

For this afternoon:

... if you're participating

- Please download GABE
 - cosmo.kenyon.edu/gabe.html
- If you have a MAC
 - Try to run the instructions on that page, new(ish) operating systems should work without issue
- If you have LINUX
 - Please install fftw3 (enabling regular and long-double: openmp and threads for each)
- If you plan to use the remote version, please email me (giblinj@kenyon.edu) as soon as possible!
- **Please also have Jupyter (or similar) for plotting. I will distribute a .ipynb this afternoon!**



Kenyon College

Numerical Preheating

Thoughts and Exercises: Morning

Tom Giblin

RIKEN Interdisciplinary Theoretical and Mathematical Sciences Program

Saitama, Japan

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My goal

- I'm not going to try to sell you anything (although if you want to buy it...)
- There are some (beautiful) subtleties associated with doing nonlinear dynamics
 - This will not be an exhaustive list
- **Please talk with us (all of the people who do numerical work)**
 - We're (always) happy to share ideas on figuring out whether results are *physical* or *numerical*

Simulating Preheating

Has a long history

- Perhaps the first real simulations are from 1996/7
 - S. Yu. Khlebnikov and Igor Tkachev
 - Phys. Rev. Lett. 77, 219 (1996)
- Made 'famous' by LatticeEasy around 2000
 - Gary Felder and Igor Tkachev
 - Comput.Phys.Commun.178:929-932.2008
- LatticeEasy opened the door to (any of us) to look at nonlinear dynamics of *any* inflationary model

There are many (many) codes

...many of which are open-source or available

- LatticeEasy, Gary Elder 2000 (<https://www.felderbooks.com/latticeeasy/index.html>)
 - CLUSTEREasy, arXiv:0712.0813 (<https://www.felderbooks.com/latticeeasy/index.html>)
- DEFrost, Andrei Frolov, 2008, arXiv:0809.4904, (<https://www.sfu.ca/physics/cosmology/defrost/>)
- CUDAEasy, Jani Sanio, 2009, arXiv:0911.5692
- PSpectRe, Richard Easther, Hal Finkle, Nathaniel Roth, arXiv:1005.1921
- HLATTICE, Zhiqi Huang, arXiv: 1102.0227 (<https://www.cita.utoronto.ca/~zqhuang/hlat/>)
- GABE, JTG, Hillary Child, J. Tate Deskins, arXiv:1305.0561, (<https://cosmo.kenyon.edu/gabe.html>)
- PyCOOL, Jani Sainio, arXiv:1201.5029
- CosmoLattice, 2020, Daniel G. Figueroa, Adrien Florio, Francisco Torrenti, Wessel Valkenburg, arXiv:2006.15122, (<https://cosmolattice.net/>)

AND I'm missing some

- This list doesn't include programs
 - (like CACTUS) that were designed for simulating scalar fields in other contexts
 - That were written for Numerical Relativity and can handle scalar fields
 - Etc, etc
 - I apologize in advance for any citations or contributions that I've left off!!

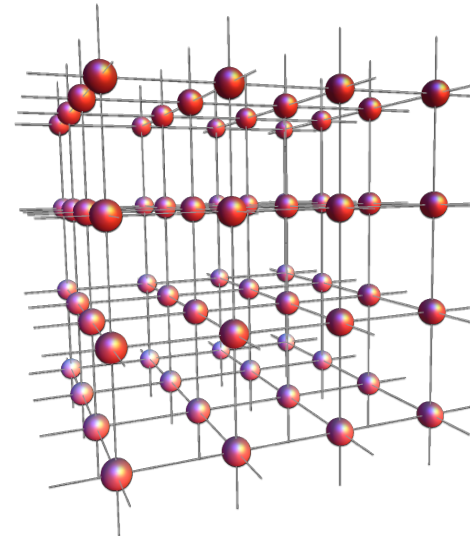
What's the primary take-away?

The discrete system *is* a physical system

...but it's not the same physical system as the continuum



$$\phi(\vec{x})$$



$$\phi(\vec{x}_i)$$

The simplest (and most relevant) example

...the ideal numerical system*

- Consider the wave equation: $\ddot{\phi} - \nabla^2 \phi = 0$

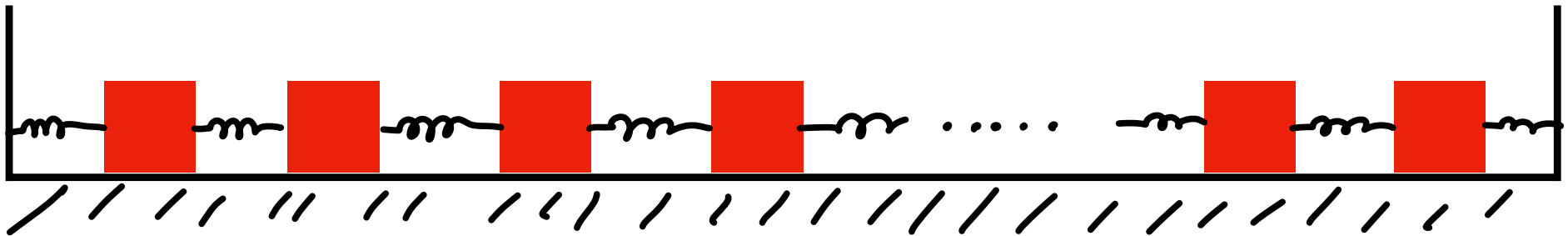
- Which is the coupled system of first-order PDEs: $\dot{\phi} = \omega$
 $\dot{\omega} = \nabla^2 \phi$

- Which are, in the discrete limit, a set of N (or N^3) coupled ODE's**:

$$\dot{\phi}(\vec{x}_i) = \omega(\vec{x}_i) \quad \dot{\omega}(\vec{x}_i) = \frac{1}{\Delta x^2} [\phi(\vec{x}_{i+1}) - 2\phi(\vec{x}_i) + \phi(\vec{x}_{i-1})]$$

*due to its strong hyperbolicity **written in 1-d

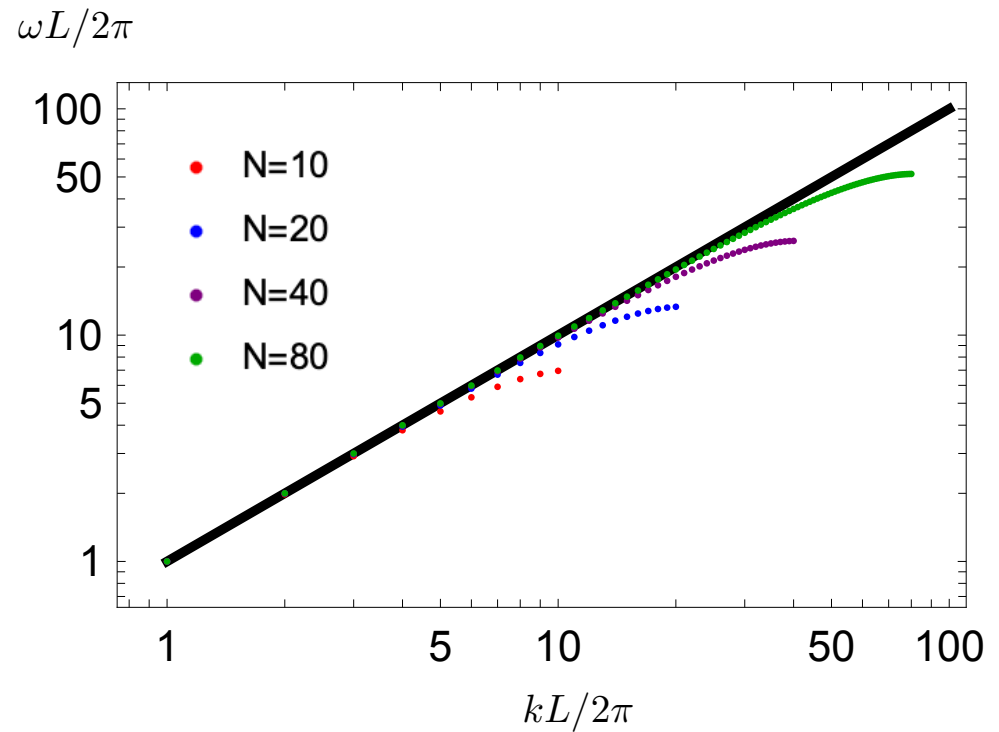
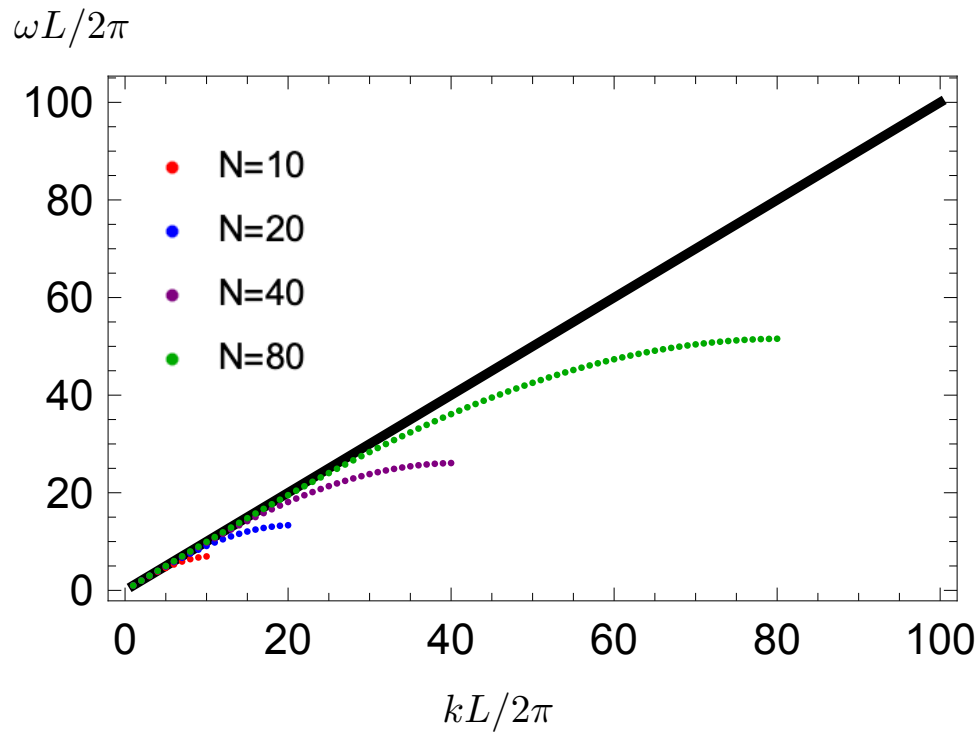
Because you remember blocks and springs



- This system of discrete blocks approximates a continuum system
- However, the frequencies of the normal modes of the system come from find the eigenvalues of the system of equations

