



Lecture 2

FDTD: THE FINITE DIFFERENCE TIME DOMAIN METHOD

The **finite-difference time-domain (FDTD) method** is a classical technique for the modeling of circuits, antennas and scattering problems.

The FDTD method is easy to implement and very flexible: it can apply to the modeling of **complex regions**, with **inhomogeneous medium** and the **mixed/time-dependent boundary conditions**. The solution is calculated in the **time domain**, thus permitting the study of transient behavior with pulsed source.

Originally proposed in the 1920's by A. Thom to solve hydrodynamic equations, it was applied to the electromagnetic modeling by **K. Yee** in **1966** for the first time.

The FDTD method is based on the **approximation of the partial derivatives appearing in the differential equations as finite differences.**

The implementation of the FDTD method involves three steps:

1. definition of a **grid of nodes** in the solution region
2. approximation of the partial derivatives as **finite differences**
3. **iterative solution** of the equation starting from the initial conditions

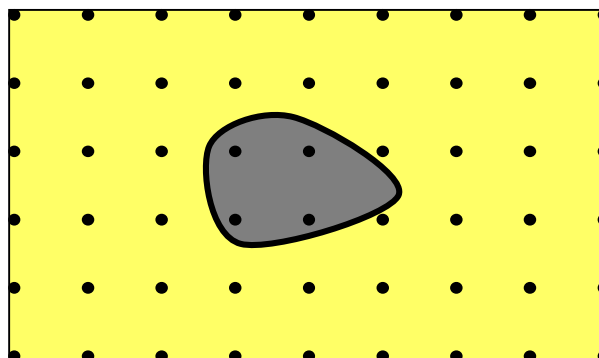
Reference books:

1. M.N.O. Sadiku, *Numerical techniques in Electromagnetics*, CRC Press, 2000.
2. A. Taflove and S.C. Hagness, *Computational Electrodynamics: The Finite-Difference Time-Domain Method*, Artech House, 2005.

The application of the FDTD method requires the definition of a grid of nodes in the whole investigation domain.

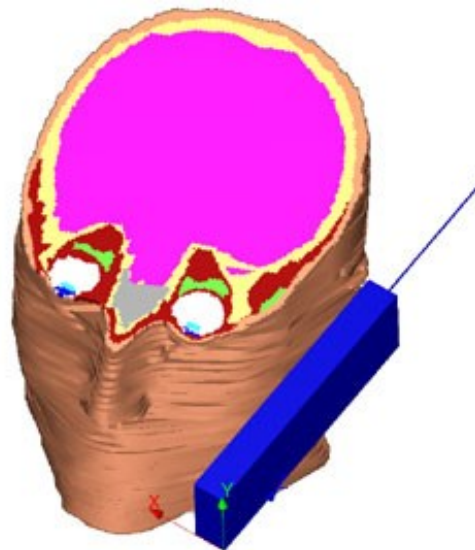
The spacing of the grid nodes is related to the **geometry** and to the **variation of the electromagnetic field**.

In general accurate results are obtained when the grid spacing is around $\lambda/20$ at the maximum frequency of interest.

