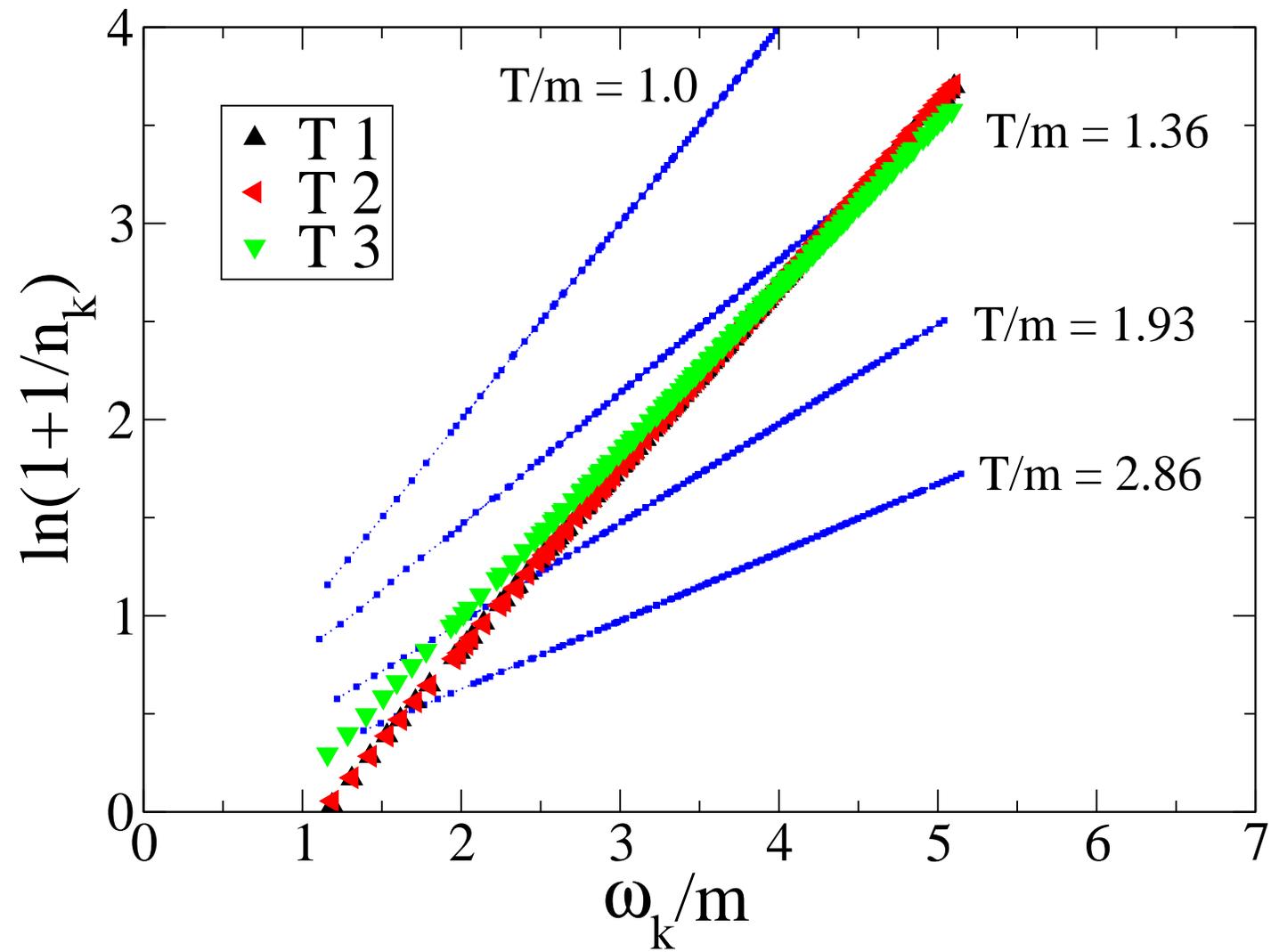
Symmetric phase, including basketball



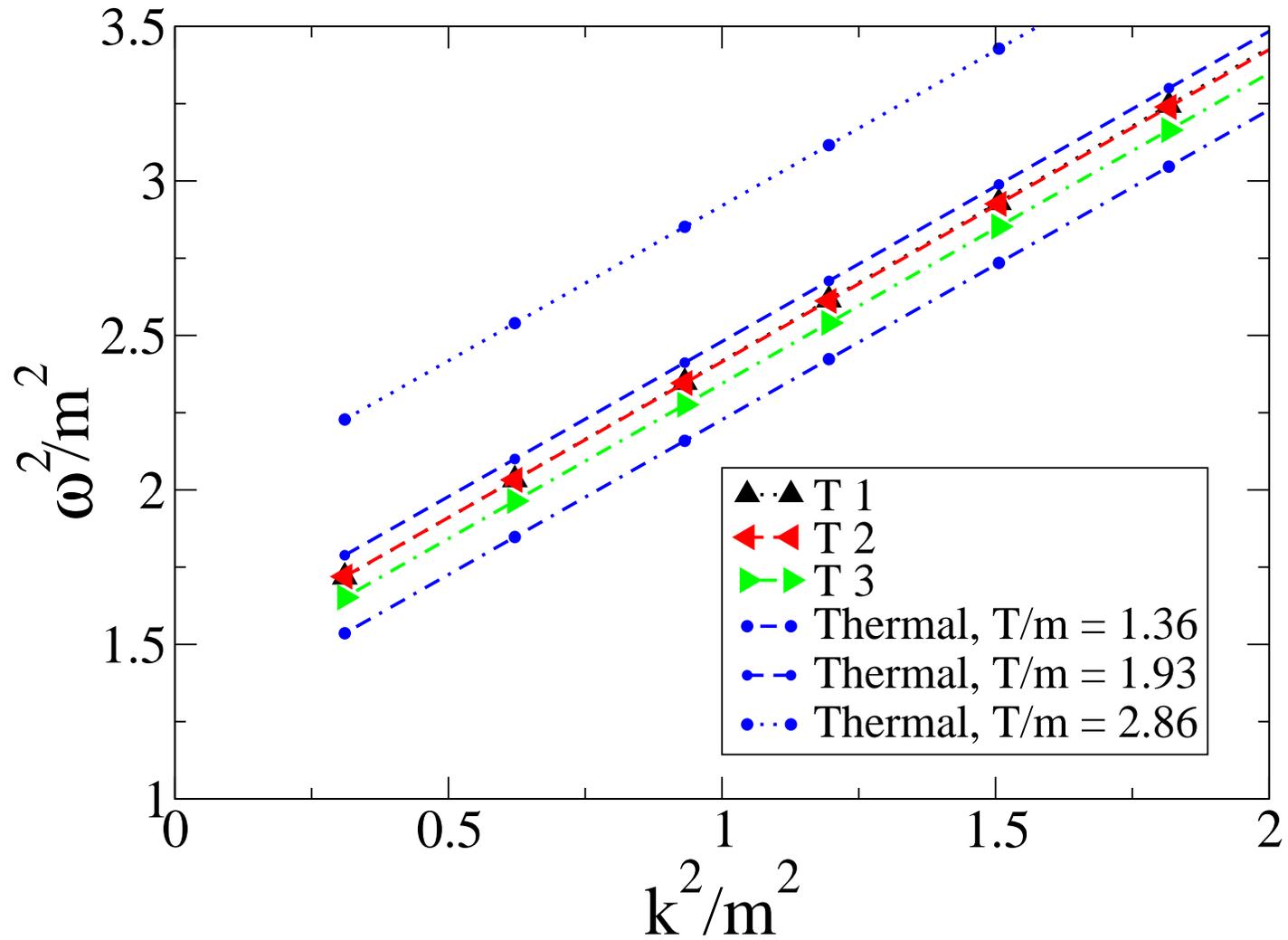