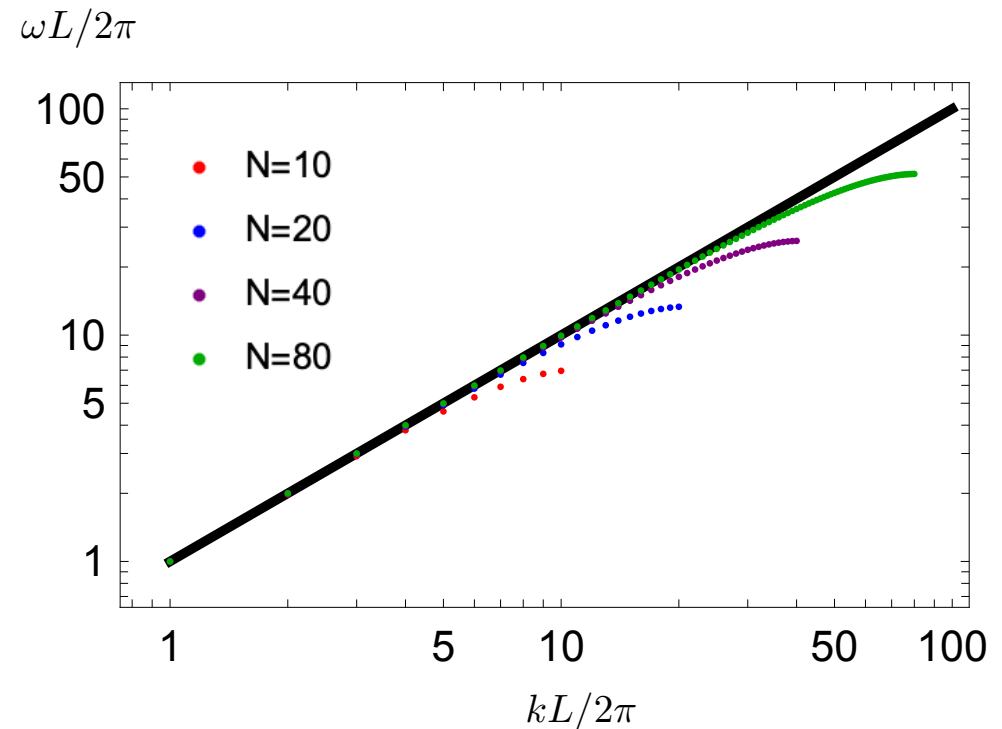
$$\ddot{\vec{y}} = \frac{k}{m} \begin{pmatrix} -2 & 1 & 0 & 0 & & \\ 1 & -2 & 1 & 0 & & \\ 0 & 1 & -2 & 1 & & \\ 0 & 0 & 1 & -2 & & \\ & & & & \ddots & \\ & & & & & \ddots \end{pmatrix} \vec{y}$$

The dispersion relation for this system



The dispersion relation for this system

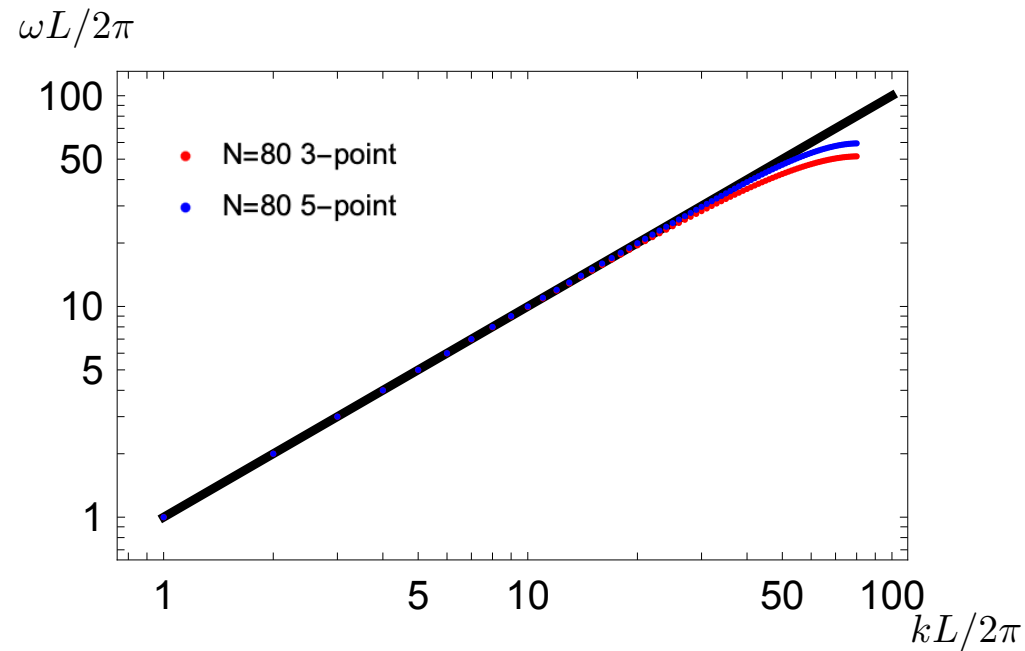
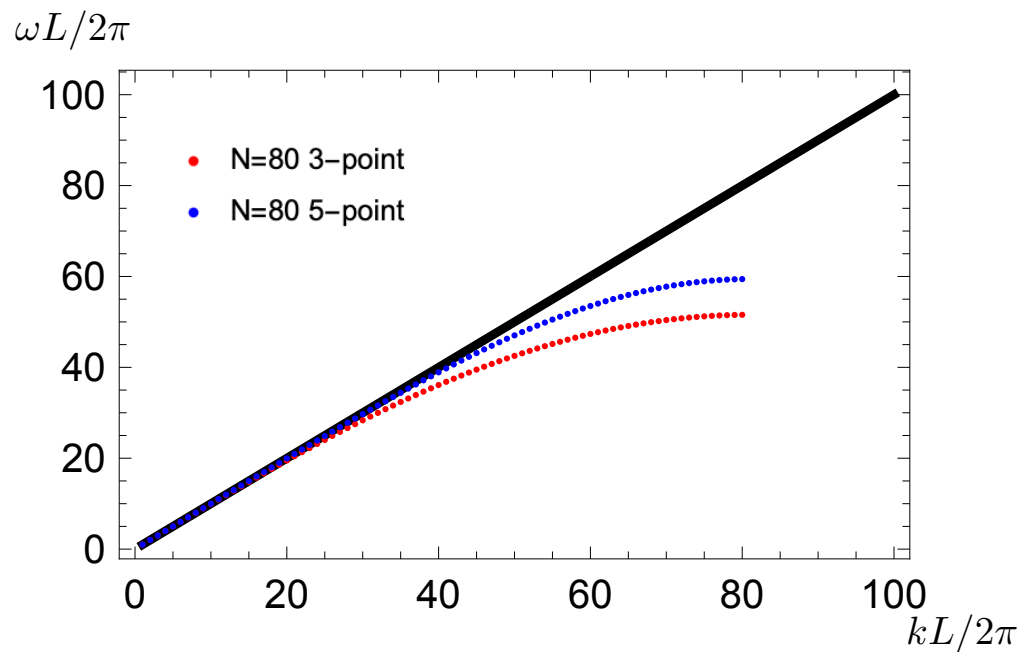
- Remember that the more *points* you have, the larger the *wavenumber* you can resolve!
- The box-size sets the *minimum* (non-zero) wavenumber that you can resolve
- This also means that all wavelengths larger than this are “included” in the zero-bin



You can play some 'tricks'

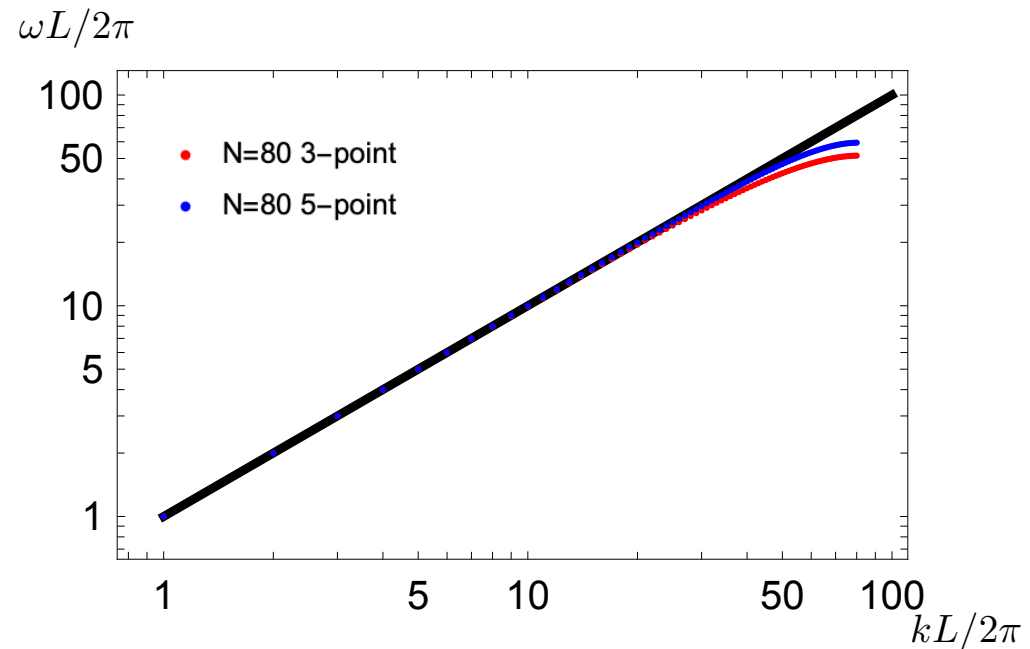
...like stencils

$$\dot{\omega}(\vec{x}_i) = \frac{1}{12\Delta x^2} [-\phi(\vec{x}_{i-2}) + 16\phi(\vec{x}_{i+1}) - 30\phi(\vec{x}_i) + 16\phi(\vec{x}_{i-1}) - \phi(\vec{x}_i)]$$



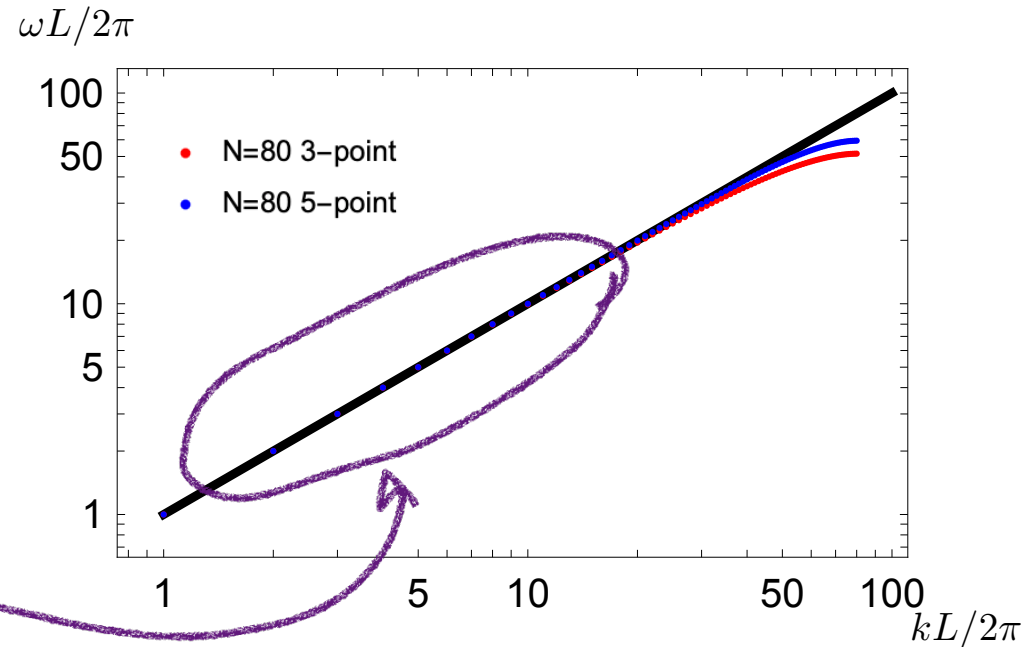
You have to look where you can *trust*

- To quantify how much you “trust” you need to do more sophisticated tests
 - More on these later
- But in general, you plan to have the physics you care about in the lower-half of the log-modes



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A little bit about GABE

Grid and Bubble Evolver (GABE)

- A discretized lattice of N^3 points
- 2nd-order Runge-Kutta Integration scheme
- Natively it handles n -scalar fields
 - Which are treated completely non-linearly
- Gravity is treated homogeneously via the Friedman constraint
- It's “easy” to change potentials
 - And *possible* to change equations of motion



