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- In particular, by means of the perturbation method the equations have been **linearised**, making them amenable to analytical investigation.
- However, to obtain solutions in the general case, it is necessary to solve the **full nonlinear system**.
- In numerical weather prediction (NWP) the fully nonlinear primitive equations are solved by **numerical means**.
- In the atmosphere, the nonlinear **advection** process is a dominant factor.
- To get some idea of the methods used, we look at the simple problem of formulating time-integration algorithms for the solution of the **simple advection equation**.

Recap. of Discretization Methods

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- An *analytical* problem becomes an *algebraic* one.
- A problem with an *infinite* degree of freedom is replaced by one with a *finite* degree of freedom.
- A *continuous* problem goes over to a *discrete* one.

The Finite Difference Method

We start by looking at the *Taylor expansion* of $f(x)$:

$$f(x + \Delta x) = f(x) + f'(x).\Delta x + \frac{1}{2}f''(x)\Delta x^2 + [O(\Delta x^3)] \quad (1)$$

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We neglect these and obtain approximations for the derivative of $f(x)$ as follows:

$$f'(x) = \frac{f(x + \Delta x) - f(x)}{\Delta x} + O(\Delta x) = f'_F + O(\Delta x)$$

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Keeping only leading terms, we incur **errors of order** $O(\Delta x)$.

We can do better than this: subtracting (2) from (1) yields:

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Adding (1) and (2) gives the corresponding expression for the second derivative:

$$f''(x) = \frac{f(x + \Delta x) - 2f(x) + f(x - \Delta x)}{\Delta x^2} + O(\Delta x^2)$$

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Fourth-order accurate schemes are sometimes used in NWP, but *second order accuracy is more popular*.

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Consider the function $f(x) = A \sin(kx)$.

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- Compare these to the true derivative $f'(x)$ and investigate their behaviour for small Δx .
- Demonstrate thus that the **centered difference is of higher order accuracy**.

Grid Resolution and Accuracy

The size of the **gridstep** Δx determines the **accuracy** of the numerical scheme.

For the simple sine function the error depended on $k\Delta x = 2\pi\Delta x/L$, that is, on the ratio of the grid size Δx to the wavelength L .

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The higher the resolution, that is, the smaller the grid-size, **the heavier the computational burden.**

There is a *trade-off between resolution and accuracy.*

The Leapfrog Method

We consider the equation describing the conservation of a quantity $Y(x, t)$ following the 1D motion of a fluid flow:

$$\frac{dY}{dt} \equiv \left(\frac{\partial Y}{\partial t} + u \frac{\partial Y}{\partial x} \right) = 0.$$

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It is analogous to a factor of the wave equation:

$$\left(\frac{\partial^2}{\partial t^2} - c^2 \frac{\partial^2}{\partial x^2} \right) Y = \left(\frac{\partial}{\partial t} + c \frac{\partial}{\partial x} \right) \left(\frac{\partial}{\partial t} - c \frac{\partial}{\partial x} \right) Y = 0,$$

and its general solution is $Y = Y(x - ct)$.

Since the **advection equation is linear**, we can construct a general solution from Fourier components

