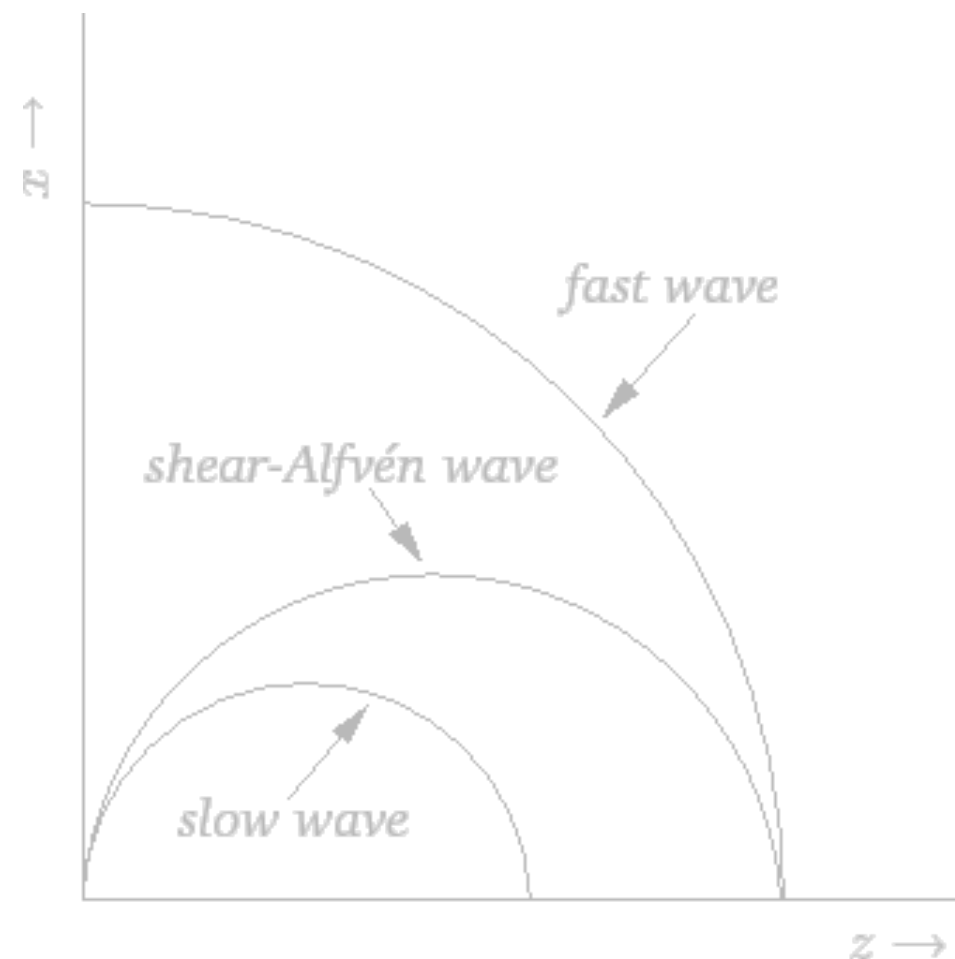


# How to Write a 1-D MHD Code using Finite Difference?



# Outline

- **Review of 1-D MHD equations and normalization**
- **Initial and Boundary Conditions**
- **Put everything together**
- **Once through the mhd.m code**

# 1-D MHD equations

Put the equations together:

Plasma variables

$$\rho, u_x, u_y, p$$

Field variables

$$B_y, E_z, J_z$$

$$\frac{\partial}{\partial t} \rho = - \nabla \cdot \rho \mathbf{u}$$

$$\rho \frac{D\mathbf{u}}{Dt} = - \nabla p + \mathbf{J} \times \mathbf{B}$$

$$p = \frac{\beta_0}{2} \rho^\gamma$$

$$\frac{\partial \mathbf{B}}{\partial t} = - \nabla \times \mathbf{E}$$

$$\mathbf{E} = - \mathbf{u} \times \mathbf{B}$$

$$\mathbf{J} = \nabla \times \mathbf{B}$$

$$\frac{\partial}{\partial t} \rho = - \frac{\partial}{\partial x} (\rho u_x)$$

$$\frac{\partial u_x}{\partial t} = - u_x \frac{\partial u_x}{\partial x} + \frac{1}{\rho} \left( - \frac{\partial p}{\partial x} - J_z B_y \right)$$

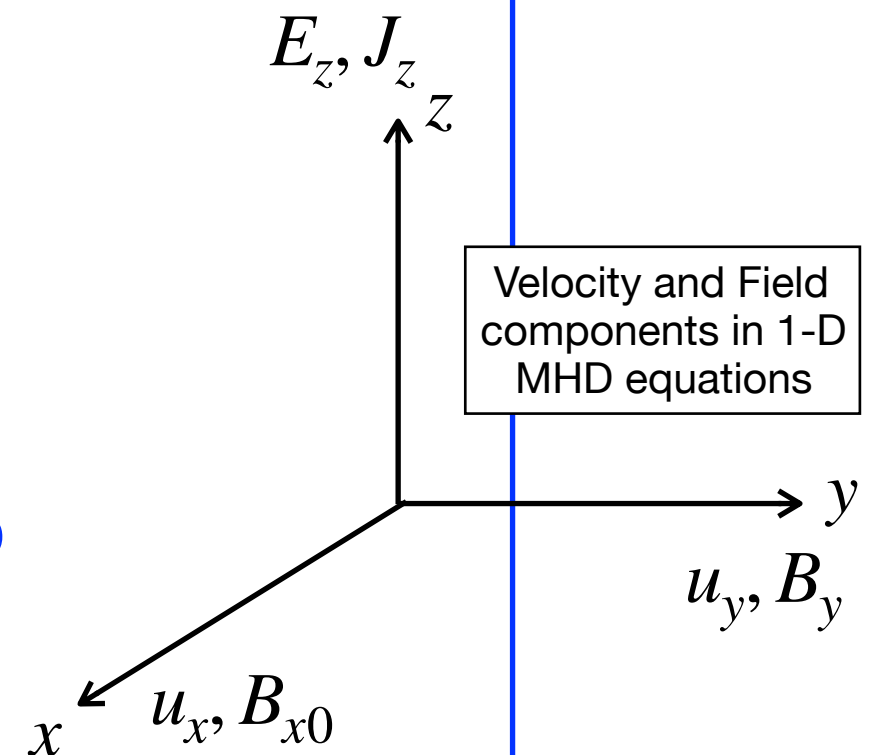
$$\frac{\partial u_y}{\partial t} = - u_x \frac{\partial u_y}{\partial x} + \frac{1}{\rho} J_z B_{x0}$$

$$p = \frac{\beta_0}{2} \rho^\gamma$$

$$\frac{\partial B_y}{\partial t} = \frac{\partial E_z}{\partial x}$$

$$E_z = - u_x B_y + u_y B_{x0}$$

$$J_z = \frac{\partial B_y}{\partial x}$$



# Spatial Derivatives - Finite Difference

1-D MHD equations

$$\frac{\partial}{\partial t} \rho = - \frac{\partial}{\partial x} (\rho u_x)$$

$$\frac{\partial u_x}{\partial t} = - u_x \frac{\partial u_x}{\partial x} + \frac{1}{\rho} \left( - \frac{\partial p}{\partial x} - J_z B_y \right)$$

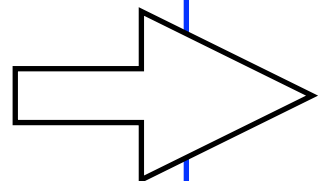
$$\frac{\partial u_y}{\partial t} = - u_x \frac{\partial u_y}{\partial x} + \frac{1}{\rho} J_z B_{x0}$$

$$\frac{\partial B_y}{\partial t} = \frac{\partial E_z}{\partial x}$$

$$E_z = - u_x B_y + u_y B_{x0}$$

$$J_z = \frac{\partial B_y}{\partial x}$$

$$p = \frac{\beta_0}{2} \rho^\gamma$$



Finite-Difference Approximations

$$\left. \frac{\partial \rho}{\partial t} \right|_i = \frac{u_{x,i+1}^n \rho_{i+1}^n - u_{x,i-1}^n \rho_{i-1}^n}{2\Delta x}$$

$$\left. \frac{\partial u_x}{\partial t} \right|_i = - u_{x,i} \frac{u_{x,i+1}^n - u_{x,i-1}^n}{2\Delta x} + \frac{1}{\rho_i^n} \left( \frac{p_{i+1}^n - p_{i-1}^n}{2\Delta x} - J_{z,i}^n B_{y,i}^n \right)$$

$$\left. \frac{\partial u_y}{\partial t} \right|_i = - u_{x,i} \frac{u_{y,i+1}^n - u_{y,i-1}^n}{2\Delta x} + \frac{1}{\rho_i^n} J_{z,i}^n B_{x0}$$