The actual Gabriel

Setting up the physical system

- Set the *scale(s)* of the problem by defining dimensionless variables
- Make choices for the *physical* parameters of the system
- Make choices for the *numerical* parameters of the system
 - Chose a scheme to discretize the system
 - Be aware that you will need to vary these parameters in order to validate your numerical work

Defining dimensionless variables

...the hard part

- The two (structural) parts of the system are:

- Spacetime:

$$dx^\mu = \frac{dx_{\text{pr}}^\mu}{B}$$

- Gravity:

$$\left(\frac{\dot{a}}{a}\right)^2 = B^2 \left(\frac{a'}{a}\right)^2 = \frac{8\pi}{3m_{\text{pl}}^2} \rho$$

$$\rho_{\text{pr}} = \frac{\rho}{m_{\text{pl}}^2 B^2}$$

$$' \equiv \frac{\partial}{\partial t_{\text{pr}}}$$

Rescaling the fields

...is also not a choice

- The fields are always in Planck masses,

$$\phi_{\text{pr}} = \frac{\phi}{m_{\text{pl}}}$$

- So *derivatives* of fields have a factor of B as well as a factor of m_{pl} , e.g.

$$\dot{\phi} = B m_{\text{pl}} \frac{\partial \phi_{\text{pr}}}{\partial t_{\text{pr}}}$$

For scalar fields only,

Setting the scale(s) of the problem

- Rescaling the energy density is really just rescaling the potential (model)

$$\rho_{\text{pr}} = \frac{\frac{1}{2}\dot{\phi}^2 + \frac{(\nabla\phi)^2}{2a^2} + V(\phi)}{B^2 m_{\text{pl}}^2} = \frac{1}{2}\phi'_{\text{pr}}{}^2 + \frac{(\nabla_{\text{pr}}\phi_{\text{pr}})^2}{2a^2} + \frac{V}{B^2 m_{\text{pl}}^2}$$

- So the scale of the box is really set by the parameters of the potential, e.g.

$$V_{\text{pr}} = \frac{V}{B^2 m_{\text{pl}}^2} = \frac{1}{B^2 m_{\text{pl}}^2} \frac{1}{2} m^2 \phi^2 = \frac{1}{2} \phi_{\text{pr}}^2 \quad \text{where} \quad B = m$$

If the potential is more complicated

Setting the scale(s) of the problem

- For example, Axion Monodromy,

$$V_{\text{pr}} = \frac{m^2 M^2}{B^2 m_{\text{pr}}^2} \left(\sqrt{1 + \frac{\phi^2}{M^2}} - 1 \right)$$

- There appear to be many choices of B that could simplify this potential; however, the choice $B = m$ leads to

$$V_{\text{pr}} = \left(\frac{M}{m_{\text{pl}}} \right)^2 \left(\sqrt{1 + \frac{\phi_{\text{pr}}^2}{(M/m_{\text{pl}})^2}} - 1 \right) \approx \frac{1}{2} \phi_{\text{pr}}^2 + \mathcal{O}(\phi_{\text{pr}}^4)$$

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Where the (small amplitude) oscillation of the homogeneous field still sets the clock

$$V_{\text{pr}} = \left(\frac{M}{m_{\text{pl}}} \right)^2 \left(\sqrt{1 + \frac{\phi_{\text{pr}}^2}{(M/m_{\text{pl}})^2}} - 1 \right) \approx \frac{1}{2} \phi_{\text{pr}}^2 + \mathcal{O}(\phi_{\text{pr}}^4)$$

You only get one

Setting the scale(s) of the problem

- However, the choice of B "uses up" the freedom to set other scales of the problem.
 - In broad strokes, the number of parameters (beyond one) that you need to specify are the root of the numerical challenges
- Other codes (most notably LatticeEasy) have more freedom in choosing dimensionless variables

$$\phi_{\text{pr}} = Aa^r \phi \quad d\vec{x}_{\text{pr}} = B d\vec{x} \quad dt_{\text{pr}} = Ba^s dt \quad s = 2r - 3$$

- However, nothing is "free" — these choices change couplings to gravity, e.g., which just shift around where you're making your choices!

The physical parameters

Choosing the physics

- This could mean choosing masses, couplings, etc,

$$V = \frac{1}{2}m^2\phi^2 + \frac{g^2}{2}\phi^2\chi^2 \quad m = 10^{-6} m_{\text{pl}} \quad g^2 = 2.5 \times 10^{-7}$$

- Or the initial conditions (say, at the end of inflation)

$$\begin{aligned} \phi_0 &= 0.193 m_{\text{pl}} & \dot{\phi}_0 &= -0.142 m m_{\text{pl}} \\ \phi_0^{\text{pr}} &= 0.193 & \dot{\phi}_0^{\text{pr}} &= -0.142 \end{aligned}$$

- Which can also give other *physical* quantities of interest, e.g. H

The physical choices *inform* the numerical ...but they do not define them

- For vanilla preheating, we find

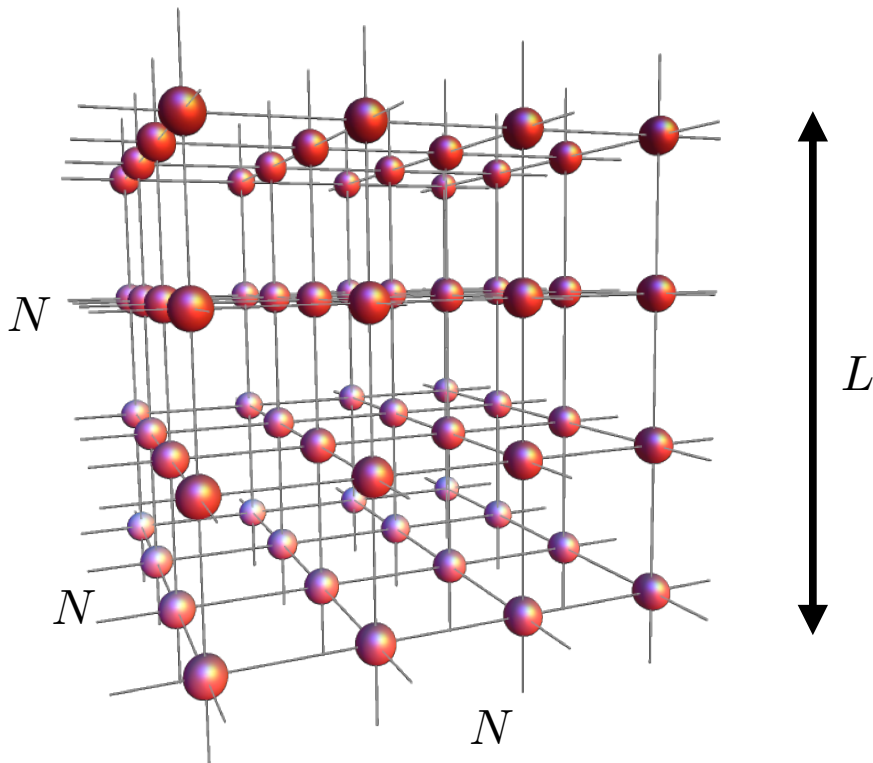
$$H_{\text{pr}} \approx 0.49 \sim 0.5$$

- Which means that we *should* think about

$$L_{\text{pr}} \sim 2$$

- BUT: this only suggests that the physics we're interested in are on *this scale*, and that choices of $L_{\text{pr}} = 10$ or $L_{\text{pr}} = 0.5$ should give similar phenomenology

Basic Numerical Parameters



$$k_{\min} = \frac{2\pi}{L} \longleftrightarrow \lambda_{\max} = L$$

$$k_{\max} = \frac{\sqrt{3}N}{2} \frac{2\pi}{L} \longleftrightarrow \lambda_{\min} = \frac{L}{N} \frac{2}{\sqrt{3}} \approx \Delta x$$

So what should L be?

...what it needs to be

- You have to *vary* L and N (among others) to ensure that the *physics* you are looking for is independent of these choices.
- This is known as a *convergence test* — that the *physics* converges as the simulation more closely approximates the continuum
- This step is **crucial** to convince us that the simulations are predicting outcomes from the continuum theory.
 - **Regardless of any other measure (e.g. energy conservation), you must show convergence if we are to believe that the discrete system approximates the continuum.**

And of course, Δt

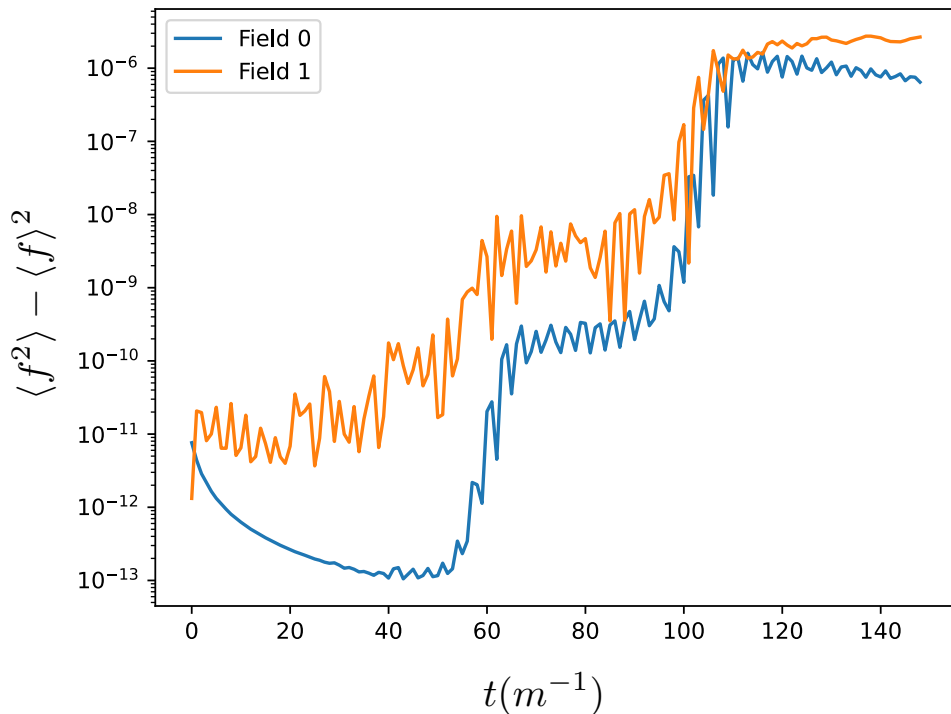
- The Courant–Friedrichs–Lewy (or just Courant) Condition give *guidance* as to how small your timestep should be
- Basically, you need enough *time* resolution to resolve the fastest-oscillating mode, e.g.

$$dt \lesssim T \lesssim \frac{2\pi}{\omega_{\max}} = 2\pi \frac{2}{\sqrt{3}} \frac{1}{2\pi} \frac{L}{N} \mathcal{O}(1) \times \Delta x$$

- But, we probably knew that already. This guidance is only a place to *start*; you still have to run a *timestep* convergence test to ensure that the timestep is small enough.