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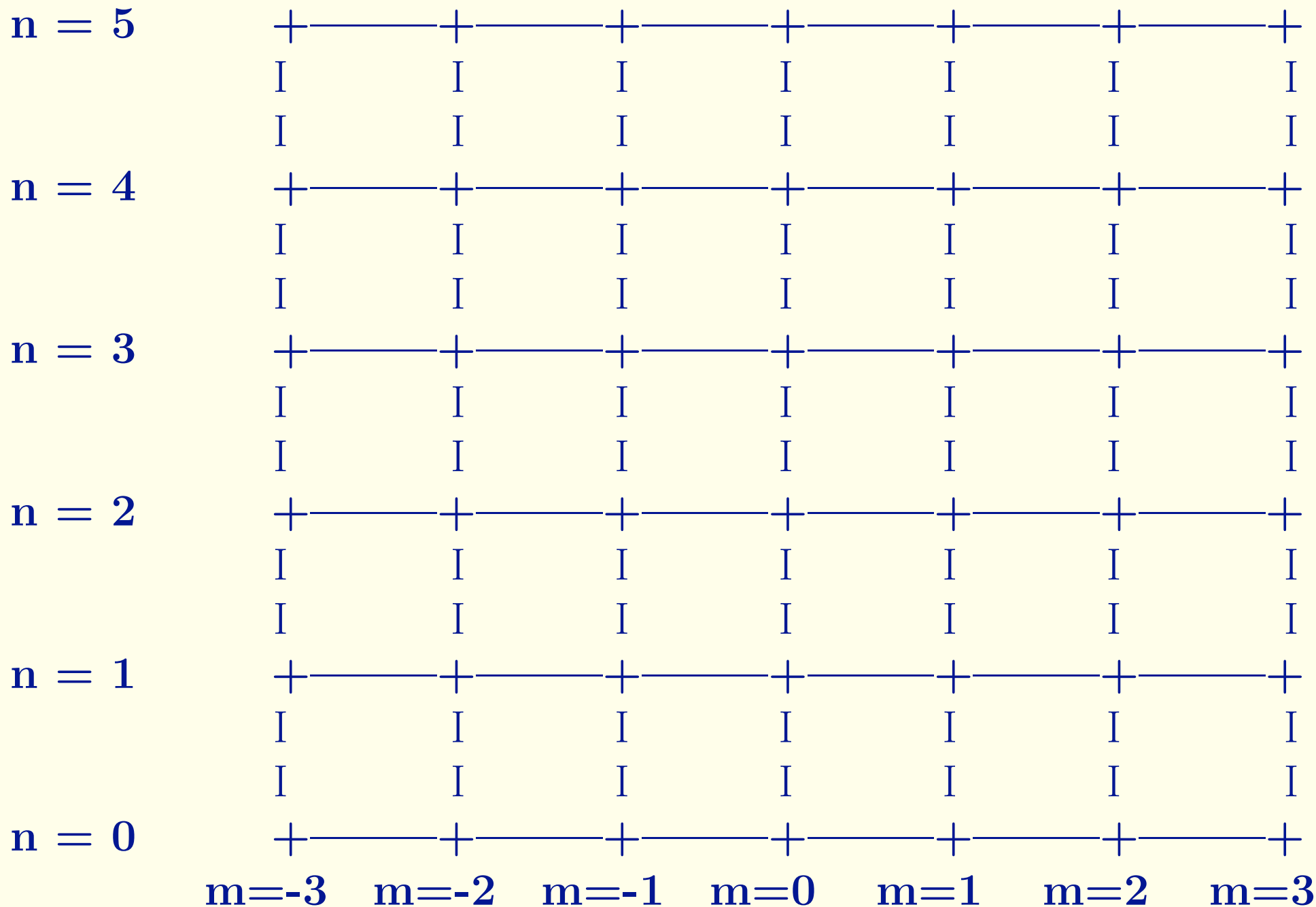
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Let the variables x and t be represented by the horizontal and vertical axes. Positive time corresponds to the upper half plane. The initial data occur on the x -axis.

Space-Time Grid:

Space axis horizontal
Time axis vertical



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Then the **(CTCS) finite difference approximation** to the differential equation may be written as follows:

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Solving for the value at time $(n+1)\Delta t$ gives