$$\left. \frac{\partial B_y}{\partial t} \right|_i = - \frac{E_{z,i+1}^n - E_{z,i-1}^n}{2\Delta x}$$

$$E_{z,i}^n = - u_{x,i}^n B_{y,i}^n + u_{y,i}^n B_{x0}$$

$$J_{z,i}^n = \frac{B_{y,i+1}^n - B_{y,i-1}^n}{2\Delta x}$$

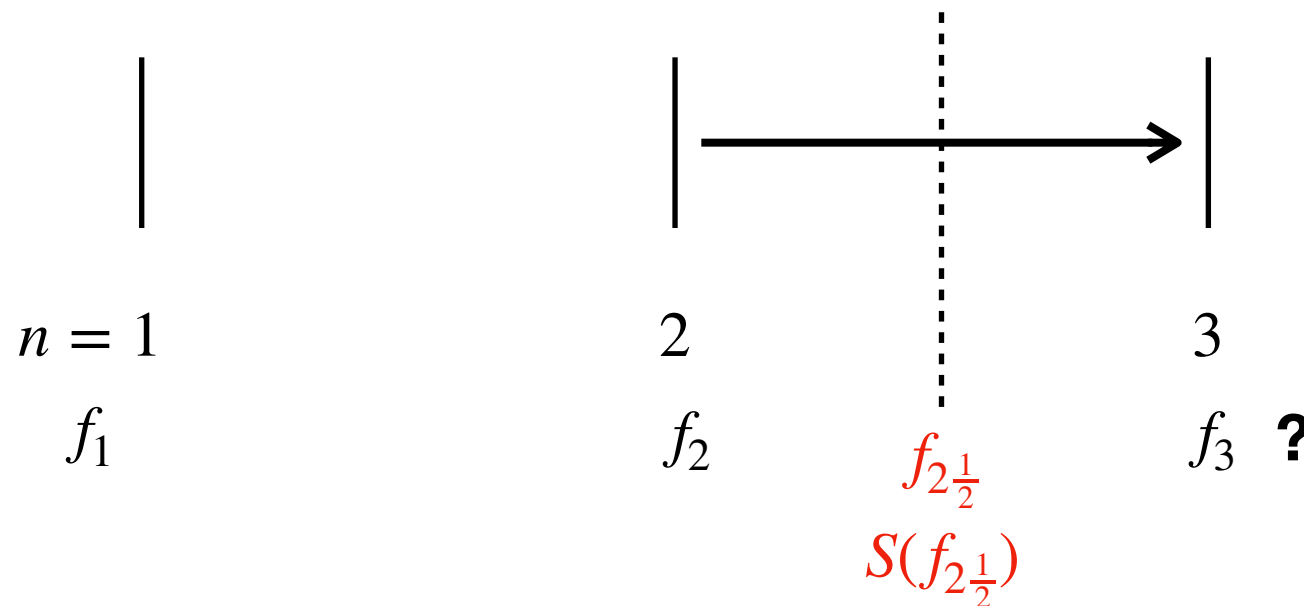
$$p_{z,i}^n = \frac{\beta_0}{2} (\rho_i^n)^\gamma$$

# Time Evolution - Leapfrog Trapezoidal

There are many choices of getting a viable second order time stepping method, we will use a kind of “predictor - corrector” scheme named “*Leapfrog Trapezoidal*” (LT) method in the following discussions, which uses two sub-steps for the time evolution

Assume that  $f$  is known at  $t_1$  and  $t_2$ , corresponding to  $n=1$  and  $n=2$ , we want to calculate  $f$  at  $t_3$  using the differential equation

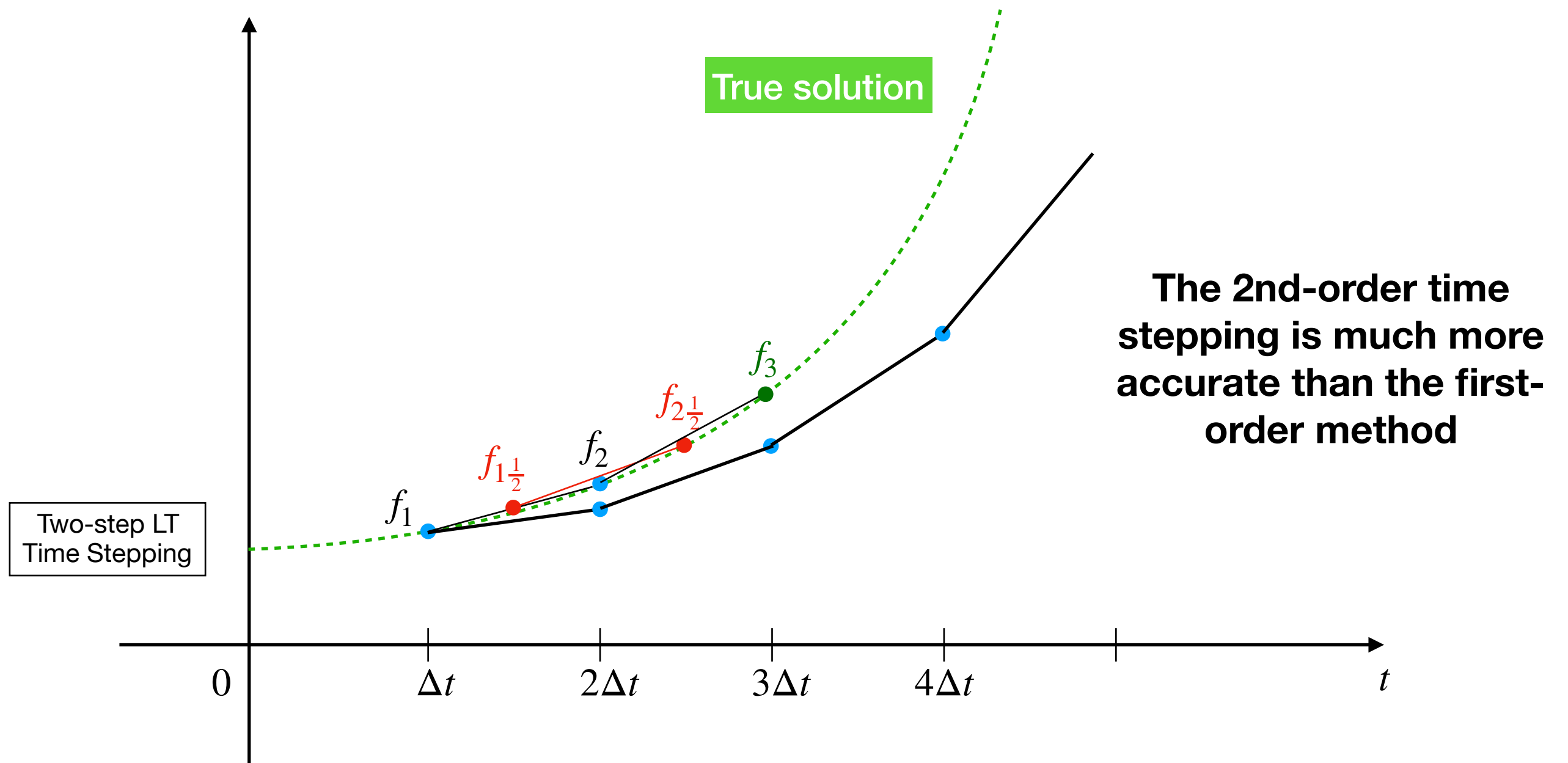
$$\frac{\partial f}{\partial t} = S(f)$$



- Going from  $f_2$  to  $f_3$  using  $f_2$  and  $S(f_2)$  gives the first order time stepping
- If we know  $f_{2.5}$  and  $S(f_{2.5})$ , we get something like a central difference scheme for  $f_3$ :

$$\frac{f_i^{n+1} - f_i^n}{\Delta t} = S(f_i^{n+\frac{1}{2}}) \implies f_i^{n+1} = f_i^n + \Delta t S(f_i^{n+\frac{1}{2}})$$