Even when you win, it might only be temporary

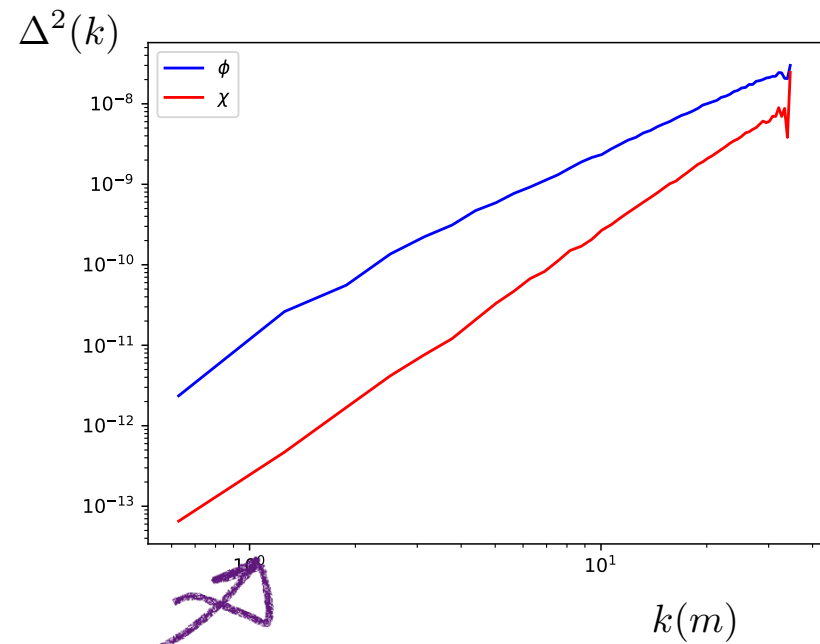
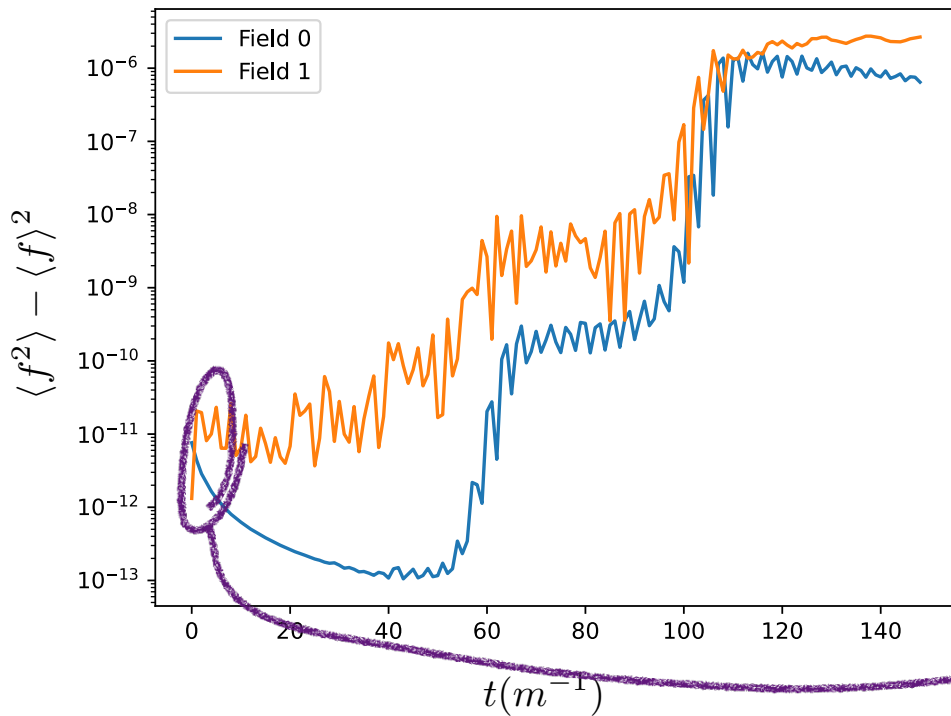
...consider vanilla preheating



- This is a plot of the variance, bumpiness, of the fields as a function of time for the model/parameters that come “shipped” with GABE
- This is the classic, ‘vanilla preheating’ model where we have the three stages of preheating
- So we can look at the modes (particle production) over the course of this run

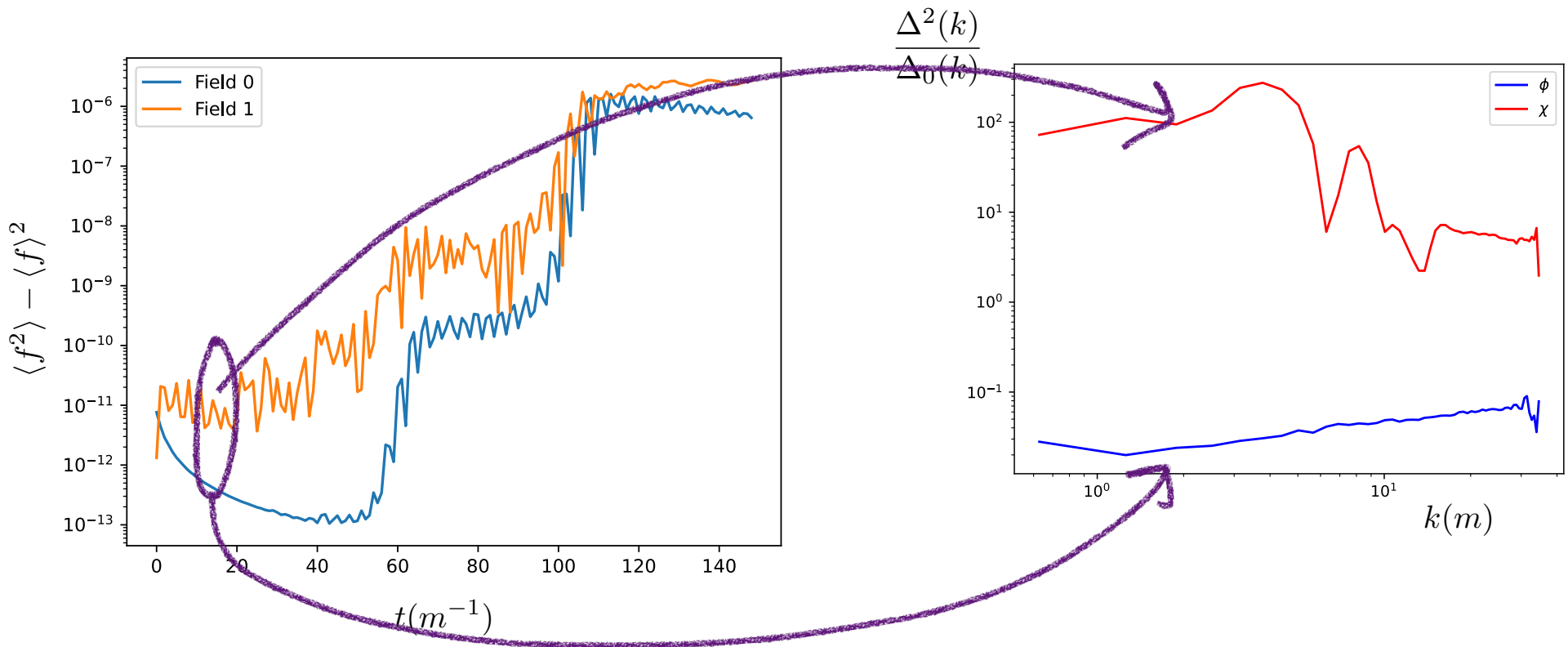
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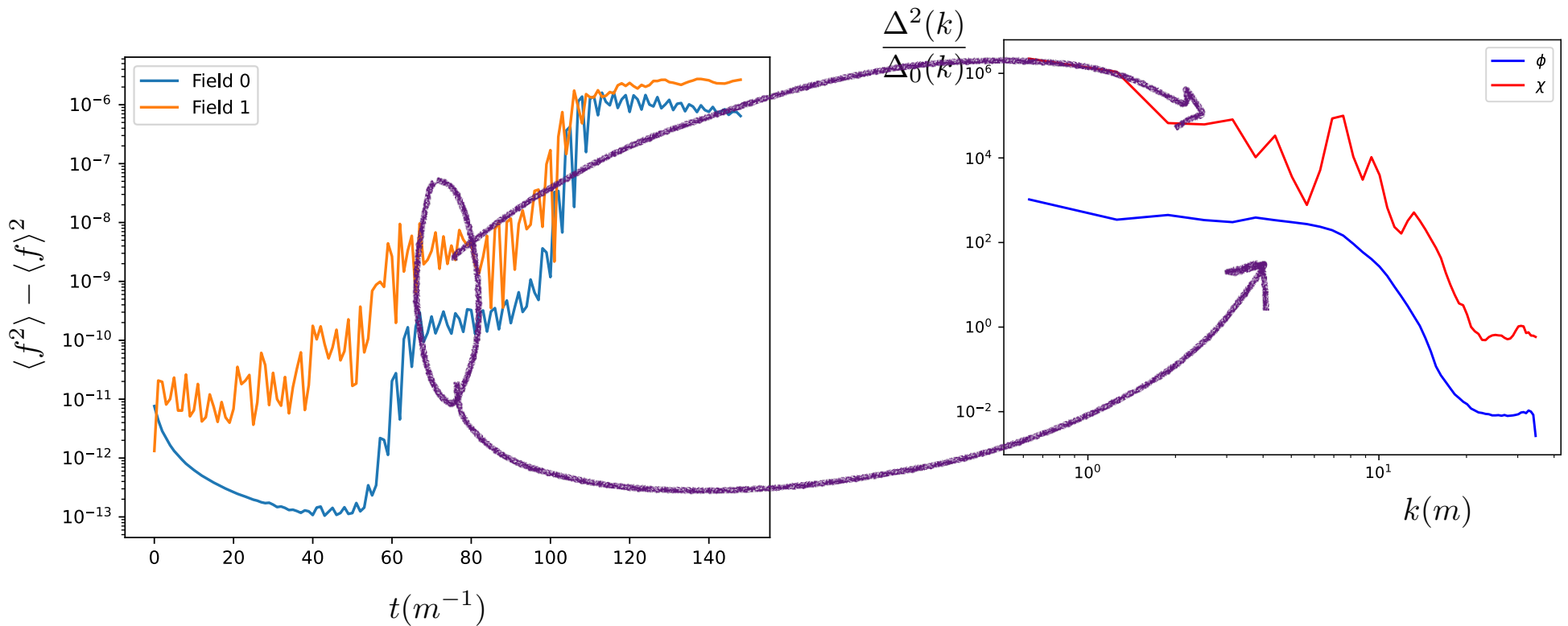
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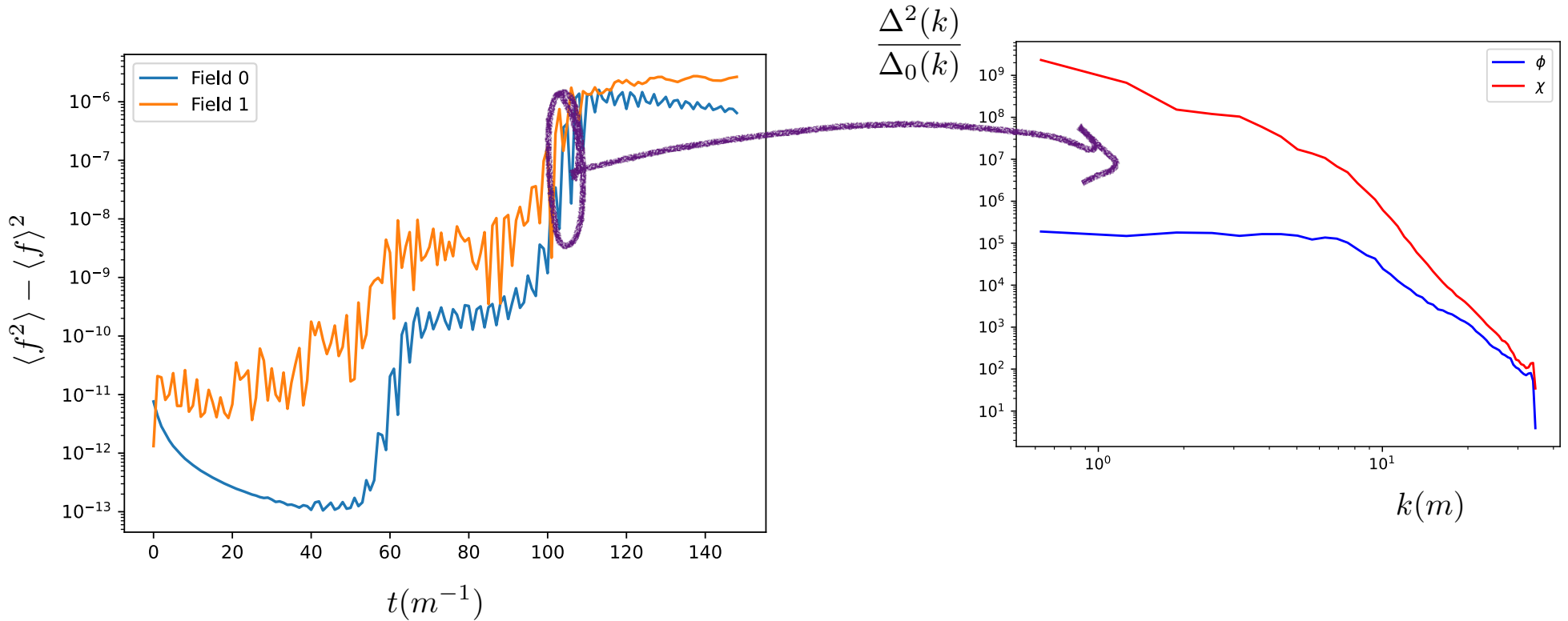
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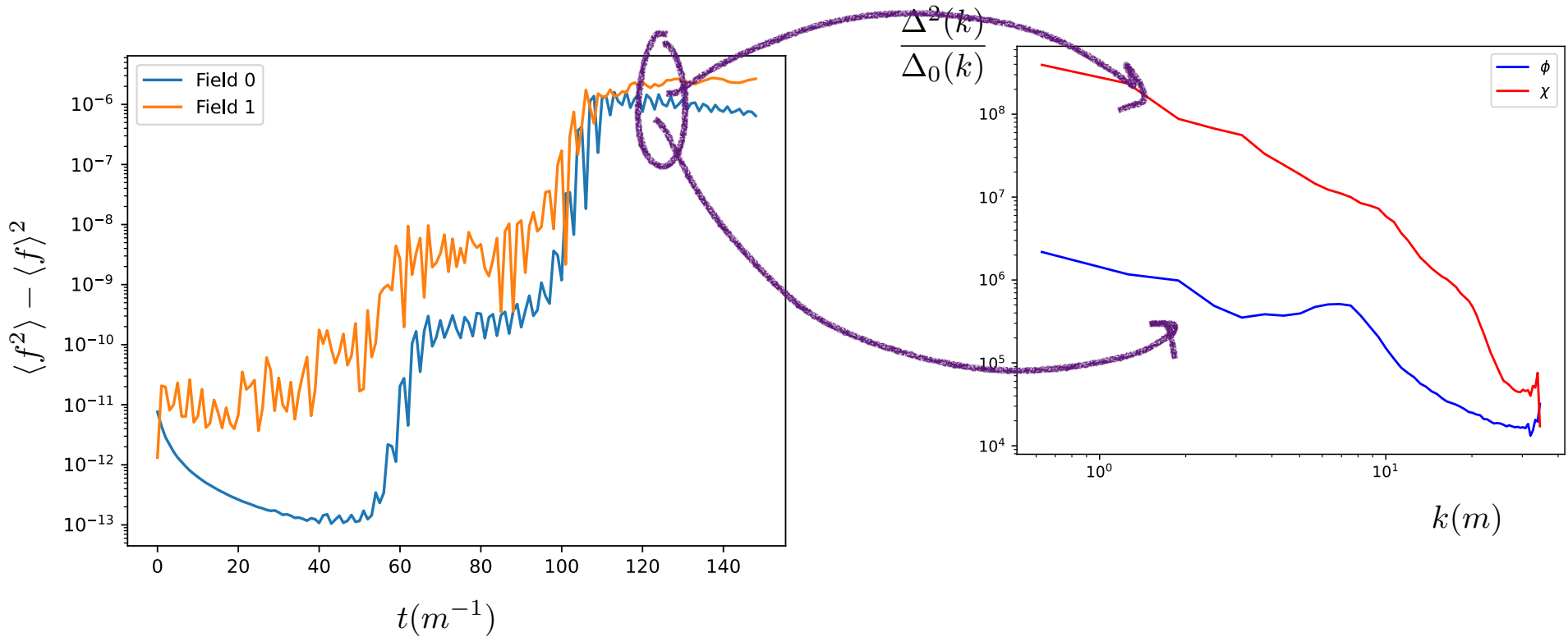
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Once power gets into those high modes

