

On stability of numerical schemes

How to discretize time for some problems

18.303 Linear Partial Differential Equations: Analysis and Numerics

Behavior of time-dependent problems

Let us consider the differential equation

$$\alpha \frac{\partial^2 u(t, \mathbf{x})}{\partial t^2} + \mu \frac{\partial u(t, \mathbf{x})}{\partial t} = \mathcal{L}u(t, \mathbf{x})$$

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with some boundary conditions that we will not specify at this point. Here we assume $\alpha, \mu \geq 0$.

We assume that \mathcal{L} has a complete eigenbasis $\{\phi_n\}$ with

$$\mathcal{L}\phi_n(\mathbf{x}) = \lambda_n \phi_n(\mathbf{x})$$

and

$$u(t, \mathbf{x}) = \sum_n \hat{u}_n(t) \phi_n(\mathbf{x}).$$

Plugging the expansion of u in the differential equation and using the *linear independence* of ϕ_n gives

$$\alpha \frac{\partial^2 \hat{u}_n(t)}{\partial t^2} + \mu \frac{\partial \hat{u}_n(t)}{\partial t} = \lambda_n \hat{u}_n(t).$$

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Plugging this in gives the characteristic equation

$$(\alpha a_n^2 + \mu a_n - \lambda_n) \exp(a_n t) = 0,$$

which can be solved giving

$$a_n = \frac{-\mu \pm \sqrt{\mu^2 + 4\alpha\lambda_n}}{2\alpha}$$

if $\alpha \neq 0$ and

$$a_n = \frac{\lambda_n}{\mu}$$

otherwise.

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Note: remember that the eigenvalues of the Laplacian are negative (semi)definite. The modes with the zero eigenvalue are constant in time and the rest of them decrease. This is the case for the heat equation. See how important it is that the eigenvalues are negative?

Case ii: $\mu = 0$

This gives a kind of wave equation. Now the solution is

$$\exp\left(\pm\sqrt{\frac{\lambda_n}{\alpha}}t\right).$$

Again, if *all* the eigenvalues are negative, the solution will be oscillating. Otherwise there will be a solution (for each \mathbf{n}) that is decaying and a solution that grows exponentially. We also have the special case $\lambda_n = 0$ for modes that do not change in time.

$$\hat{u}_n(t) \propto \exp\left(\frac{-\mu \pm \sqrt{\mu^2 + 4\alpha\lambda_n}}{2\alpha}t\right)$$

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If we assume that $\alpha \neq 0$ we can write this as

$$\hat{u}_n(t) \propto \exp\left(\frac{-\mu \pm \sqrt{\mu^2 + 4\alpha\lambda_n}}{2}t\right) = \exp\left(\frac{-\mu t}{2}\right) \exp\left(\pm \frac{\sqrt{\mu^2 + 4\alpha\lambda_n}}{2}t\right)$$

with the rescaling $\mu \rightarrow \mu/\alpha$ and $\lambda_n \rightarrow \lambda_n/\alpha$ (if we were careful here, we would use different symbols for these new coefficients but let's allow us to be a bit lazy).

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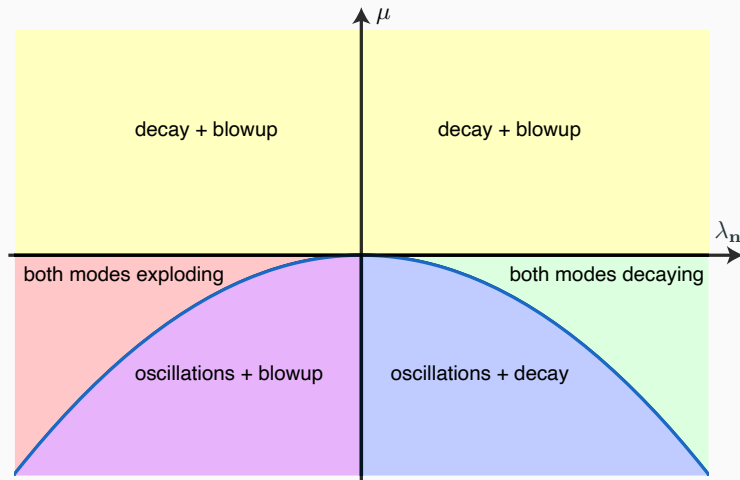


Diagram of different behavior

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- We will get back to energies later; for now we'll concentrate on stability in the sense that the solutions don't blow up

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Forward difference will give

$$\frac{u^{(k+1)} - u^{(k)}}{s} = \mathcal{L}u^{(k)}.$$

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Assume we have some *truncated* basis with N eigenvectors. We assume λ_N to be the largest eigenvalue (by absolute value). Now we get the condition

$$1 - s\lambda_N > -1 \Leftrightarrow s < 2/\lambda_N.$$

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Exercise: Assume $\mathcal{L} = \Delta$ (the Laplacian). Assume we have Dirichlet boundaries for the heat equation. Now

$$A = \frac{1}{\Delta x^2} \begin{pmatrix} -2 & 1 & & & \\ 1 & -2 & 1 & & \\ & \ddots & \ddots & \ddots & \\ & & 1 & -2 & 1 \\ & & & 1 & -2 \end{pmatrix}.$$

What is the stability condition for the time step?

The eigenvectors of A are $\phi_n(k) = \sin(nk\pi/(N+1))$ and the associated eigenvalues are

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This is called the Courant–Friedrichs–Lewy (CFL) condition. For those of you who did Pset 1, you probably noticed that the time step had to be very small. This explains why.

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This is called the Crank-Nicolson method. It's equivalent to using central difference for time at time $k + 1/2$. For the coefficients we get

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Word of warning: stability just means here that the discretized system doesn't blow up. We should still make s sufficiently small if we want to approximate the time evolution well.