# Interpretation of the LT method



$$f^{n+\frac{1}{2}} = \frac{1}{2}(f^n + f^{n-1}) + \Delta t \cdot S(f^n)$$

$$f^{n+1} = f^n + \Delta t \cdot S(f^{n+\frac{1}{2}})$$

# Leapfrog Trapezoidal time stepping

The LT algorithm for time stepping goes like

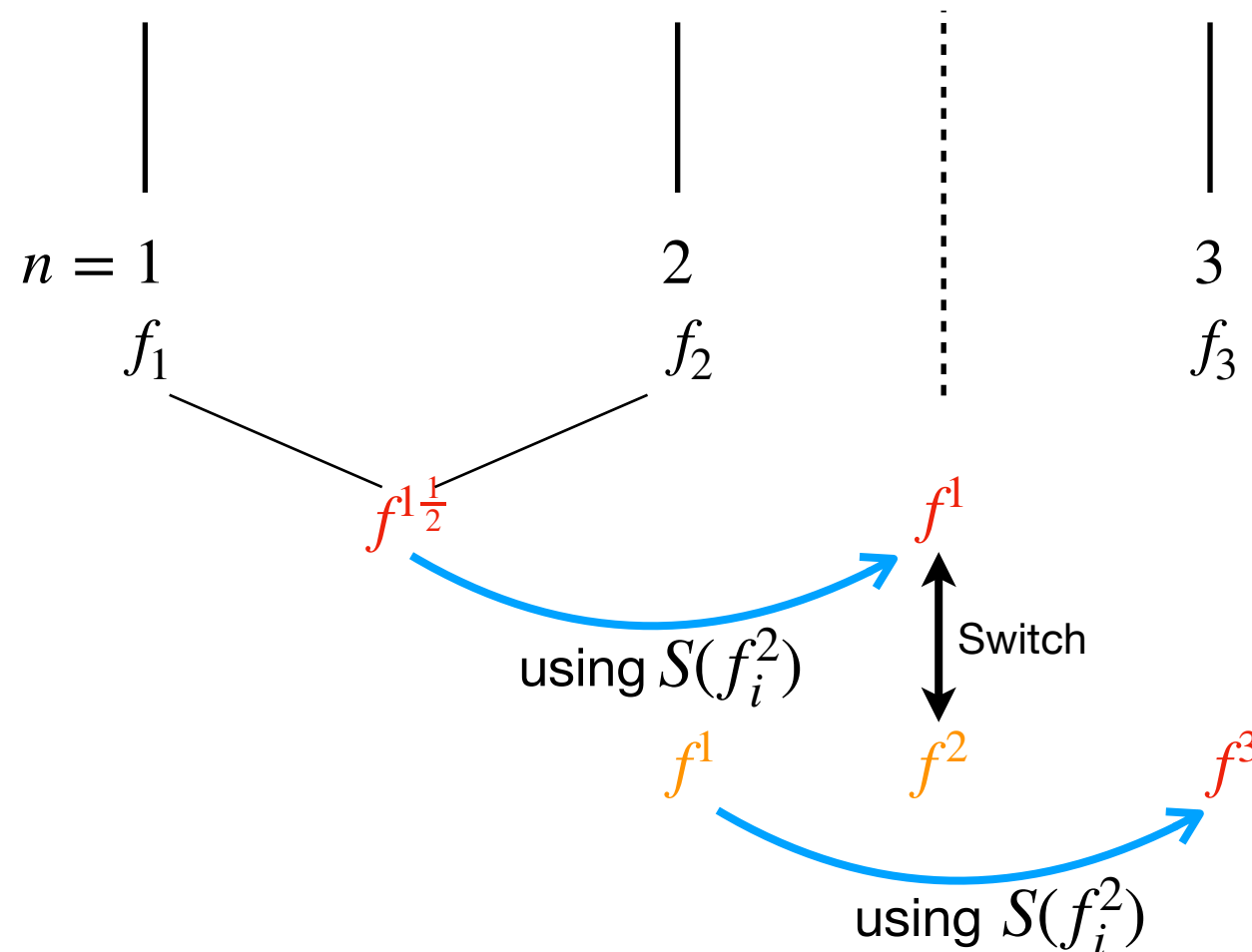
Two-step LT  
Time Stepping

$$f^{n+\frac{1}{2}} = \frac{1}{2}(f^n + f^{n-1}) + \Delta t \cdot S(f^n)$$

$$f^{n+1} = f^n + \Delta t \cdot S(f^{n+\frac{1}{2}})$$

$n = 2, 3, 4, \dots$

In general we need to store the variables (arrays) at each new time step (and each half time step like  $f^{2.5}$ ), it could be a waste of resource while the scale of the simulation is enormous. So we do something like this to get around:



Algorithm used in code

STEP 1:

$$f^1 = \frac{1}{2}(f^1 + f^2) + \Delta t \cdot S(f^2)$$

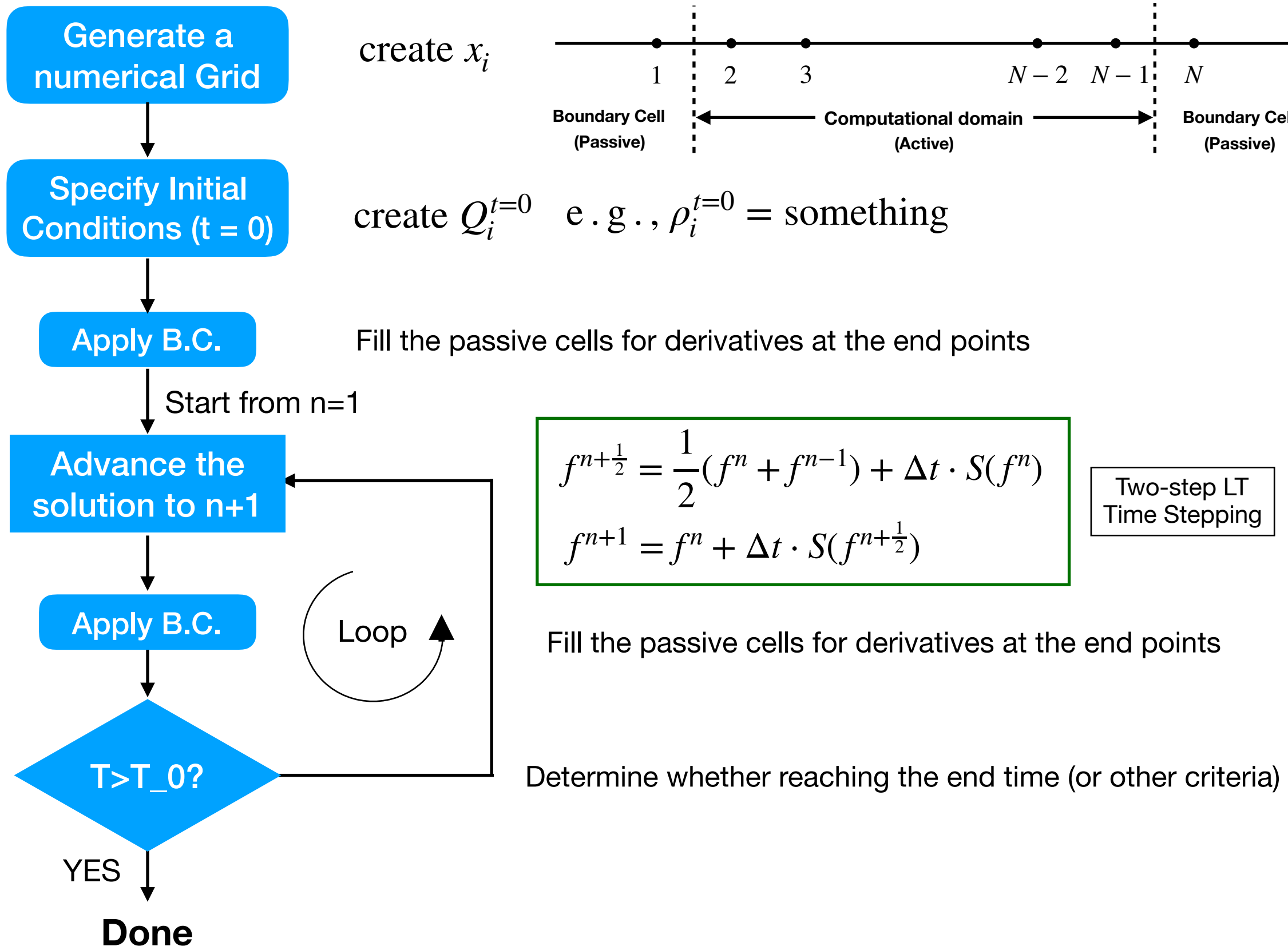
STEP 2:

switch  $f_1$  and  $f_2$

STEP 3:

$$f^2 = f^1 + \Delta t \cdot S(f^2)$$

# Build the code w/ predictor-corrector steps

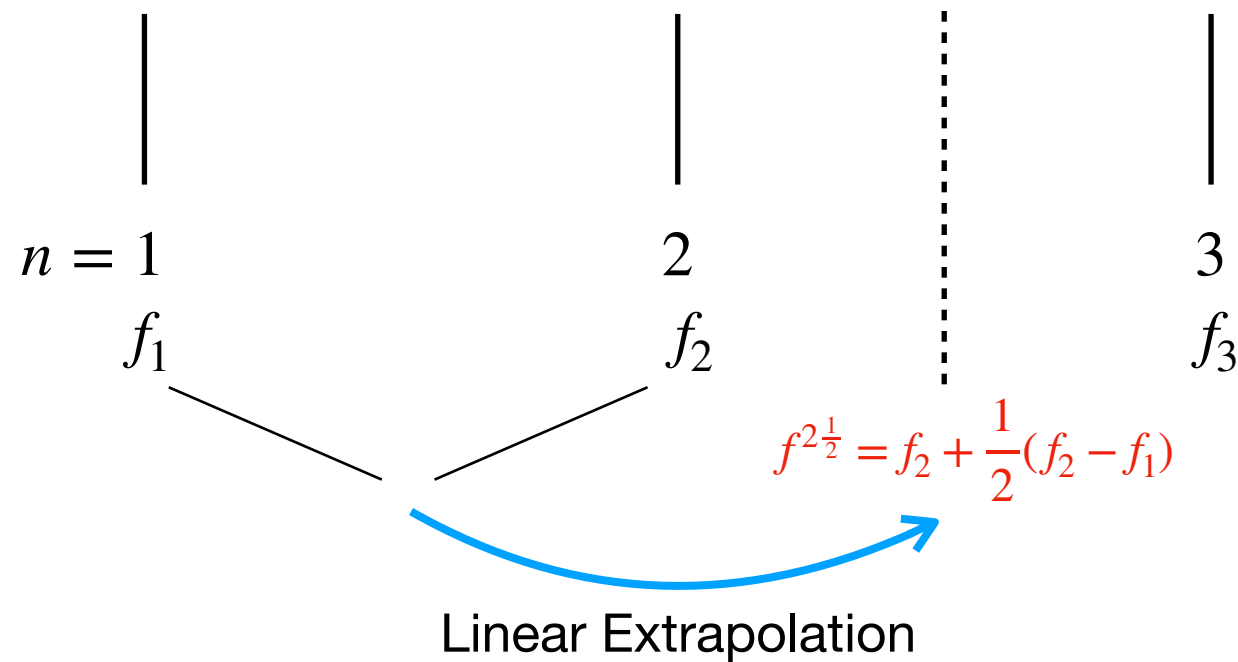




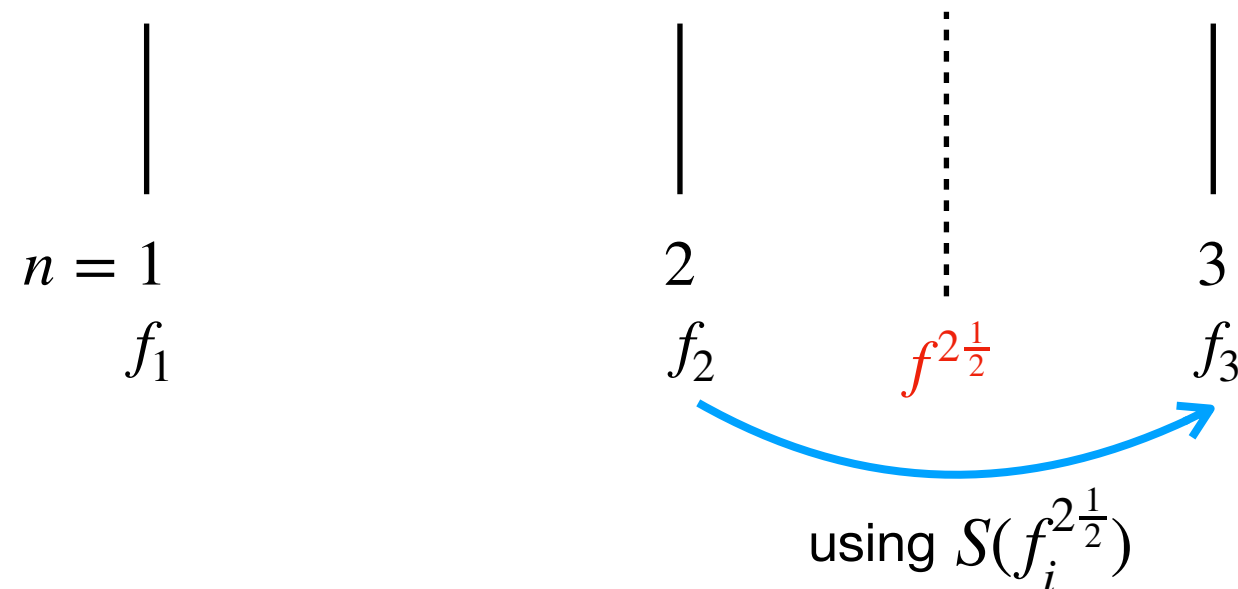
# One-Step predictor-corrector method

## Adams-Bashforth Method

The calculation of  $f_{2.5}$  is the ***predictor*** step - which is a linear extrapolation



Going from  $f_2$  to  $f_3$  is the same ***corrector*** step as leapfrog trapezoid method



# Energy in an MHD code (magnetic)

The magnetic energy equation is basically the Poynting theorem applied to the plasma:

$$\begin{aligned}
 \frac{\partial \mathbf{B}}{\partial t} = -\nabla \times \mathbf{E} & \xrightarrow{\cdot \mathbf{B}} \mathbf{B} \cdot \frac{\partial \mathbf{B}}{\partial t} = -\mathbf{B} \cdot \nabla \times \mathbf{E} \xrightarrow{\nabla \cdot (\mathbf{E} \times \mathbf{B}) = \mathbf{B} \cdot \nabla \times \mathbf{E} - \mathbf{E} \cdot \nabla \times \mathbf{B}} \\
 & \longrightarrow \frac{\partial}{\partial t} \left( \frac{1}{2} B^2 \right) = -\mathbf{E} \cdot \nabla \times \mathbf{B} - \nabla \cdot (\mathbf{E} \times \mathbf{B}) \xrightarrow[\mathbf{J} = \nabla \times \mathbf{B}]{\text{Ampere's law}} \\
 & = -\mathbf{E} \cdot \mathbf{J} - \nabla \cdot (\mathbf{E} \times \mathbf{B})
 \end{aligned}$$

So the volume-integrated magnetic energy equation is written as

$$\boxed{\text{Magnetic Energy}} \quad \int_V \frac{\partial}{\partial t} \left( \frac{1}{2} B^2 \right) dV = \int_V \left[ -\nabla \cdot (\mathbf{E} \times \mathbf{B}) - \mathbf{E} \cdot \mathbf{J} \right] dV$$

And

$$\boxed{\text{Kinetic Energy}} \quad \int_V \frac{d}{dt} \left( \frac{1}{2} \rho u^2 \right) dV = \int_V \left[ -\nabla \cdot \left( \frac{1}{2} \rho u^2 \mathbf{u} \right) - \mathbf{u} \cdot \nabla p + \mathbf{J} \cdot \mathbf{E} \right] dV$$

Converts magnetic energy to/from kinetic energy

$$\boxed{\text{Internal Energy}} \quad \int_V \frac{\partial}{\partial t} \left( \frac{p}{\gamma - 1} \right) dV = \int_V \left[ -\nabla \cdot \left( \frac{\gamma p}{\gamma - 1} \mathbf{u} \right) + \mathbf{u} \cdot \nabla p \right] dV$$

Converts internal energy to/from kinetic energy

# Energy in an MHD code (total)

Adding the three energy equations together

$$\int_V \frac{d}{dt} \left( \frac{1}{2} \rho u^2 + \frac{p}{\gamma - 1} + \frac{1}{2} B^2 \right) dV = \int_V \left[ -\nabla \cdot \left( \frac{1}{2} \rho u^2 \mathbf{u} \right) - \nabla \cdot \left( \frac{\gamma p}{\gamma - 1} \mathbf{u} \right) - \nabla \cdot (\mathbf{E} \times \mathbf{B}) \right] dV$$

Total Energy

$$= - \int_V \nabla \cdot \left[ \left( \frac{1}{2} \rho u^2 \mathbf{u} \right) + \left( \frac{\gamma p}{\gamma - 1} \mathbf{u} \right) + (\mathbf{E} \times \mathbf{B}) \right] dV$$

Divergence  
Theorem

$$\xrightarrow{\int_V \nabla \cdot \mathbf{A} dV = \oint_S \mathbf{A} \cdot d\mathbf{S}}$$

$$= - \oint_S d\mathbf{S} \cdot \left[ \left( \frac{1}{2} \rho u^2 \mathbf{u} \right) + \left( \frac{\gamma p}{\gamma - 1} \mathbf{u} \right) + (\mathbf{E} \times \mathbf{B}) \right]$$

Given certain boundary conditions the surface integral  $\oint_S$  may go to zero, leading to total **energy conservation**

$$\mathcal{E}_T = \frac{1}{2} \rho u^2 + \frac{p}{\gamma - 1} + \frac{1}{2} B^2$$