- Remember that gravitational waves

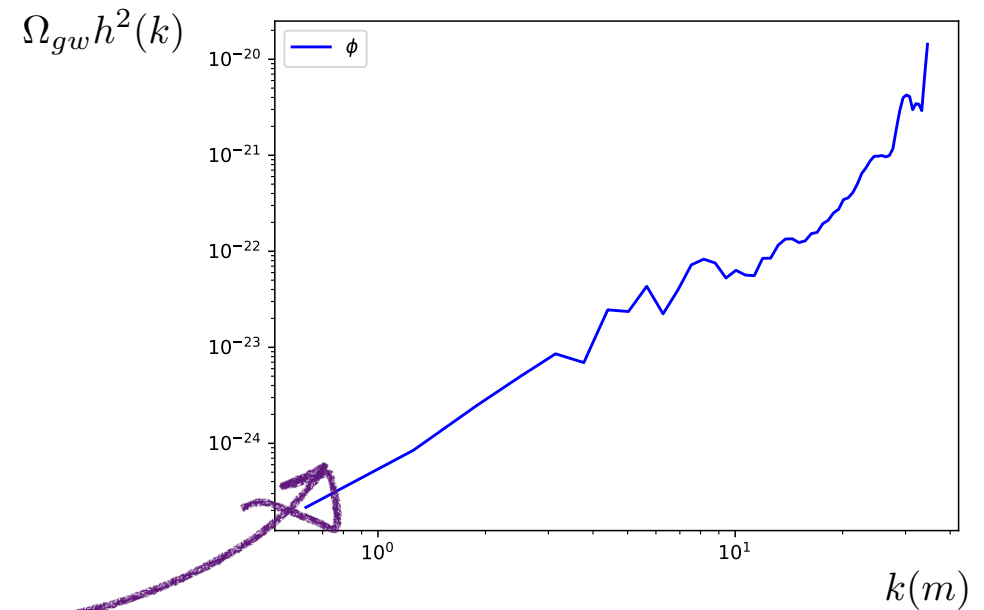
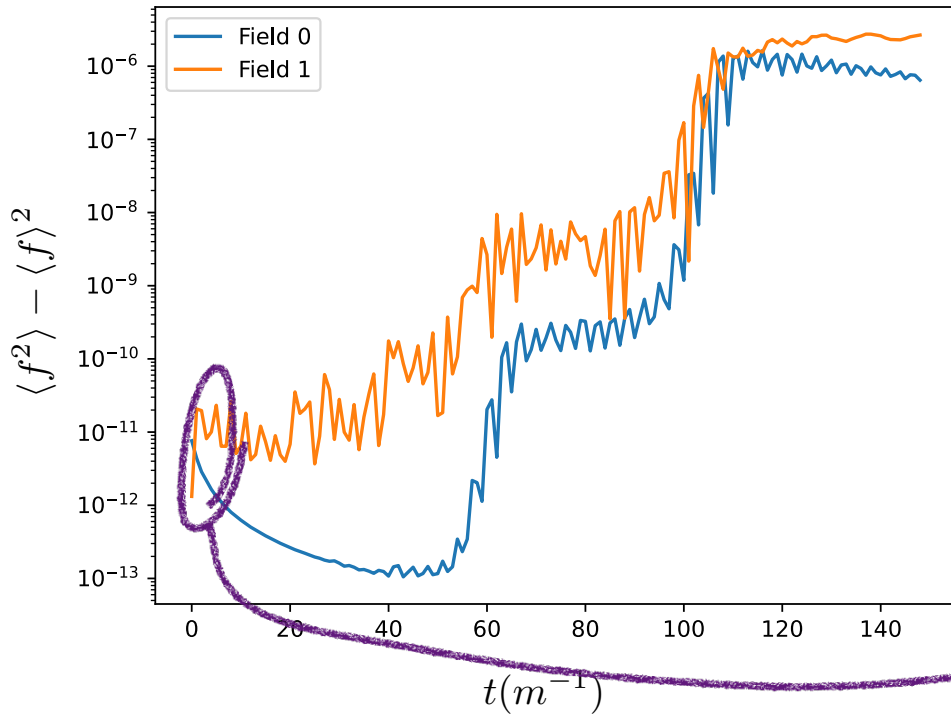
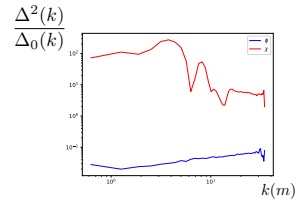
$$\ddot{h}_{ij} + 3H\dot{h}_{ij} - \nabla^2 h_{ij} = \frac{18\pi G}{a^2} S_{ij}^{\text{TT}}$$

- Are sourced by non-linear combinations of derivatives,

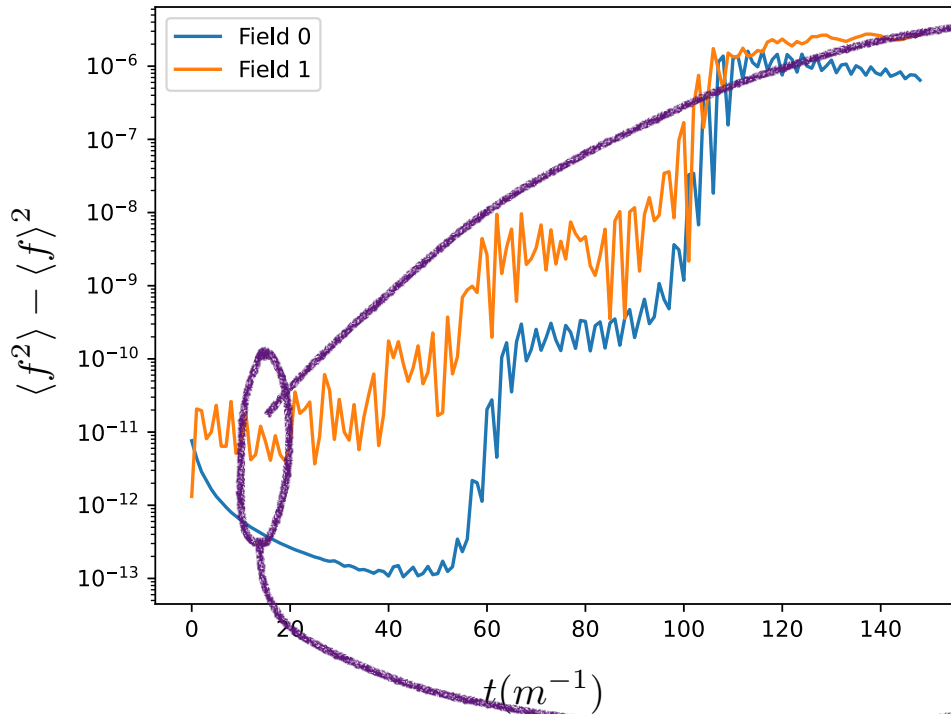
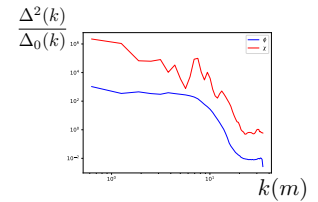
$$S_{ij} \sim \partial_i \phi \partial_j \phi$$

- They are very sensitive to any errors in high-frequency modes!