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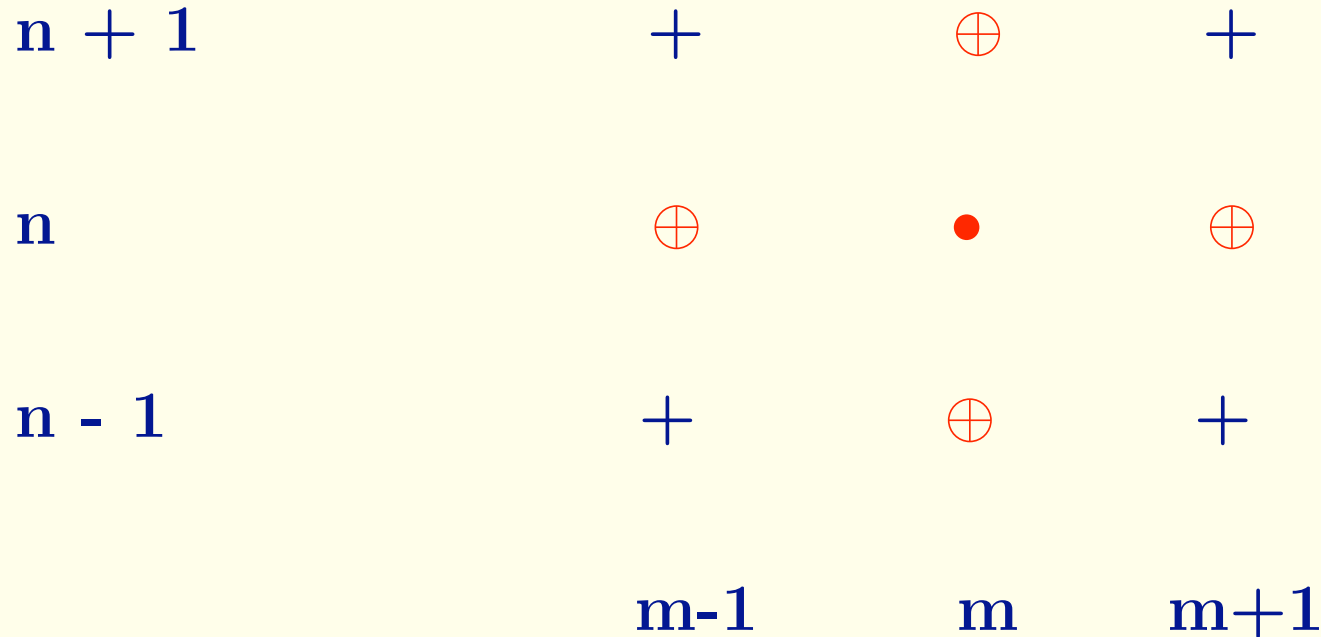
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The ratio $\mu \equiv \frac{c\Delta t}{\Delta x}$ will be found to be crucial.