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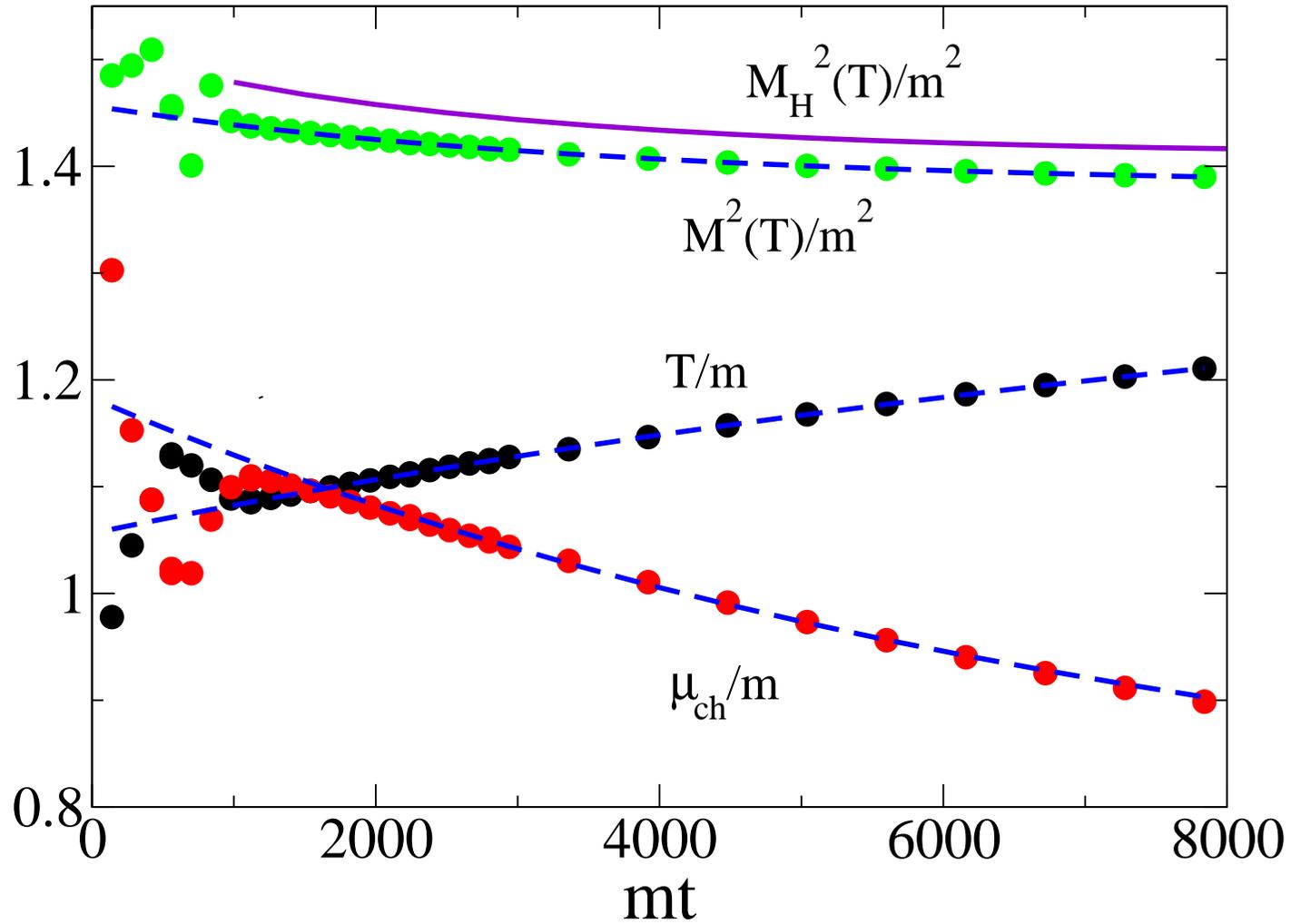
Kinetic equilibration