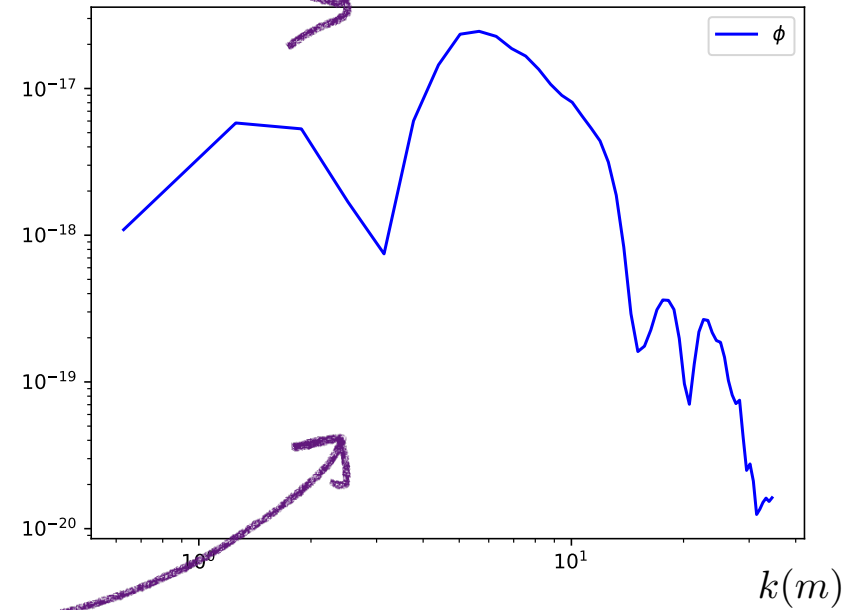
Even when you win, it might only be temporary ...consider vanilla preheating...with gravitational waves



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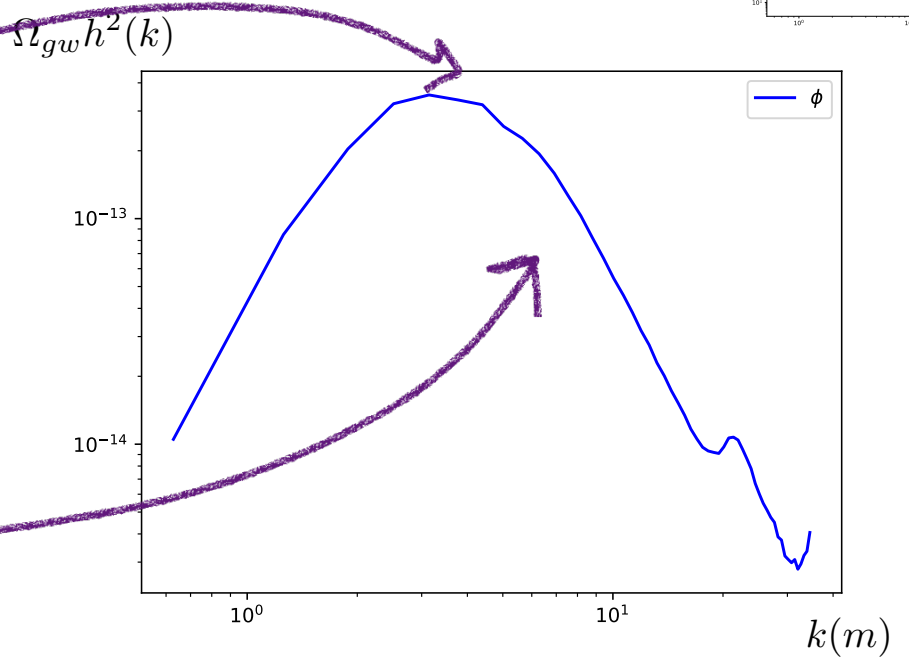
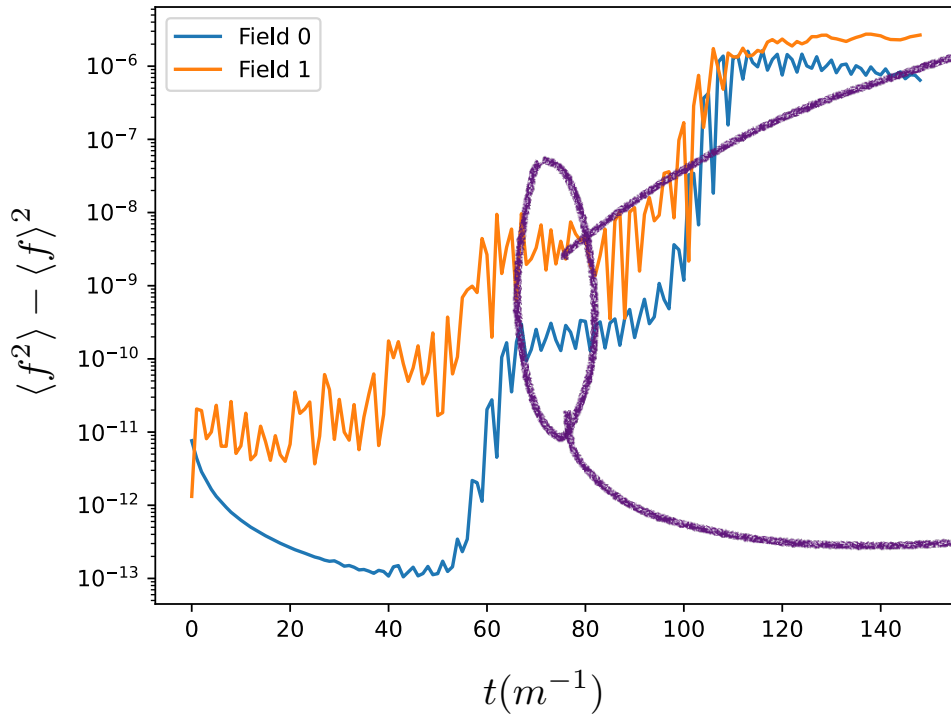
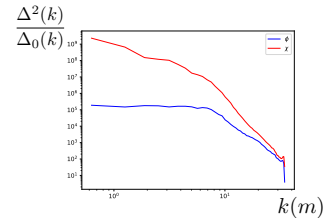
$\Omega_{gw} h^2(k)$



$k(m)$

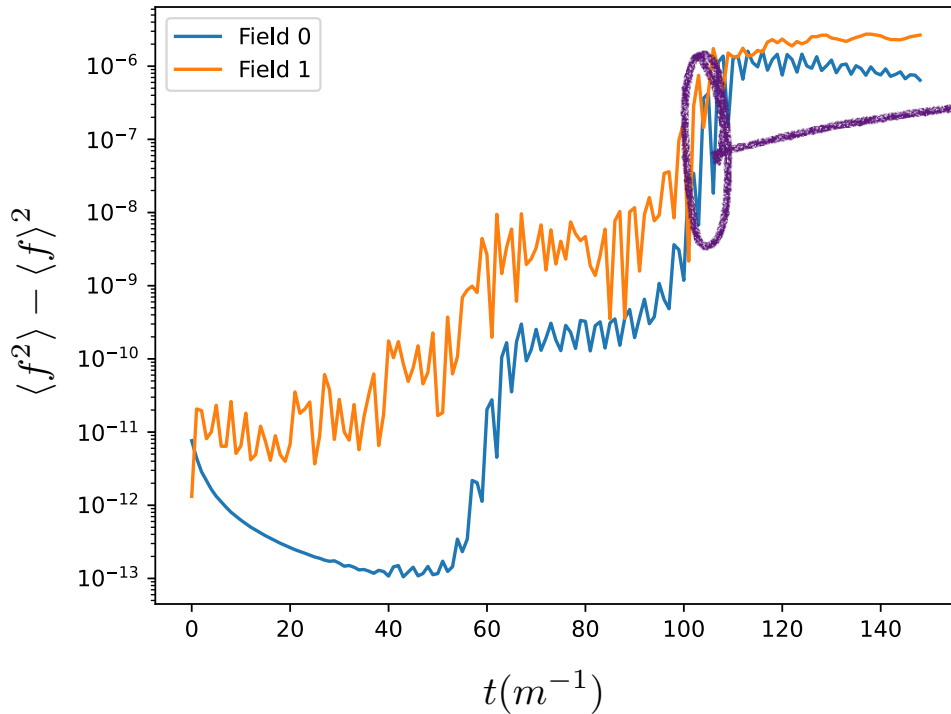
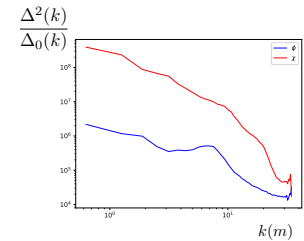
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...consider vanilla preheating...with gravitational waves

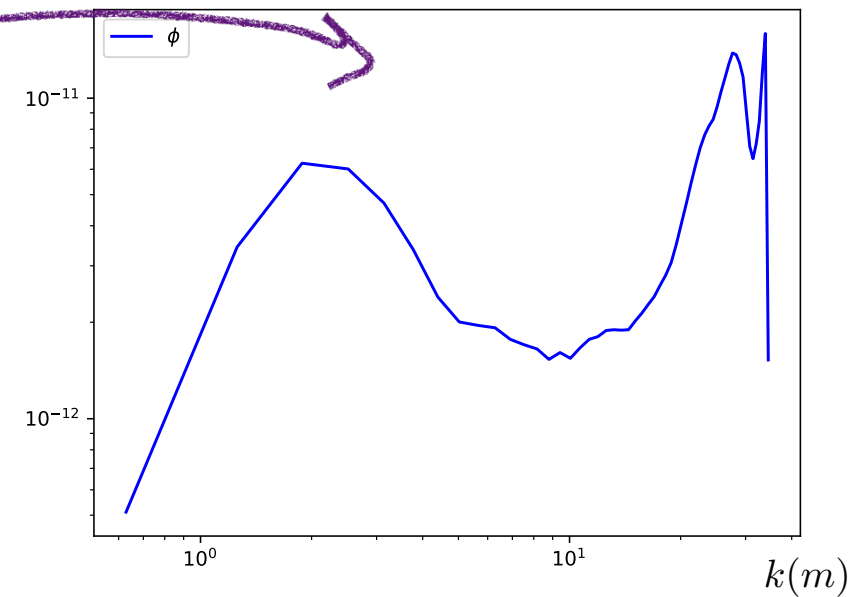


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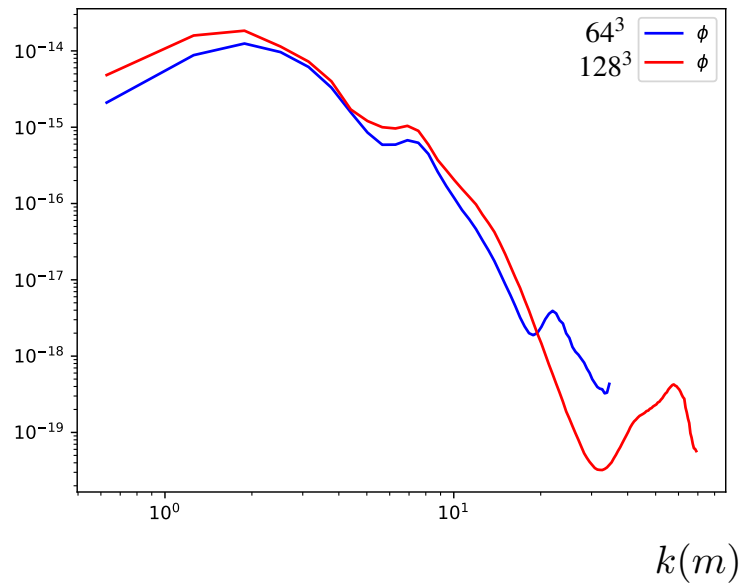
$$\Omega_{gw} h^2(k)$$



How do we know what's real??