Inter-dependency of Points

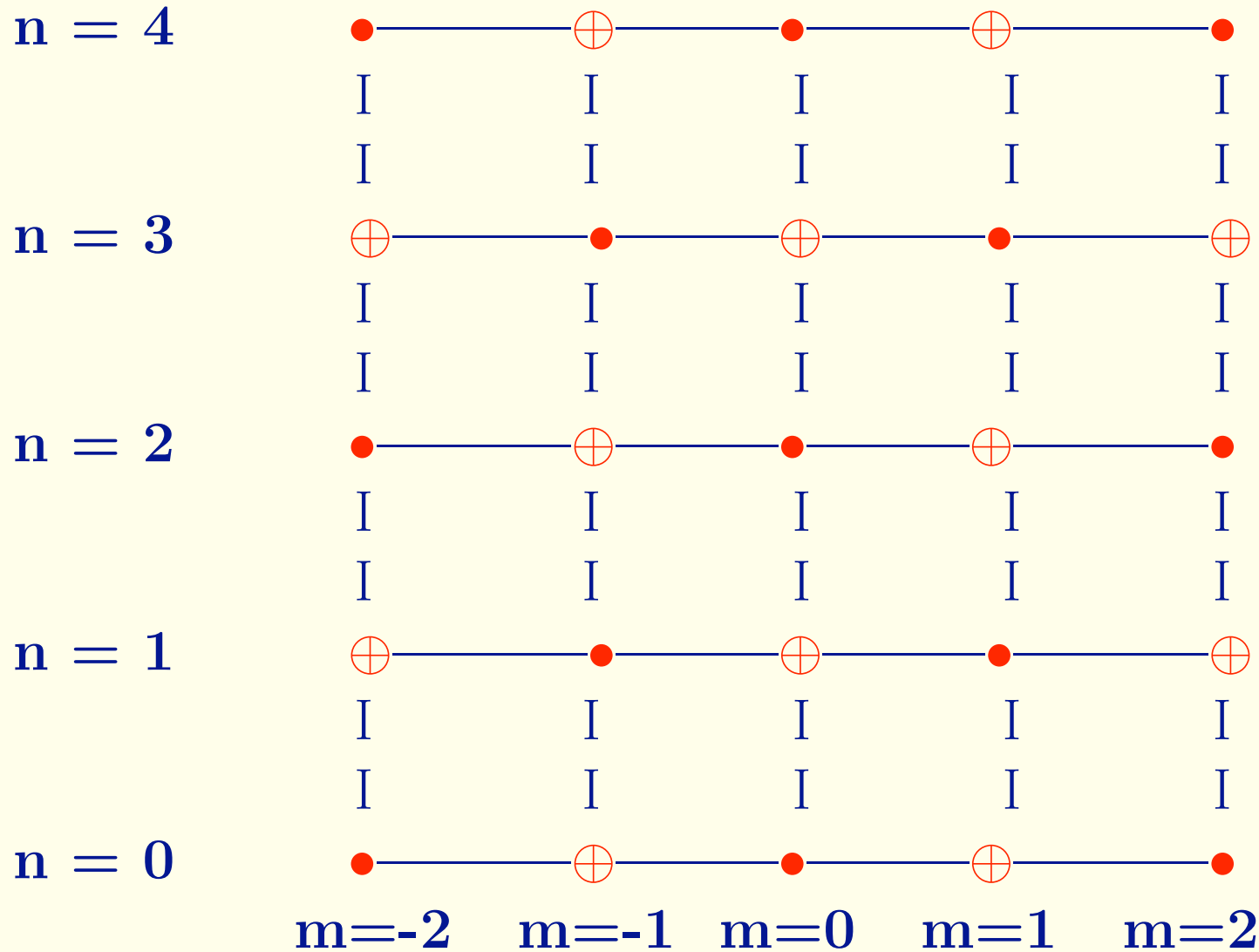


The evaluation of the equation at point \bullet involves values of the variable at points \oplus . Solving for Y_m^{n+1} thus requires

$$Y_m^{n-1}, \quad Y_{m-1}^n \quad \text{and} \quad Y_{m+1}^n.$$

The leapfrog scheme *splits the grid* into two independent sub-grids.

Grid Splitting



The finite difference grid splits into two sub-grids.

Steps must be taken to avoid divergence of the two solutions.

Recall the **(CTCS)** finite difference approximation:

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Assuming that we know the solution up to time $n\Delta t$, the values at time $(n + 1)\Delta t$ can be calculated, and the **solution advanced by one timestep** in this way.

Then the whole procedure can be repeated to advance the solution to $(n + 2)\Delta t$, and so on.

Question:

Under what conditions does the solution of the finite difference equation approximate that of the original differential equation?



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Substituting this in the equation, defining $\mu = c\Delta t/\Delta x$ and dividing by a common factor gives

$$A^2 + (2i\mu \sin k\Delta x)A - 1 = 0$$

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Substituting this in the equation, defining $\mu = c\Delta t/\Delta x$ and dividing by a common factor gives

$$A^2 + (2i\mu \sin k\Delta x)A - 1 = 0$$

This is a quadratic for the amplitude A , with solutions

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We write

$$A_{\pm} = -i\sigma \pm \sqrt{1 - \sigma^2} \quad \text{where} \quad \sigma \equiv \mu \sin k\Delta x$$

We consider in turn the two cases.

Case I: $|\mu| \leq 1$

The quantity under the square-root sign is positive, so the modulus of A is given by

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The two values of the phase are

$$\psi_1 = -\arcsin \sigma$$

and

$$\psi_2 = \pi - \psi_1.$$

The solution of the equation may now be written

$$Y_m^n = \left[D \exp(i\psi_1 n) + E \exp[i(-\psi_1 + \pi)n] \right] \exp(ikm\Delta x)$$

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Exercise:

Check in detail the algebra leading to this solution.

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If the ratio μ is small, the **physical mode solution** is given approximately by

$$Y \approx a \exp[ik(m\Delta x - cn\Delta t)]$$

which is just the **analytical solution**.

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In this simple case, we can **eliminate the computational mode**. In general, it is much more difficult.

Case II: $|\mu| > 1$

Recall that the roots of the quadratic are

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This phenomenon is called **computational instability.**

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We thus require that $|\mu| \leq 1$. This condition for stability is known as the CFL Criterion:

$$\frac{c\Delta t}{\Delta x} \leq 1$$

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Thus, **halving** the grid size in a two dimensional domain results in an **eightfold increase** in computation time.

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The semi-Lagrangian algorithm has enabled us to integrate the primitive equations using a **time step of 15 minutes**.

This can be compared to a typical timestep of 2.5 minutes for Eulerian schemes.

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We discuss semi-Lagrangian schemes in a later lecture.

End of §3.2.2