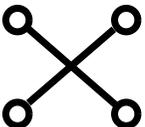
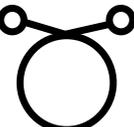
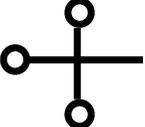
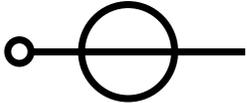
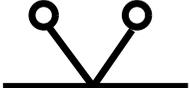


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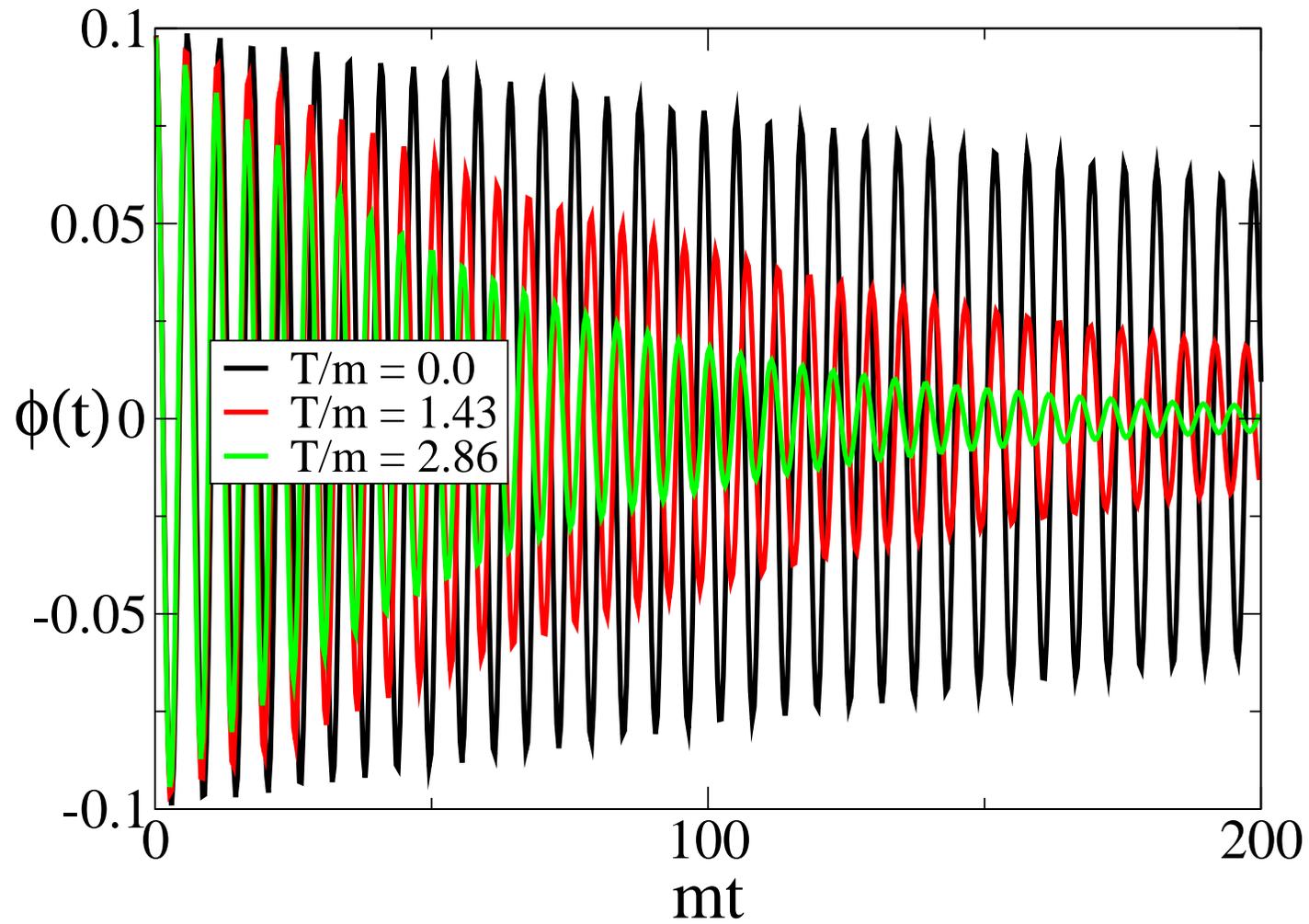