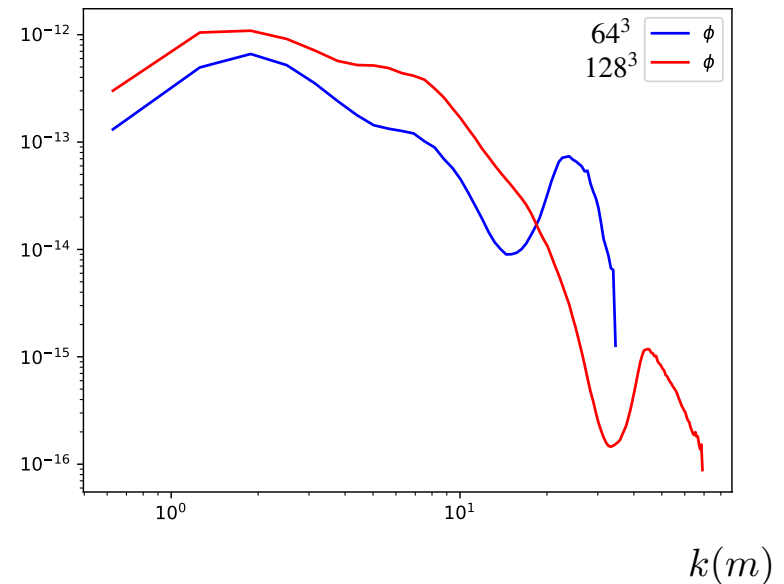
These spectra *change* when you change the numerical parameters!

$$\Omega_{gw} h^2(k)$$



$$t = 112 m^{-1}$$

$$\Omega_{gw} h^2(k)$$



$$t = 117 m^{-1}$$

Some things I find confusing

Initial conditions

The vacuum

- Assuming that modes are *sub-horizon* (or nearly sub-horizon, we assume that the fields have Bunch-Davies initial conditions,

$$\mathcal{P}_k = \frac{1}{2\omega_k} \quad \omega_k = \sqrt{k^2 + m_{\text{eff}}^2}$$

- On the lattice, we need to translate this to the *discrete system*. For the most part this has a straightforward definition,

$$\phi(\vec{k}) = \frac{1}{(2\pi)^{3/2}} \int d^3x \phi(\vec{x}) e^{-i\vec{k}\cdot\vec{x}} \quad \Phi(\vec{k}_j) = \frac{1}{(2\pi)^{3/2}} (\Delta x)^3 \sum_i \phi(\vec{x}_i) e^{-i\vec{k}_j\cdot\vec{x}_i}$$

We chose the 2-pt correlation function

to be the object that is invariant between the continuum and the discrete

$$\begin{aligned}\langle \phi(\vec{x})\phi(\vec{y}) \rangle &= \frac{1}{(2\pi)^3} \int d^3k d^3p \langle \phi(\vec{k})\phi(\vec{p}) \rangle e^{i\vec{k}\cdot\vec{x}} e^{-i\vec{p}\cdot\vec{y}} = \frac{1}{(2\pi)^3} \int d^3k \mathcal{P}_k e^{i\vec{k}(\vec{x}-\vec{y})} \\ &= \frac{1}{(2\pi)^3} (\Delta k)^6 \sum_i \sum_j \langle \Phi(\vec{k}_l)\Phi(\vec{p}_m) \rangle e^{i\vec{k}_l\cdot\vec{x}_i} e^{-i\vec{p}_m\cdot\vec{y}_j} = \frac{1}{(2\pi)^3} (\Delta k)^3 \sum_l \mathcal{P}_k e^{i\vec{k}_l(\vec{x}_i-\vec{y}_j)}\end{aligned}$$

So there's a modification of the momentum-space 2-pt correlation function:

$$\begin{aligned}\langle \phi(\vec{k})\phi(\vec{p}) \rangle &= \mathcal{P}_k \delta(\vec{k} - \vec{p}) \\ \langle \Phi(\vec{k}_l)\Phi(\vec{p}_m) \rangle &= (\Delta k)^3 \mathcal{P}_k \delta_{lm} = \left(\frac{L}{2\pi}\right)^3 \mathcal{P}_k \delta_{lm}\end{aligned}$$

This has an impact on the initial conditions

- When we convert the power spectrum to *dimensionless* units,

$$\langle |\phi_{\text{pr}}(k_j)| \rangle = \left(\frac{B^6}{m_{\text{pl}}^2} \right) \left(\frac{L_{\text{pr}}}{B^2 2\pi} \right)^3 \mathcal{P}_k \quad \mathcal{P}_k^{\text{pr}} = \frac{B^3}{m_{\text{pl}}^2} \mathcal{P}_k$$

- Which means for Bunch-Davies....

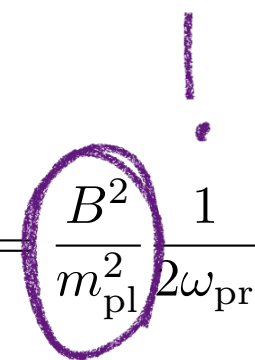
$$\mathcal{P}_{k,\text{BD}}^{\text{pr}} = \frac{B^3}{m_{\text{pl}}^2} \mathcal{P}_{k,\text{BD}} = \frac{B^2}{m_{\text{pl}}^2} \frac{1}{2\omega_{\text{pr}}}$$

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Thoughts on strong hyperbolicity

...make it wavy

- The tricks we play (as theorists) to reduce the number of degrees of freedom can negatively affect numerical stability
- Adding degrees of freedom keep equations strongly hyperbolic (that is, wave-like) which means you need to store (and evolve) more information than you *have to*.
- But it makes the problem solvable. Examples include:
 - Using Lorenz gauge, and keep track of constraints
 - Yesterday, I talked about BSSN and how this works for Numerical Relativity
 - Stiff equations of motion can be stabilized with extra degrees of freedom

Looking at Scalar Galileons

We can

- We start with a stiff, derivatively coupled equation of motion,

$$\square\pi + \frac{1}{3\Lambda^3} ((\square\pi)^2 - (\partial_\mu\partial_\nu\pi)^2) = -\frac{T}{3m_{\text{pl}}}$$

- By identifying

$$H_{\mu\nu} = \partial_\mu\partial_\nu\pi \qquad A_\mu = \partial_\mu\pi$$

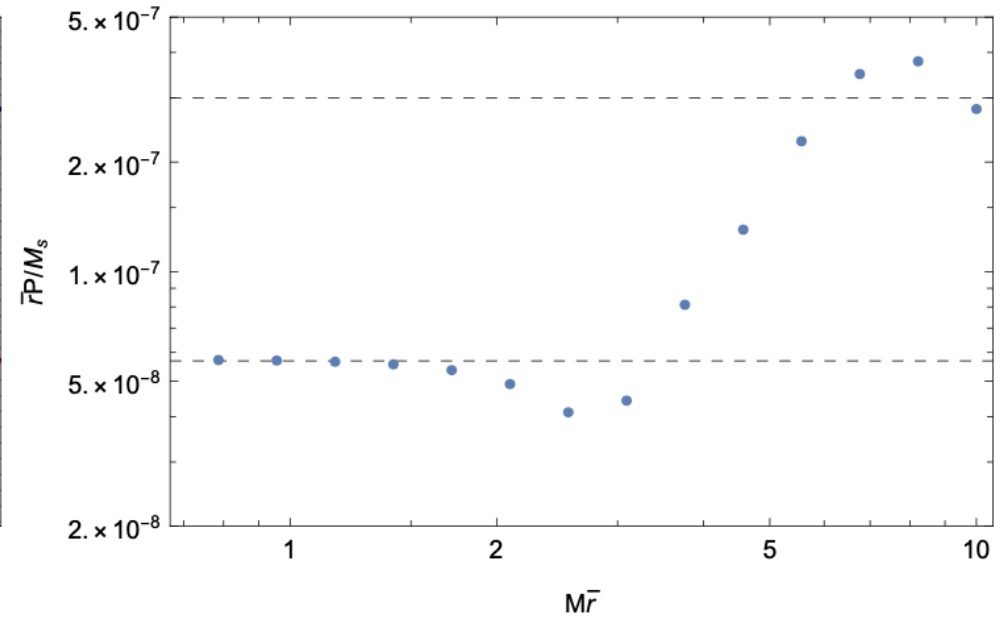
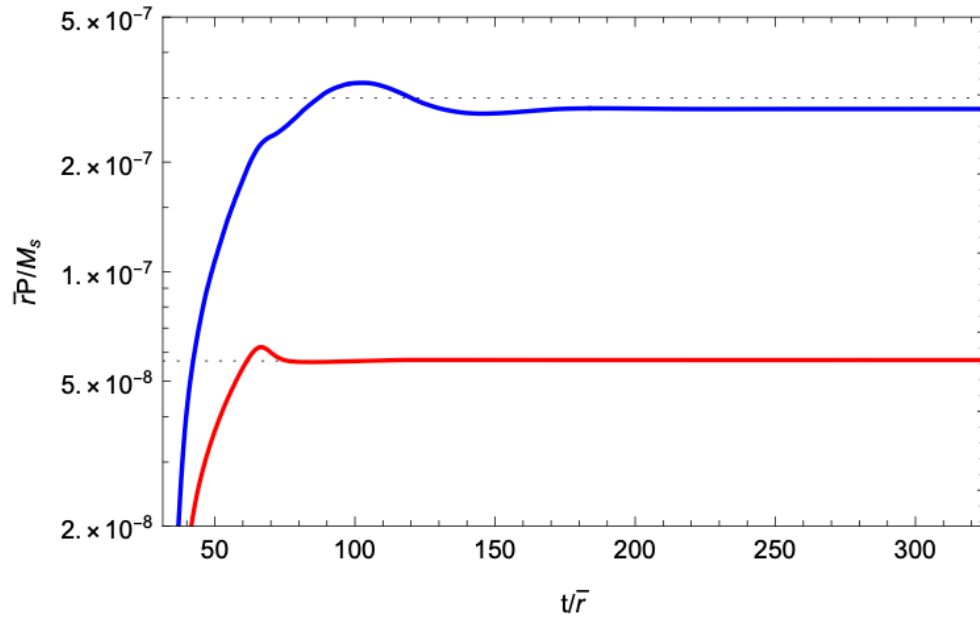
- We get a more complicated system, but one that's strongly hyperbolic

$$\square\pi + \frac{1}{3\Lambda^3} (H^{\mu\nu}H_{\mu\nu} - (H^\nu{}_\nu)^2) = -\frac{T}{3m_{\text{pl}}}$$

$$\square A_\mu - \frac{1}{\tau}\partial_t A_\mu - M^2 A_\mu = -M^2\partial_\mu\pi$$

$$\square H_{\mu\nu} - \frac{1}{\tau}\partial_t H_{\mu\nu} - M^2 H_{\mu\nu} = -\frac{M^2}{2} (\partial_\mu A_\nu + \partial_\nu A_\mu)$$