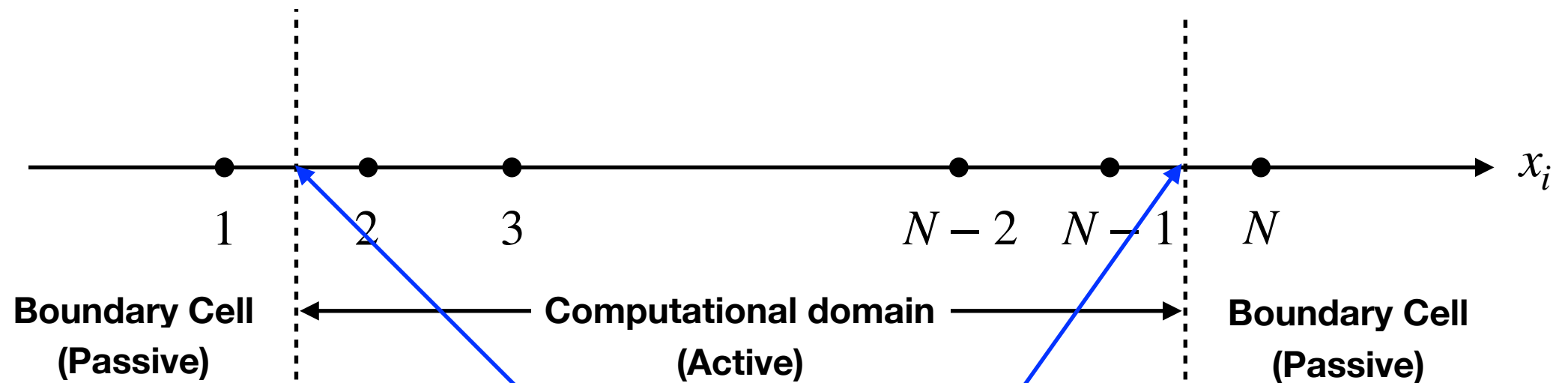
Total Energy

And the following term is called the total energy flux

$$\mathbf{F}_{\mathcal{E}} = \frac{1}{2} \rho u^2 \mathbf{u} + \frac{\gamma p}{\gamma - 1} \mathbf{u} + \mathbf{E} \times \mathbf{B}$$

Total Energy Flux

# Energy Conservation in 1-D MHD code



$$\frac{d}{dt} \int_V \mathcal{E}_T dV = - \oint_S d\mathbf{S} \cdot \left( \frac{1}{2} \rho u^2 \mathbf{u} + \frac{\gamma p}{\gamma - 1} \mathbf{u} + \mathbf{E} \times \mathbf{B} \right) \quad \text{Total Energy Flux}$$

If the total energy flux is zero on the boundary grid cells, then the total energy is conserved

$$\frac{d}{dt} \int_V \mathcal{E}_T dV = 0$$

This is called energy-conserving boundary conditions for MHD

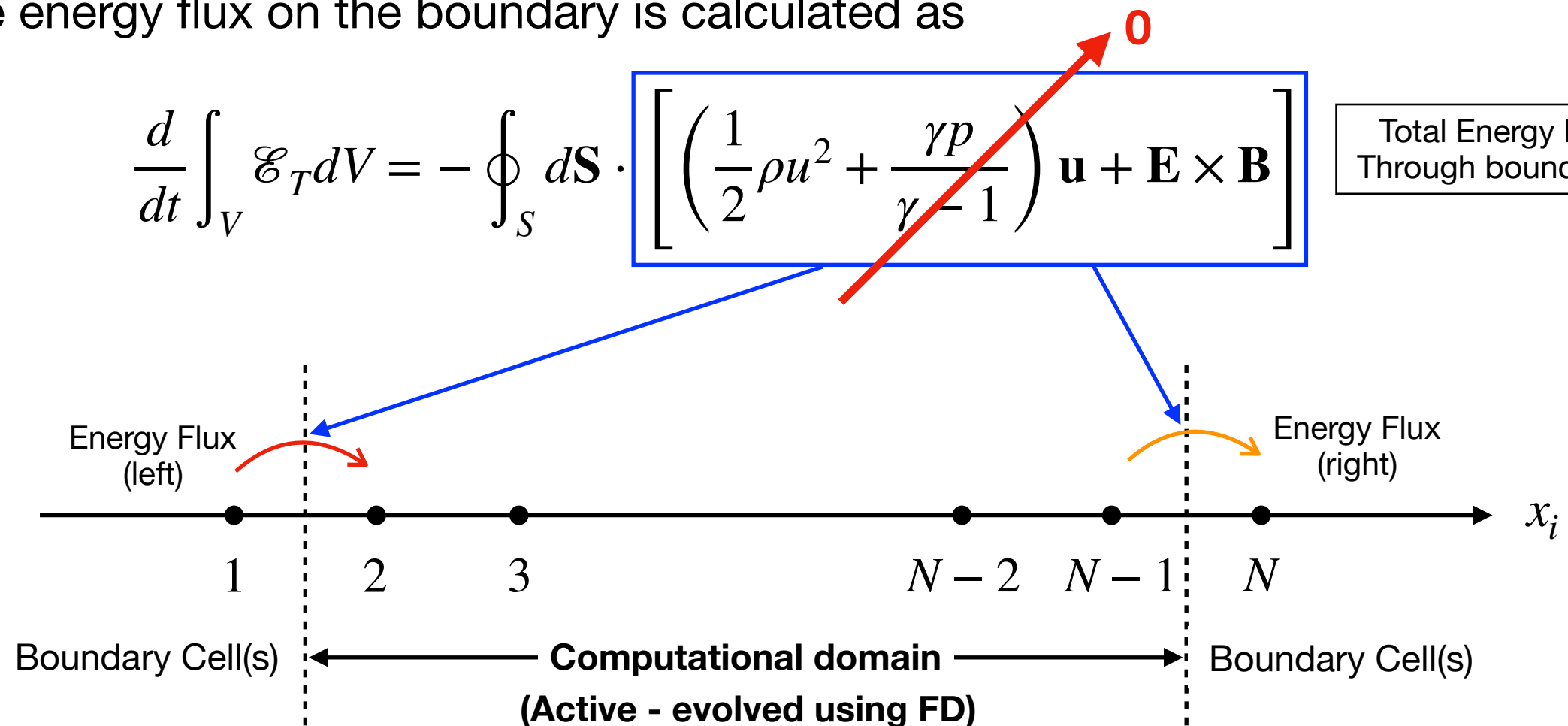
# Boundary Conditions in 1-D MHD code

Another type of boundary conditions that conserves total energy is called “**Hard Wall**”

Recall the energy flux on the boundary is calculated as

$$\frac{d}{dt} \int_V \mathcal{E}_T dV = - \oint_S d\mathbf{S} \cdot \left[ \left( \frac{1}{2} \rho u^2 + \frac{\gamma p}{\gamma - 1} \right) \mathbf{u} + \mathbf{E} \times \mathbf{B} \right]$$

Total Energy Flux Through boundaries



So if we have:

The diagram shows a 1-D computational domain along the  $x$  axis. The domain is divided into three regions: Boundary Cell(s) on the left, the Computational domain in the middle, and Boundary Cell(s) on the right. The computational domain is bounded by dashed vertical lines. The left boundary is at the interface between cell 1 and cell 2, and the right boundary is at the interface between cell 2 and cell 3. The cells are labeled 1, 2, 3. A red arrow labeled  $\mathbf{E} \equiv 0$  points from the left boundary into the domain. A red arrow labeled  $\mathbf{u} \equiv 0$  points from the left boundary into the domain. A large white arrow points from the diagram to the equation:

$$\left[ \left( \frac{1}{2} \rho u^2 + \frac{\gamma p}{\gamma - 1} \right) \mathbf{u} + \mathbf{E} \times \mathbf{B} \right]_{1\frac{1}{2}} \equiv 0$$

Total energy flux is zero on boundary

# What does “Hard-wall” boundary mean?

The “hard-wall” boundary puts a constraint on the normal component of the velocity, i.e., there’s no flow velocity across the boundary interface (wall). Thus the mass flux through the interface is also expected to be zero - the total mass within the computation domain is conserved (to the scheme accuracy)

Let’s take a look at the mass equation

$$\frac{\partial}{\partial t} \rho = - \nabla \cdot \rho \mathbf{u} \xrightarrow{\int_V dV} \int_V \left( \frac{\partial}{\partial t} \rho = - \nabla \cdot \rho \mathbf{u} \right) dV$$