Chemical equilibration

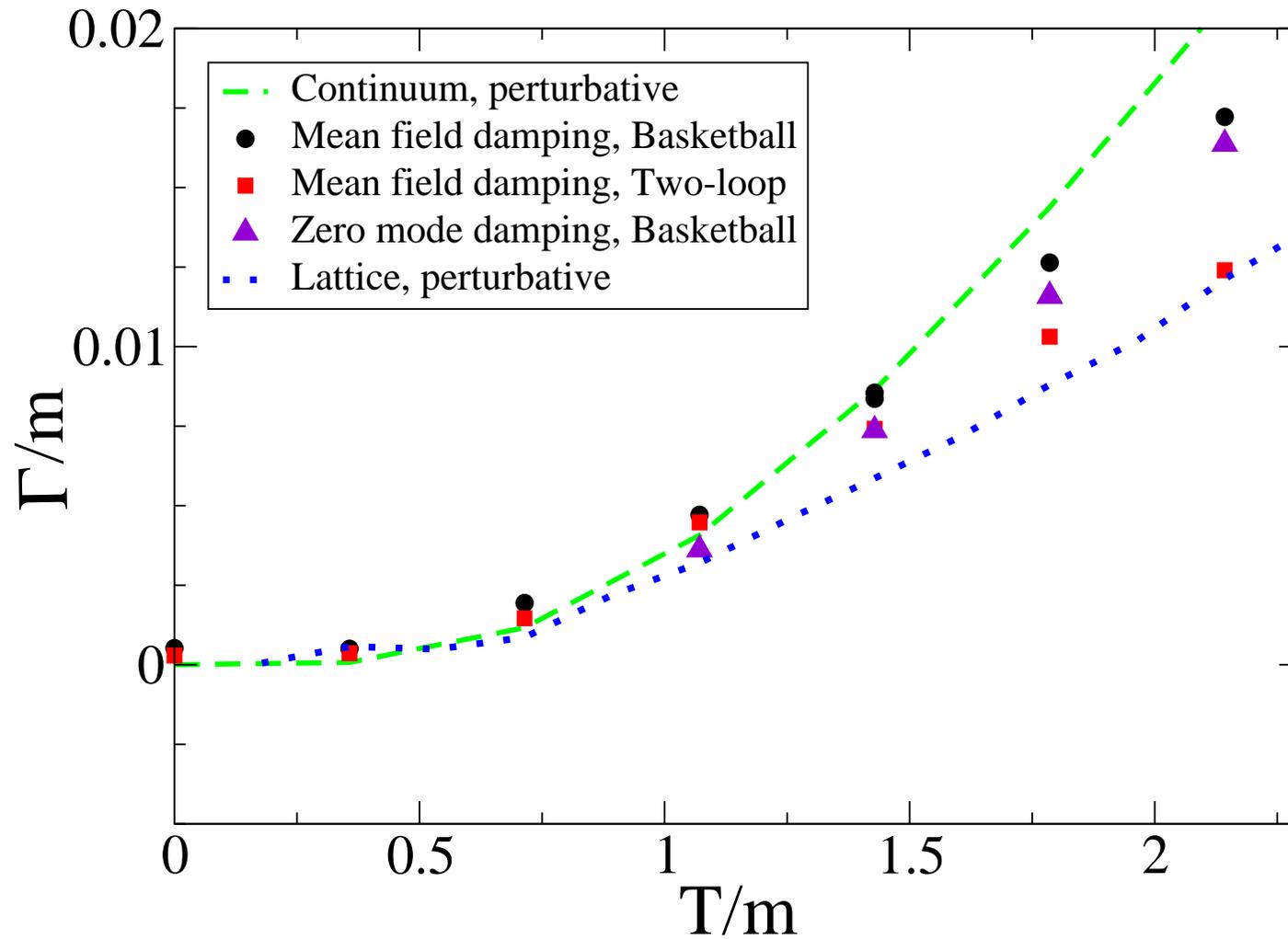


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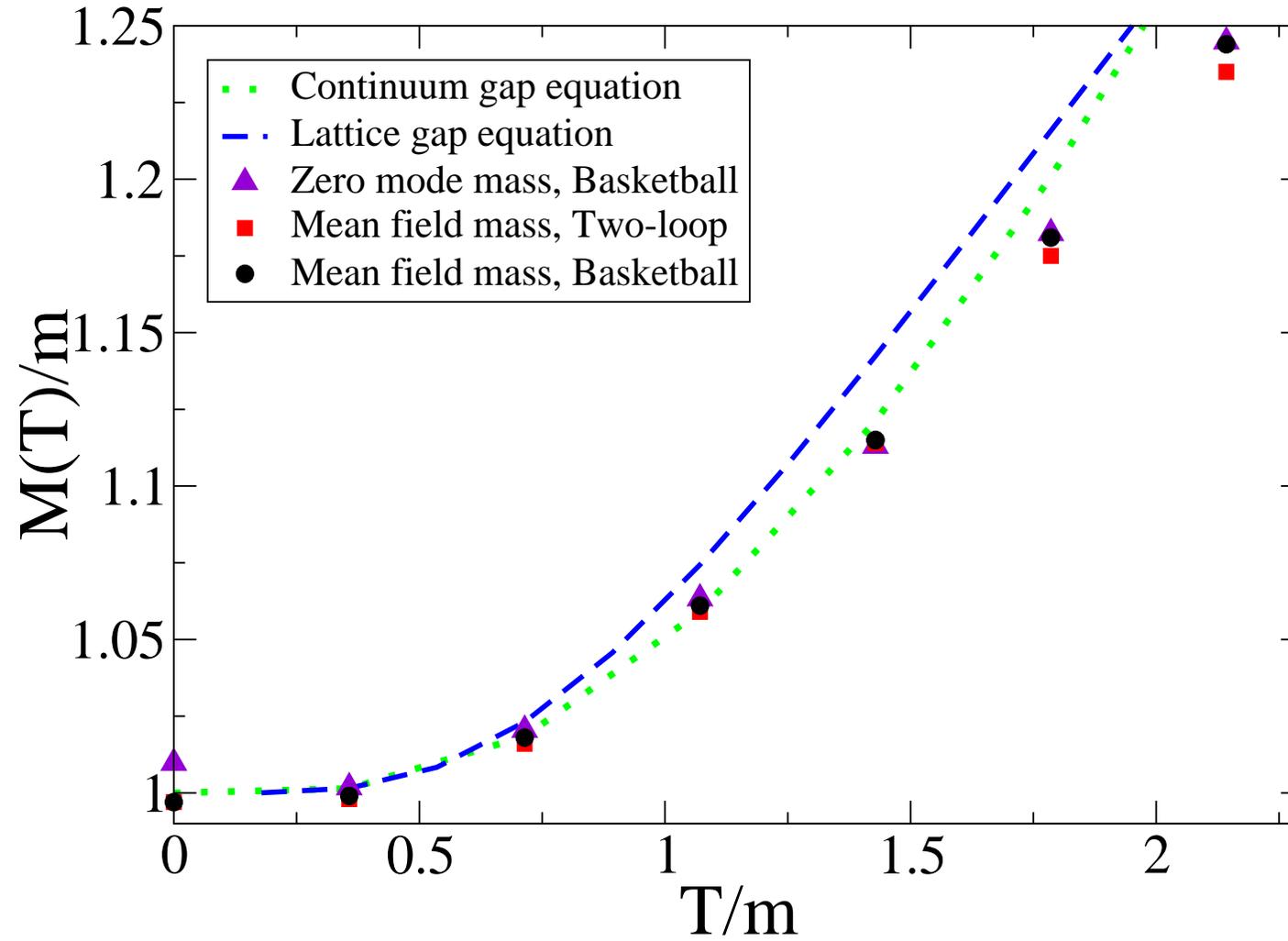
Mean field damping