We get the behavior of the full system ...but much more reliably



A quick example

...in the case of oscillons

- Many potentials create oscillons; we've been interested in α -attractor models which, in the absence of a coupled field, are great oscillon producers
- An open question has been, *do oscillons decay* which can be studied by considering an explicit coupling to another field

$$\mathcal{L}_{\text{int}} = \frac{g^2}{2} \phi^2 \chi^2$$

scalar



Peter Krosniak
'27



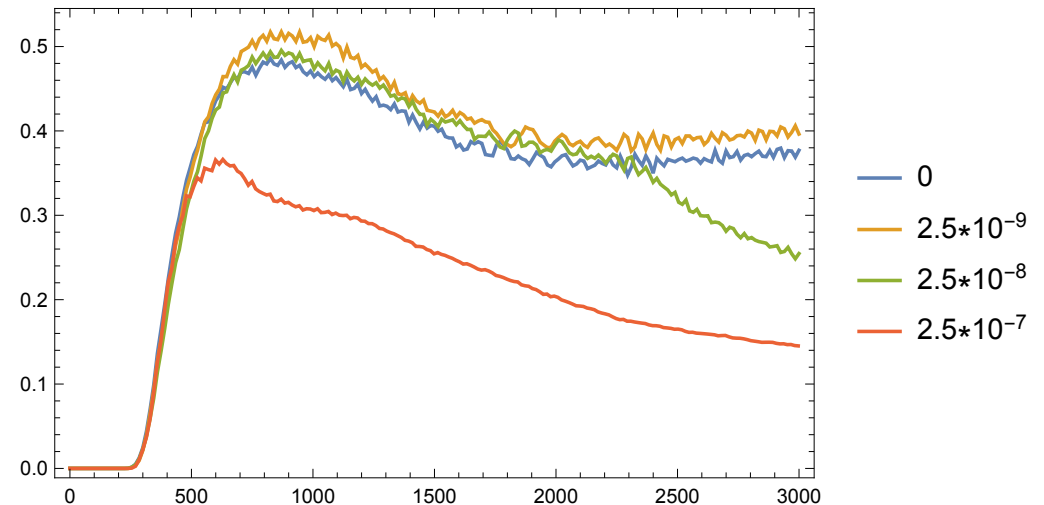
J'sun Gardner
'26

The End of the Oscillon

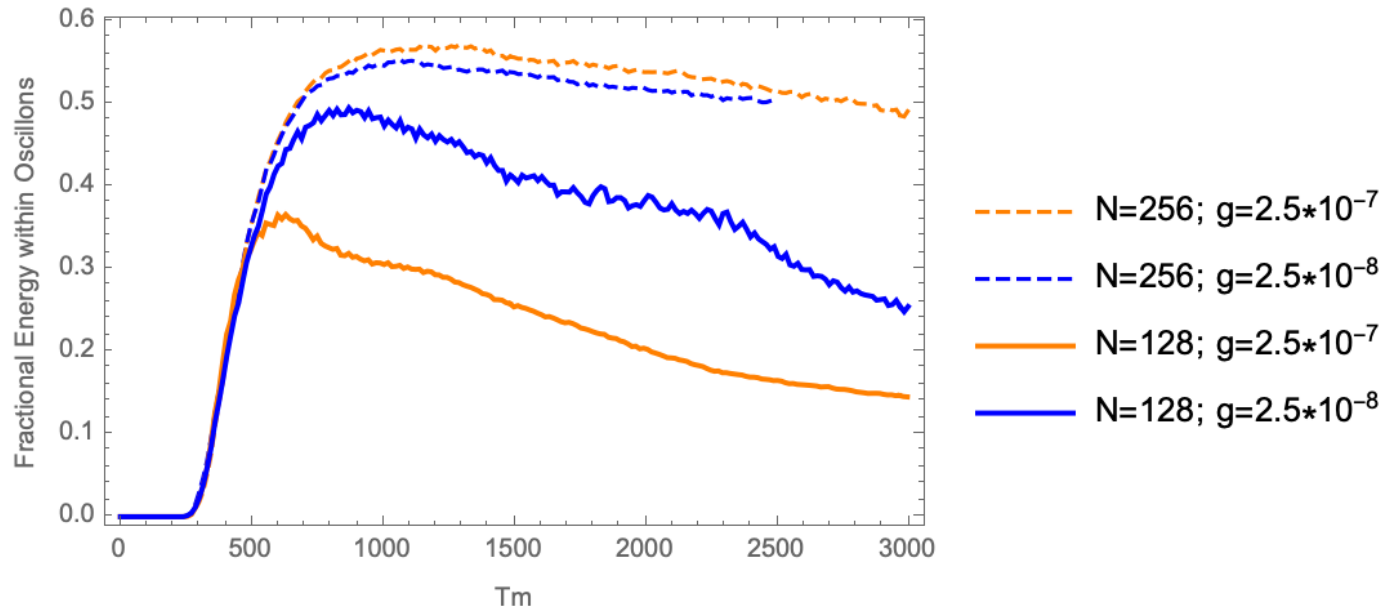
- Consider the E-model α -attractor

$$V = \frac{m^2 \mu^2}{2} \left(1 - e^{-\frac{\phi}{\mu}}\right)^2$$

- We can see that the energy in oscillons decays parametrically with the strength of the interaction
- But is it *real*?



When we look at ...the convergence test

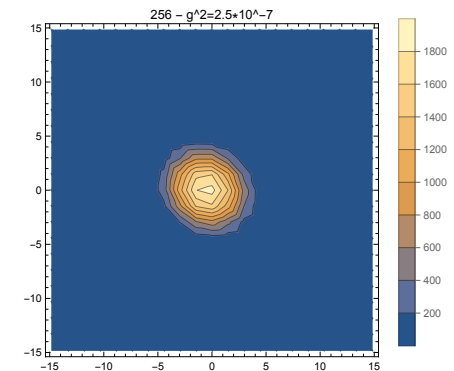
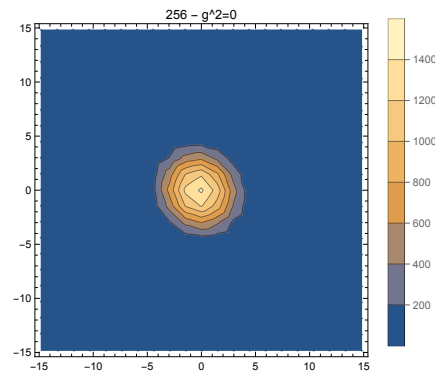
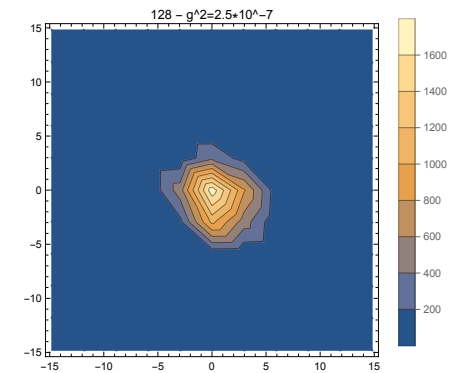
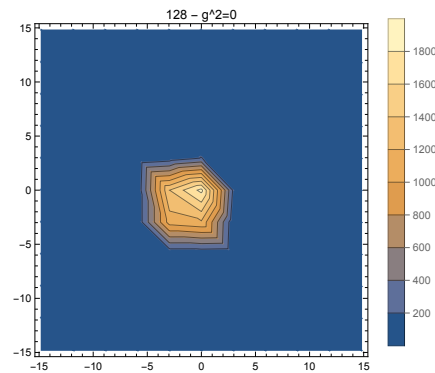


Why?

...we can actually see what's going on!

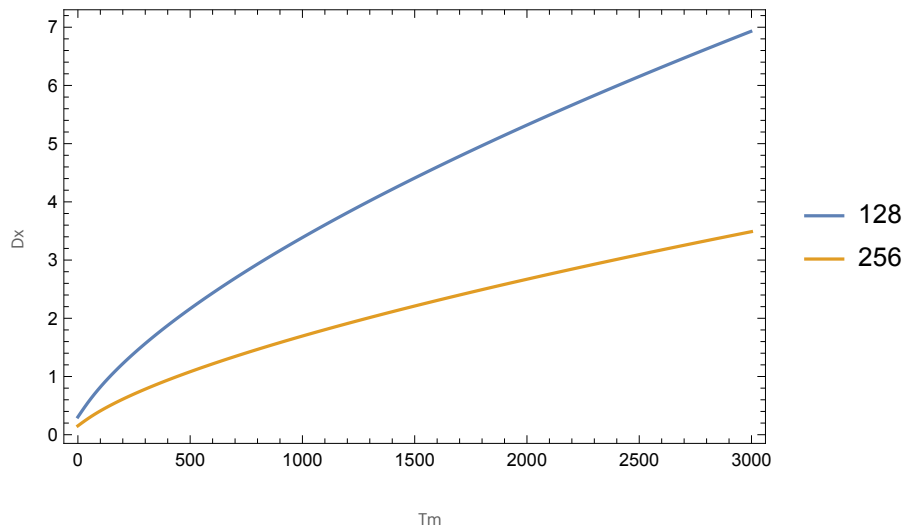
The size and shape of oscillons in 128^3 and 256^3 resolutions are different

128^3 isn't enough points to form oscillons with a circular shape

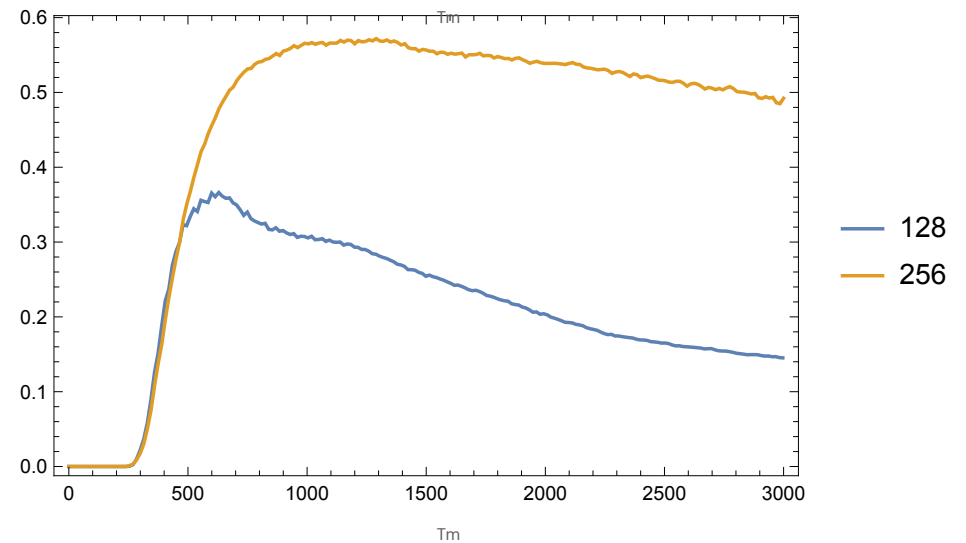


When do we approach the modes ...that we can't trust?

Physical dx of E-Model 128 and 256 $2.5 \cdot 10^{-7}$ runs vs Time



E-Model 128 and 256 $2.5 \cdot 10^{-7}$ runs vs Time



For this afternoon:

... if you're participating

- Please download GABE
 - cosmo.kenyon.edu/gabe.html
- If you have a MAC
 - Try to run the instructions on that page, new(ish) operating systems should work without issue
- If you have LINUX
 - Please install fftw3 (enabling regular and long-double: openmp and threads for each)
- If you plan to use the remote version, please email me (giblinj@kenyon.edu) as soon as possible!
- **Please also have Jupyter (or similar) for plotting. I will distribute a .ipynb this afternoon!**

Advice from my students

- Write the code yourself, so you can change anything. Seems silly but it was helpful to me to realize that I can communicate via the code (`printf("I'm here :)\n");`)
- When in doubt, plot it out - there's lots of data and it can tell you many things
- Write in *words* what you want the code to do. Translate it to a step by step list and check that this matches the order of the code
- As always - fix the bug you know :)
- The physics doesn't care about the math and we want to find the region where it doesn't care about the code either, so change code parameters to find the physics
- This is a basic idea but the code is only as smart as you let it be, so let it be dumb and change single things at a time so you know for sure what is going on.
- Go to a known regime: homogeneous, static, symmetric. Check it behaves as expected