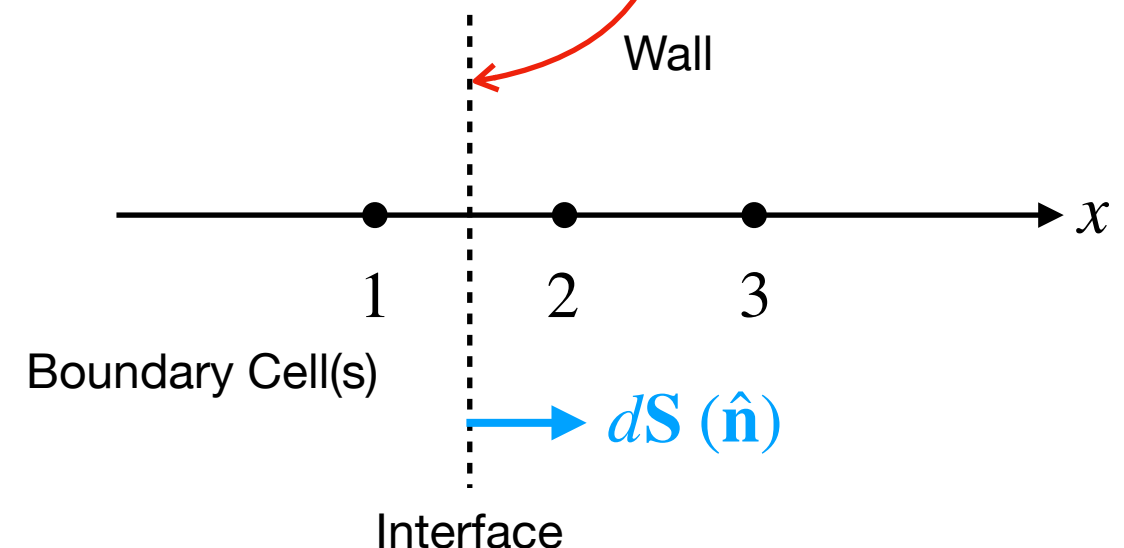
$$\longrightarrow \frac{\partial}{\partial t} \int_V \rho dV = - \int_V \nabla \cdot \rho \mathbf{u} dV \longrightarrow \frac{\partial}{\partial t} \int_V \rho dV = - \oint_S \rho \mathbf{u} \cdot d\mathbf{S}$$

So if  $\mathbf{u} \cdot d\mathbf{S} \equiv 0$

We have

$$\oint_S \rho \mathbf{u} \cdot d\mathbf{S} \equiv 0 \longrightarrow \frac{\partial}{\partial t} \int_V \rho dV = 0$$

**total mass unchanged**



# How to implement hard-wall boundary?

Hard-wall condition requires the normal component of the velocity across the wall is zero

**Hard wall:**  $u_n = \mathbf{u} \cdot \hat{\mathbf{n}} \equiv 0$   
normal

In the code, velocities are specified at grid points 1,2,3,... N. How's interface velocity calculated?

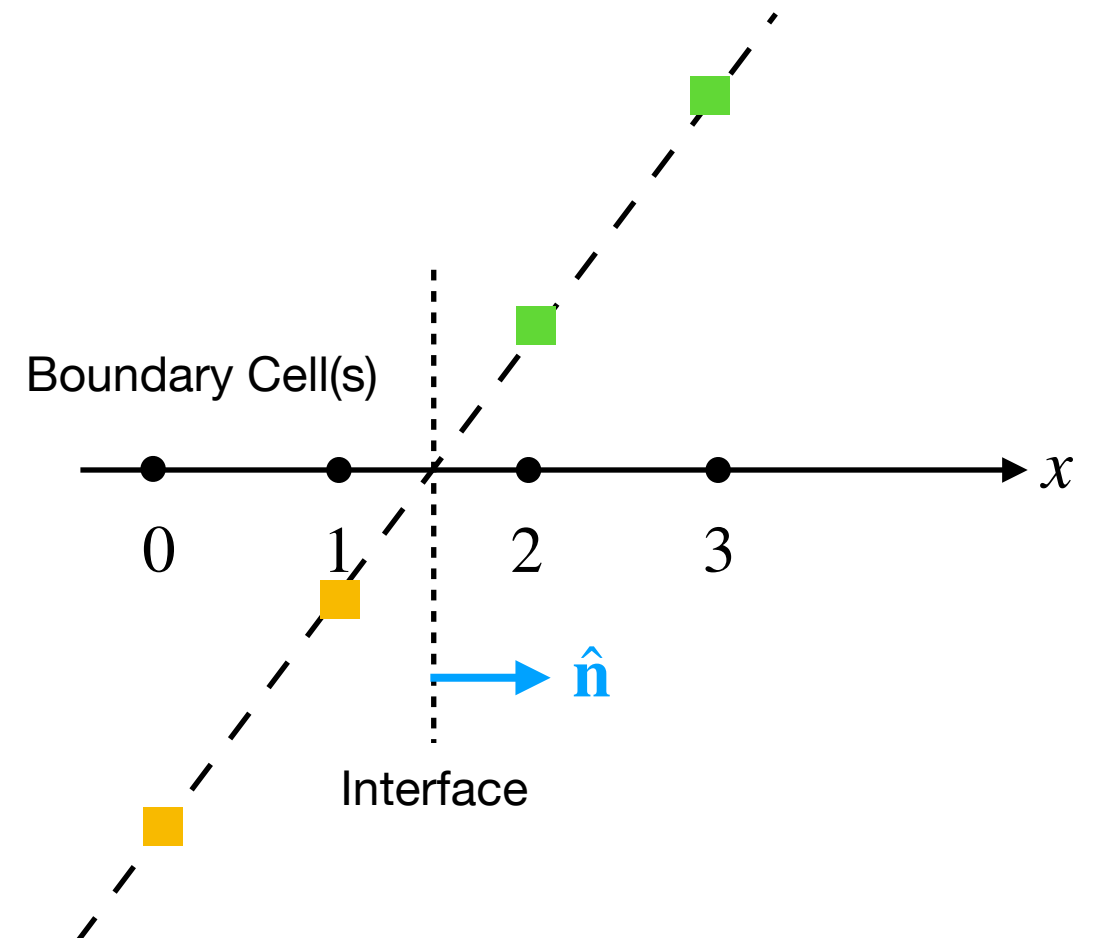
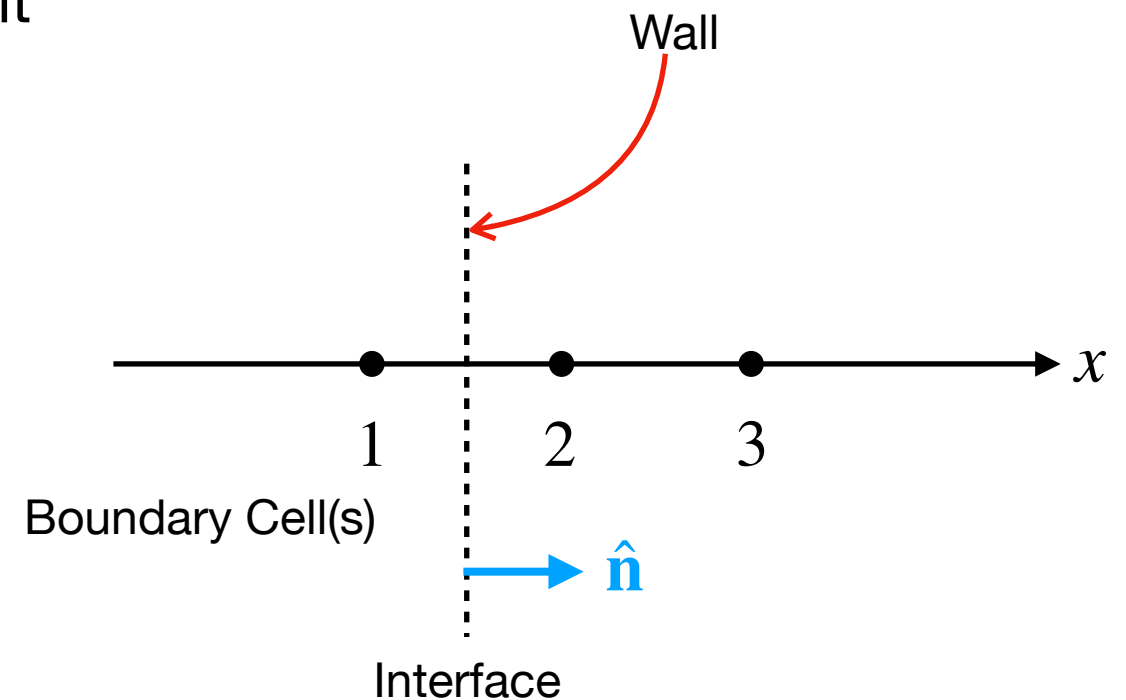
It's done based on the boundary cell(s):

If we set the boundary cells with values that are anti-symmetric with the active computational cells right next to the boundary, the “interpolated” value on the interface is going to be exactly zero (why?)