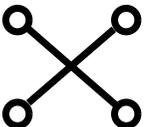
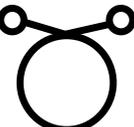
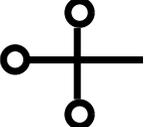
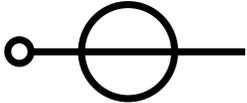
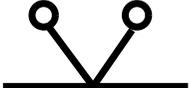
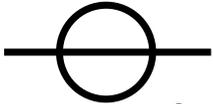


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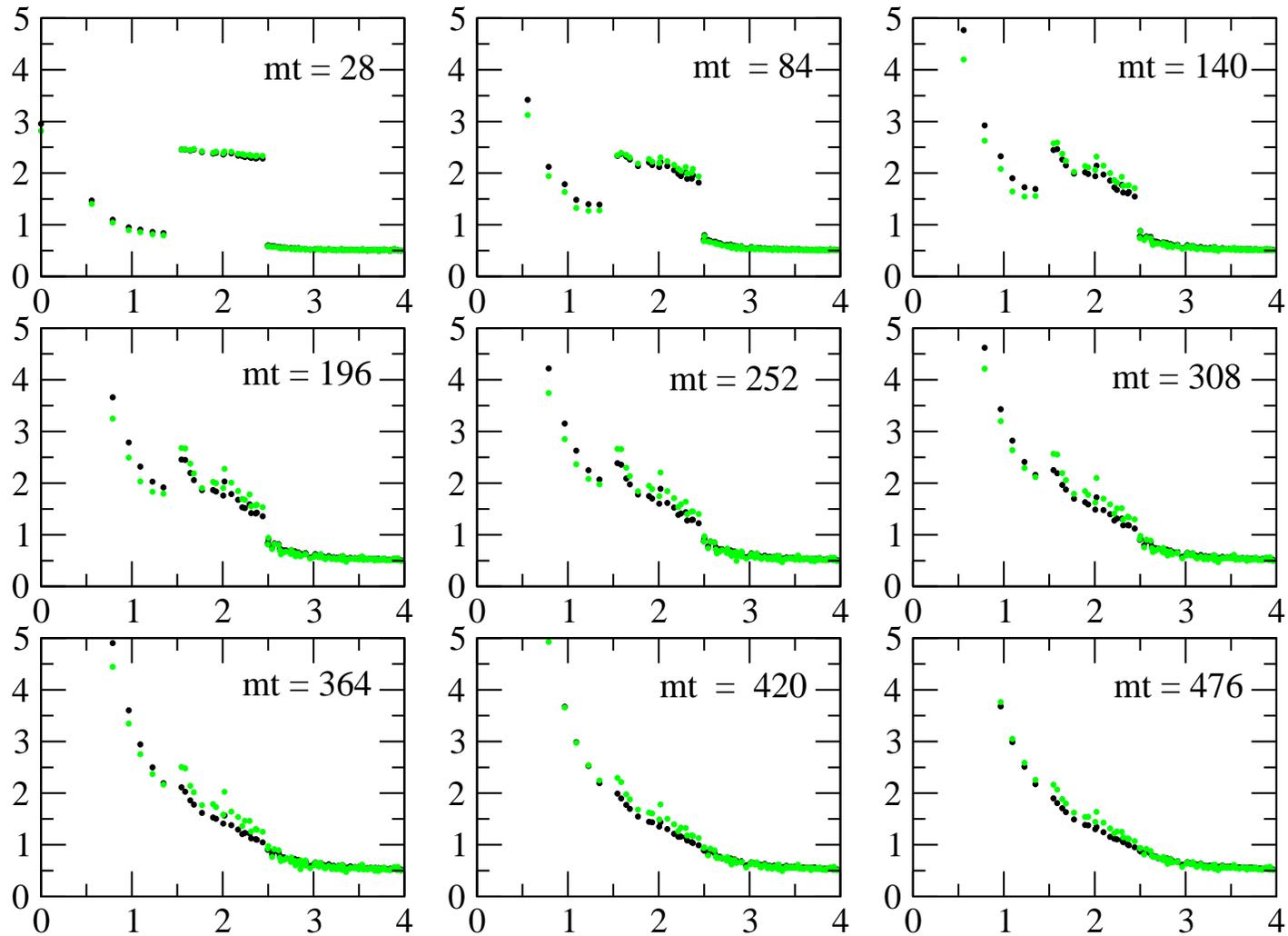


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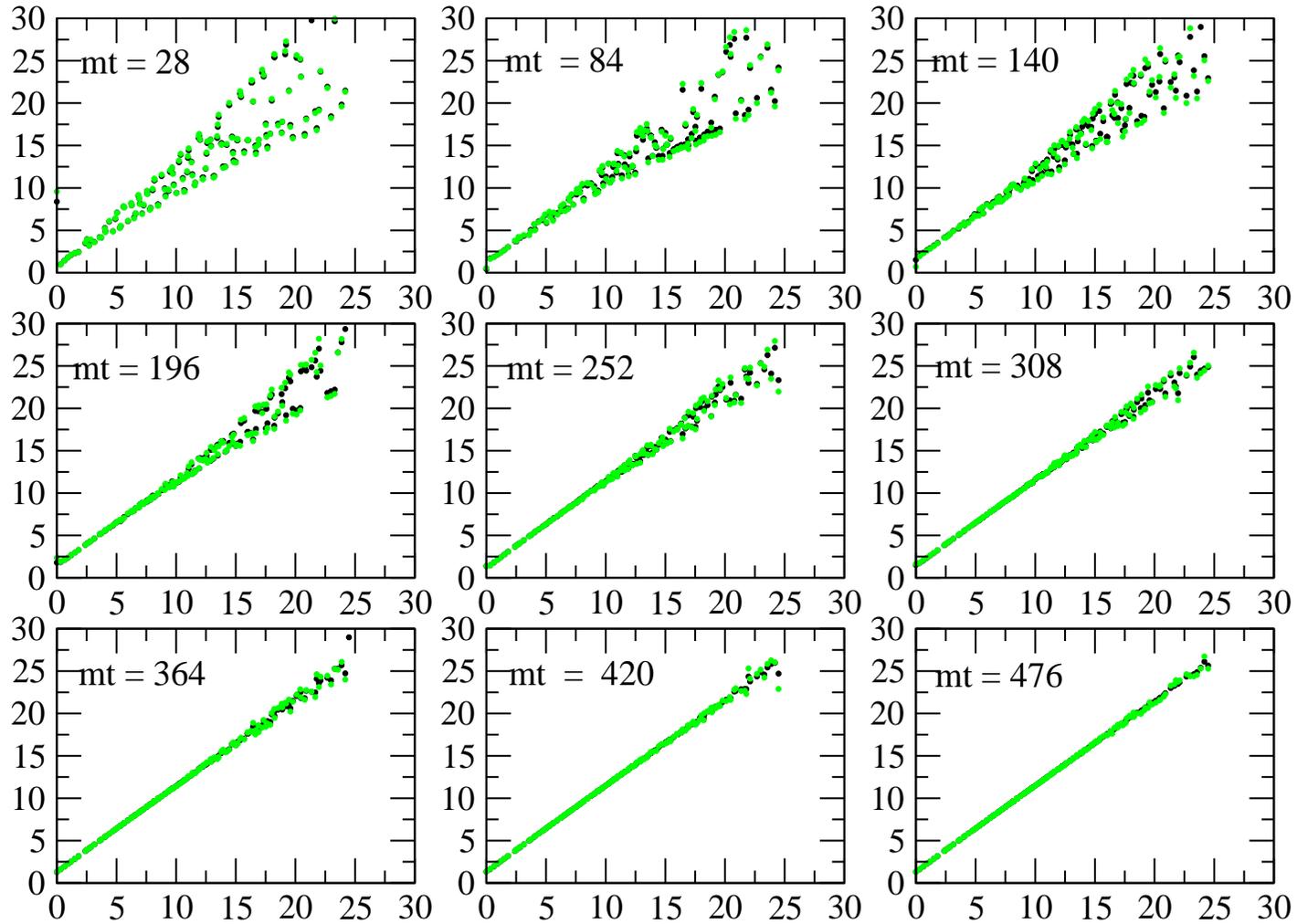