This is also called “**anti-symmetric**” boundary conditions. In the code, simply do:

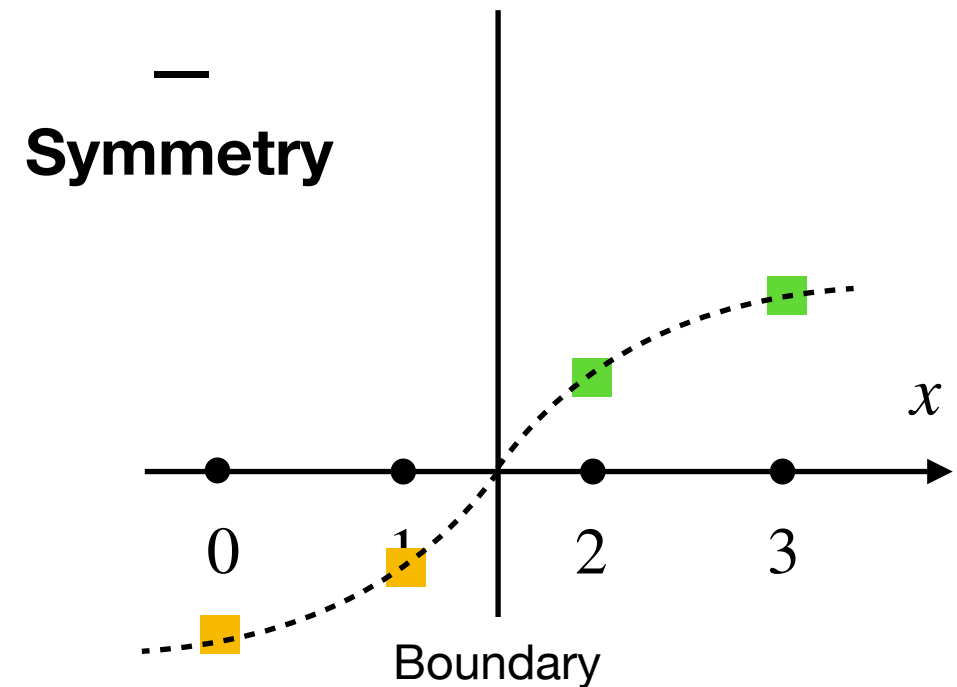
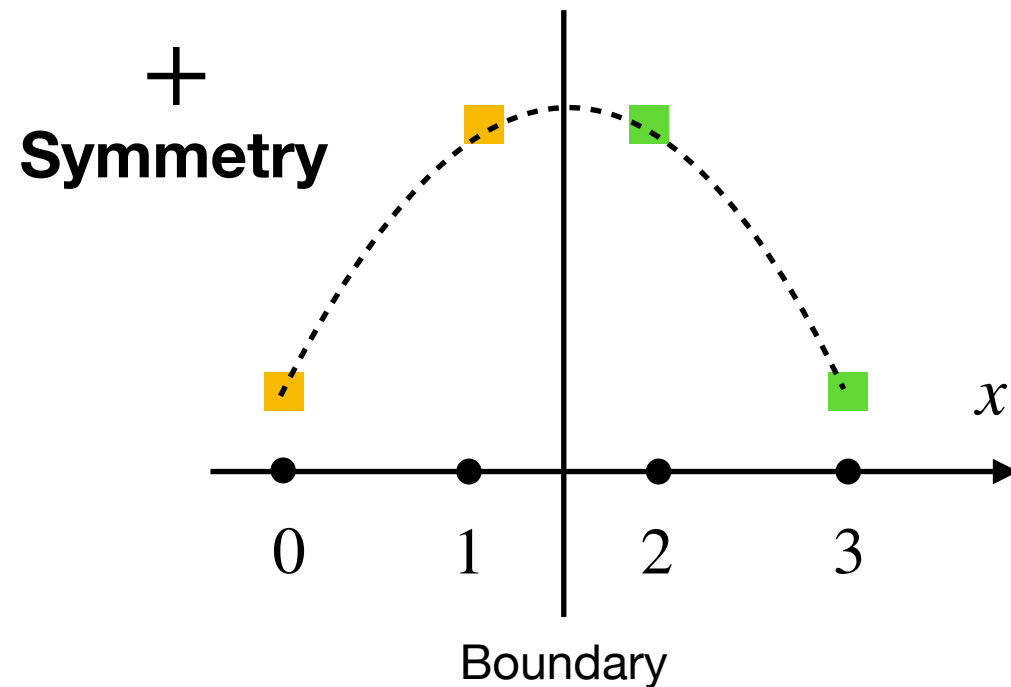
$$u_x(1) = -u_x(2)$$

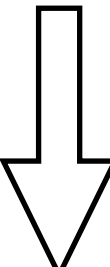
$$u_x(N) = -u_x(N-1)$$

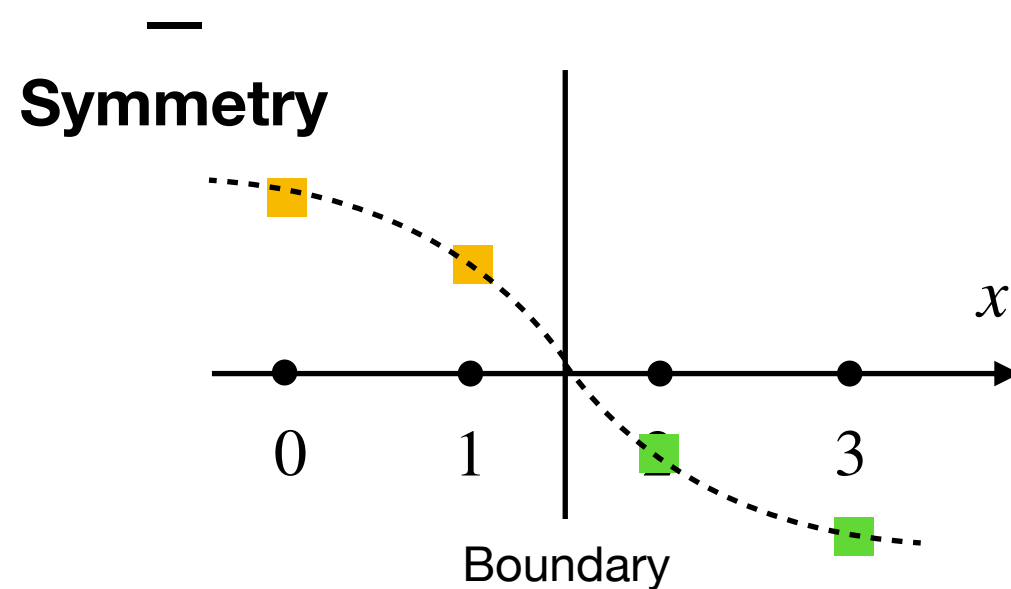


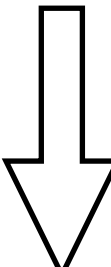
# Properties of symmetric boundaries

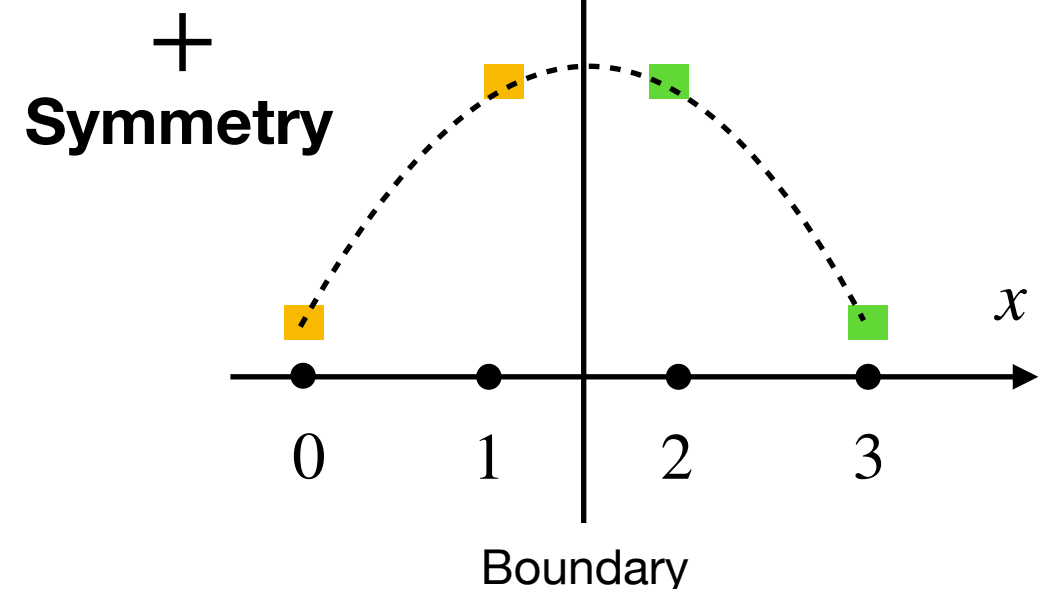
Two types of symmetries:



Consider  $\frac{\partial f}{\partial x}$   **reverses symmetry**



Consider  $\frac{\partial f}{\partial x}$   **reverses symmetry**





# Properties of symmetric boundaries

Now let's get the **full set** of boundary conditions for 1-D MHD equations, which conserves mass and energy:

**Normal velocity:** **Hard wall**  $\longrightarrow u_x (-)$

**Tangential electric field:** **Perfect conductor**  $\longrightarrow E_z (-)$

**Plasma density pressure:** **Equilibrium**  $\longrightarrow \rho, p (+)$

What about other vector components -  $B_y, u_y$

$$B_y : \quad \frac{\partial B_y}{\partial t} = \frac{\partial E_z}{\partial x} \quad \xrightarrow[\partial_x (-)]{E_z (-)} \quad B_y \sim (-)(-) \quad \longrightarrow \quad B_y (+)$$

This is because  $\frac{\partial}{\partial t}$  does not change symmetry

$$J_z : \quad J_z = \frac{\partial B_y}{\partial x} \quad \xrightarrow[\partial_x (-)]{B_y (+)} \quad J_z \sim (+)(-) \quad \longrightarrow \quad J_z (-)$$

Check the density equation:

$$\rho : \quad \frac{\partial}{\partial t} \rho = - \frac{\partial}{\partial x} (\rho u_x) \quad \xrightarrow[\partial_x (-)]{\rho (+) u_x (-)} \quad \rho \sim (+)(-)(-) \sim (+)$$

Symmetry  
consistent!

# Properties of symmetric boundaries

Also let's check the equation of  $u_x$ :

$$u_x : \frac{\partial u_x}{\partial t} = -u_x \frac{\partial u_x}{\partial x} + \frac{1}{\rho} \left( -\frac{\partial p}{\partial x} - J_z B_y \right)$$

Diagram illustrating the symmetry of the equation for  $u_x$ . The terms are grouped by symmetry signs in parentheses below them:

- $\frac{\partial u_x}{\partial t}$  is associated with  $(-)$ .
- $-u_x \frac{\partial u_x}{\partial x}$  is associated with  $(-)(-)(-)$  under a horizontal line, which is then associated with  $(-)$ .
- $\frac{1}{\rho}$  is associated with  $(+)$ .
- $-\frac{\partial p}{\partial x}$  is associated with  $(-)(+)$  under a horizontal line, which is then associated with  $(-)$ .
- $-J_z B_y$  is associated with  $(-)(+)$  under a horizontal line, which is then associated with  $(-)$ .

A large curved arrow labeled "Symmetry consistant!" points from the  $(-)$  under  $\frac{\partial u_x}{\partial t}$  to the  $(-)$  under the  $-\frac{\partial p}{\partial x}$  and  $-J_z B_y$  terms.

$$u_y : \frac{\partial u_y}{\partial t} = -u_x \frac{\partial u_y}{\partial x} + \frac{1}{\rho} J_z B_{x0}$$

Diagram illustrating the symmetry of the equation for  $u_y$ . The terms are grouped by symmetry signs in parentheses below them:

- $\frac{\partial u_y}{\partial t}$  is associated with  $(-)(-)$  under a horizontal line, which is then associated with  $(-)$ .
- $-u_x \frac{\partial u_y}{\partial x}$  is associated with  $(+)(-)$  under a horizontal line, which is then associated with  $(-)$ .
- $\frac{1}{\rho} J_z B_{x0}$  is associated with  $(+)(-)$  under a horizontal line, which is then associated with  $(-)$ .

$\Rightarrow u_y (-)$

It is straightforward to show that  $u_y$  must be anti-symmetric to make the symmetry in the equation of  $u_y$  consistent

Also can see from Ohm's law;

$$E_z = -u_x B_y + u_y B_{x0}$$

# Summary of Energy-Conserving boundaries

**Plasma variables:**  $\rho, p$  ( + )  $u_x, u_y$  ( - )

**Magnetic fields:**  $B_y$  ( + )

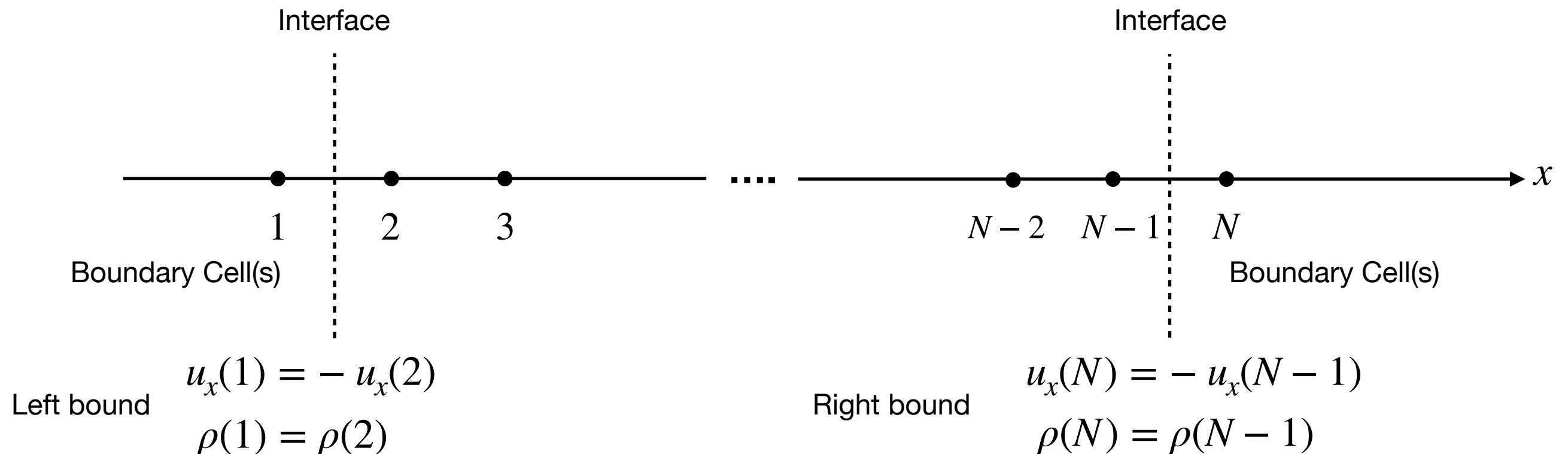
Primary  
Variables

**Electric fields:**  $E_z$  ( - )

**Currents:**  $J_z$  ( - )

Derived  
Variables

## How to implement in the code?



# Put Everything Together : 1-D MHD solver

## Finite-Difference Approximations

$$\begin{aligned}\frac{\partial \rho}{\partial t} \Big|_i &= \frac{u_{x,i+1}^n \rho_{i+1}^n - u_{x,i-1}^n \rho_{i-1}^n}{2\Delta x} \\ \frac{\partial u_x}{\partial t} \Big|_i &= -u_{x,i} \frac{u_{x,i+1}^n - u_{x,i-1}^n}{2\Delta x} \\ &\quad + \frac{1}{\rho_i^n} \left( \frac{p_{i+1}^n - p_{i-1}^n}{2\Delta x} - J_{z,i}^n B_{y,i}^n \right) \\ \frac{\partial u_y}{\partial t} \Big|_i &= -u_{x,i} \frac{u_{y,i+1}^n - u_{y,i-1}^n}{2\Delta x} + \frac{1}{\rho_i^n} J_{z,i}^n B_{x0} \\ \frac{\partial B_y}{\partial t} \Big|_i &= -\frac{E_{z,i+1}^n - E_{z,i-1}^n}{2\Delta x} \\ E_{z,i}^n &= -u_{x,i}^n B_{y,i}^n + u_{y,i}^n B_{x0} \\ J_{z,i}^n &= \frac{B_{y,i+1}^n - B_{y,i-1}^n}{2\Delta x} \\ p_{z,i}^n &= \frac{\beta_0}{2} (\rho_i^n)^\gamma\end{aligned}$$

Two-step LT  
Time Stepping

$$\frac{\partial f}{\partial t} = S(f)$$

$$\begin{aligned}f^{n+\frac{1}{2}} &= \frac{1}{2}(f^n + f^{n-1}) + \Delta t \cdot S(f^n) \\ f^{n+1} &= f^n + \Delta t \cdot S(f^{n+\frac{1}{2}})\end{aligned}$$

$n = 2, 3, 4, \dots$

Boundary  
conditions

**Plasma variables:    Magnetic fields:**

$$\begin{aligned}\rho, p \quad ( + ) \quad & u_x, u_y \quad ( - ) \\ B_y \quad ( + )\end{aligned}$$

**Electric fields:    Currents:**

$$\begin{aligned}E_z \quad ( - ) \\ J_z \quad ( - )\end{aligned}$$

Once though the mhd.m