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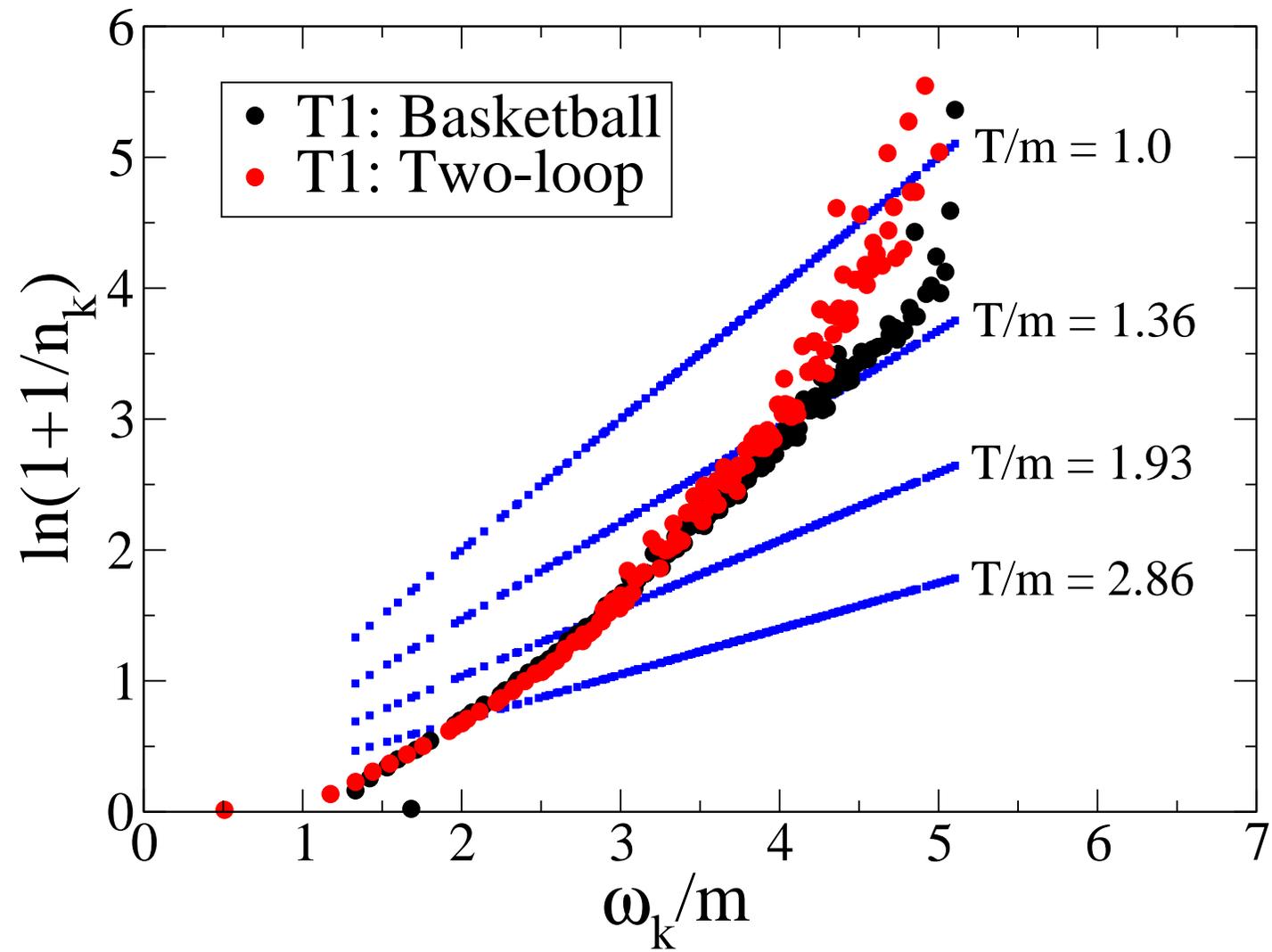
Broken phase



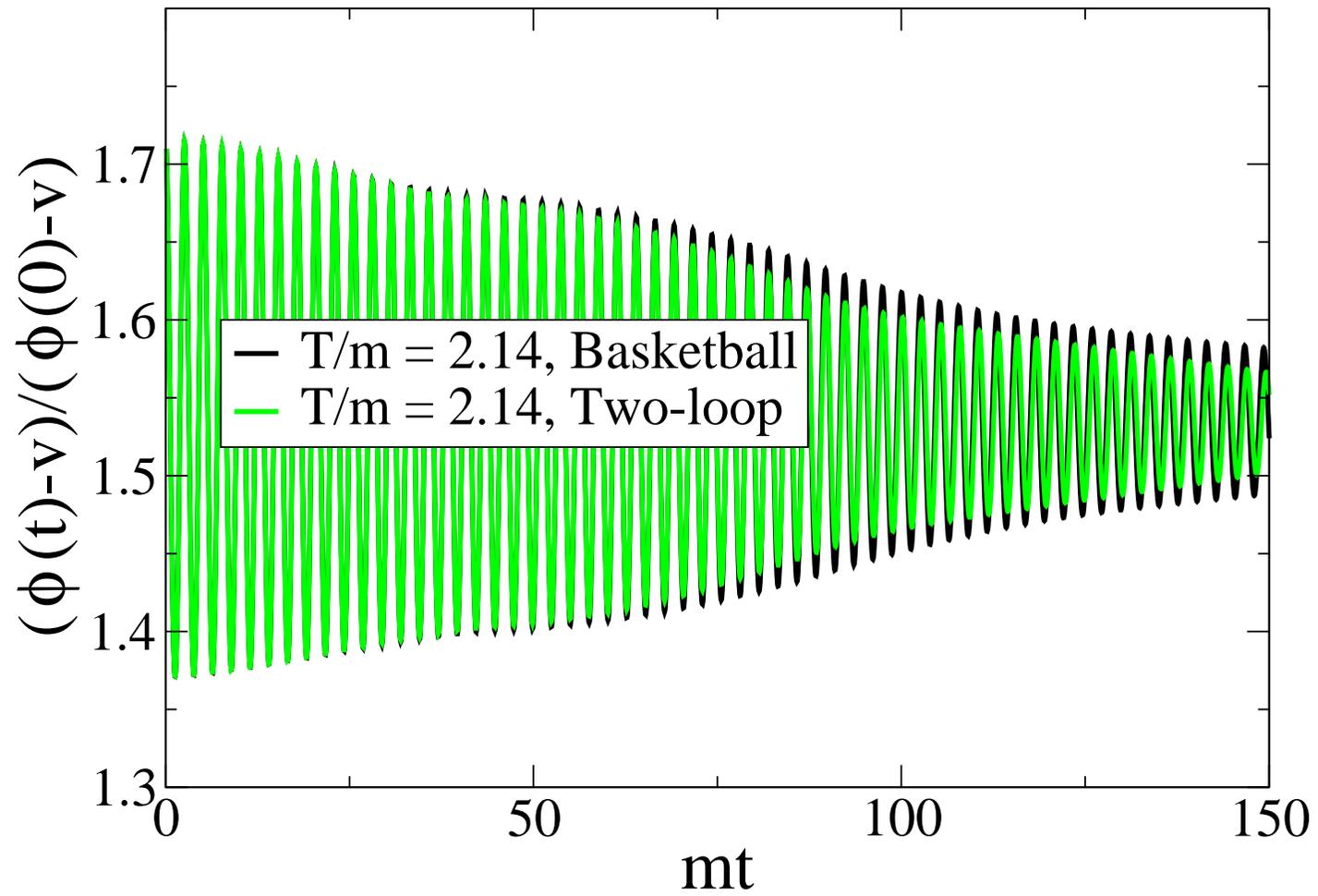
Broken phase



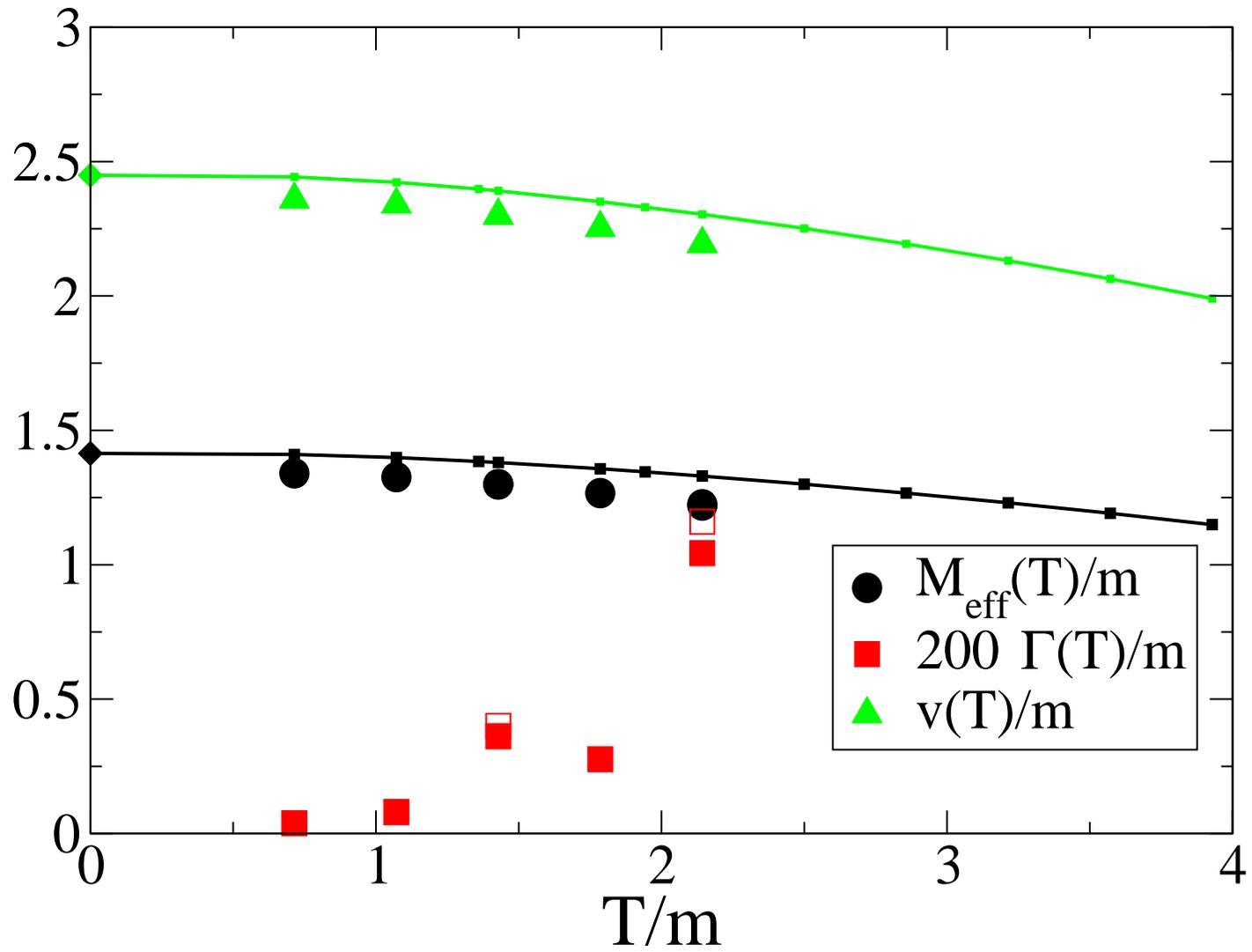
Equilibration



Mean field damping



Mean field damping



Conclusion

- In the **symmetric phase**, $\text{Re}\Sigma(\lambda^2)/\text{Re}\Sigma(\lambda) \simeq 0.3$ at $\lambda = 6$.
Expansion will break down for much larger λ . **Renormalization**.
- $\text{Im}\Sigma(\lambda^2)$ is **slightly larger** than **non-resummed** analytical calculations.
- In the **broken phase**, resummation introduces $\text{Im}\Sigma \neq 0$ at $O(\lambda)$.
Qualitatively different from perturbation theory.
- **Equilibration time**: $O(\lambda^2)/O(\lambda) \sim O(1)$.
- $O(\lambda)$ does not dominate $O(\lambda^2)$. Probably because first level of resummation is $O(\lambda^4 v^4) = O(\lambda^2)$.