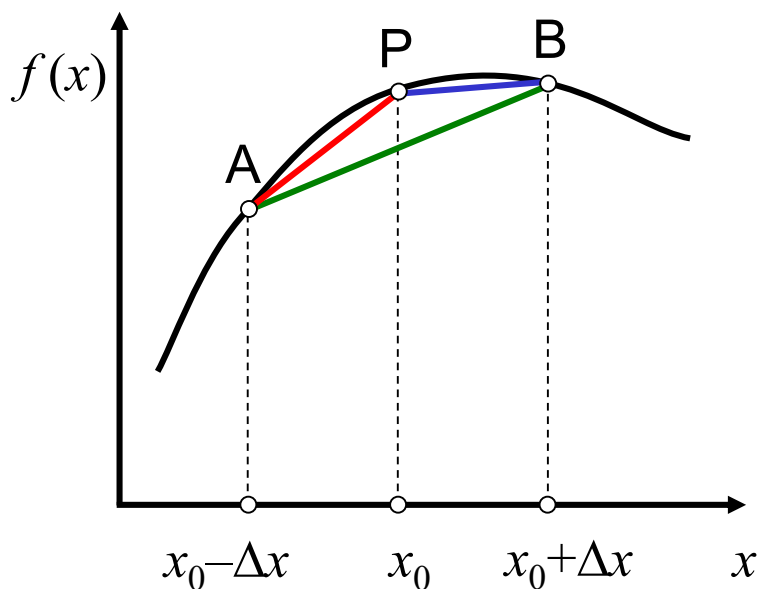


In some cases, the **sub-gridding technique** is adopted, to reduce the number of nodes.

It results very convenient in the modeling of **wire antennas** and PEC structures to arbitrarily oriented PEC geometries. This enables a much improved representation compared to common **staircase** modeling.



Given a function $f(x)$, its derivative in P can be approximated in different ways:



- FORWARD DIFFERENCE:

$$f'(x_0) \cong \frac{f(x_0 + \Delta x) - f(x_0)}{\Delta x}$$

- BACKWARD DIFFERENCE:

$$f'(x_0) \cong \frac{f(x_0) - f(x_0 - \Delta x)}{\Delta x}$$

- CENTRAL DIFFERENCE:

$$f'(x_0) \cong \frac{f(x_0 + \Delta x) - f(x_0 - \Delta x)}{2\Delta x}$$

Similarly, the **second derivative** can be approximated as:

$$\begin{aligned} f''(x_0) &\cong \frac{f'(x_0 + \Delta x/2) - f'(x_0 - \Delta x/2)}{\Delta x} \\ &\cong \frac{1}{\Delta x} \left(\frac{f(x_0 + \Delta x) - f(x_0)}{\Delta x} - \frac{f(x_0) - f(x_0 - \Delta x)}{\Delta x} \right) \\ &\cong \frac{f(x_0 + \Delta x) - 2f(x_0) + f(x_0 - \Delta x)}{(\Delta x)^2} \end{aligned}$$

By using this technique, the derivatives f' and f'' can be approximated in terms of the value of the function in a **discrete set of points**.



A more systematic approach to achieve the same result is based on the **Taylor's series**:

$$f(x_0 \pm \Delta x) = f(x_0) \pm \Delta x f'(x_0) + \frac{1}{2!} (\Delta x)^2 f''(x_0) \pm \frac{1}{3!} (\Delta x)^3 f'''(x_0) + \dots$$

Upon subtracting the expansion with sign + and the one with sign -, after dividing by $2\Delta x$, it results:

$$f'(x_0) \approx \frac{f(x_0 + \Delta x) - f(x_0 - \Delta x)}{2\Delta x} + \boxed{O(\Delta x^2)}$$

truncation error

Similarly, upon adding the expansion with sign + and the one with sign -, after dividing by $(\Delta x)^2$, it results:

$$f''(x_0) = \frac{f(x_0 + \Delta x) - 2f(x_0) + f(x_0 - \Delta x)}{(\Delta x)^2} + O(\Delta x^2)$$

It is noted that **forward and backward difference formulas have truncation error of $O(\Delta x)$** .

A simple example is useful to introduce the **concept of the iterative solution**.

Let us consider the following partial differential equation

$$k \frac{\partial \Phi}{\partial t} = \frac{\partial^2 \Phi}{\partial x^2}$$

(where k is constant).

We define a **space grid** with step Δx and a **time step** Δt (thus resulting $x=i \Delta x$ and $t=j \Delta t$, with $i, j=0,1,\dots$)

We approximate the time derivative with the **forward difference** and the space derivative with the **central difference**:

$$k \frac{\Phi(i, j+1) - \Phi(i, j)}{\Delta t} = \frac{\Phi(i+1, j) - 2\Phi(i, j) + \Phi(i-1, j)}{(\Delta x)^2}$$

It is observed that $\Phi(i, j)$ is a compact form for representing $\Phi(i \Delta x, j \Delta t)$

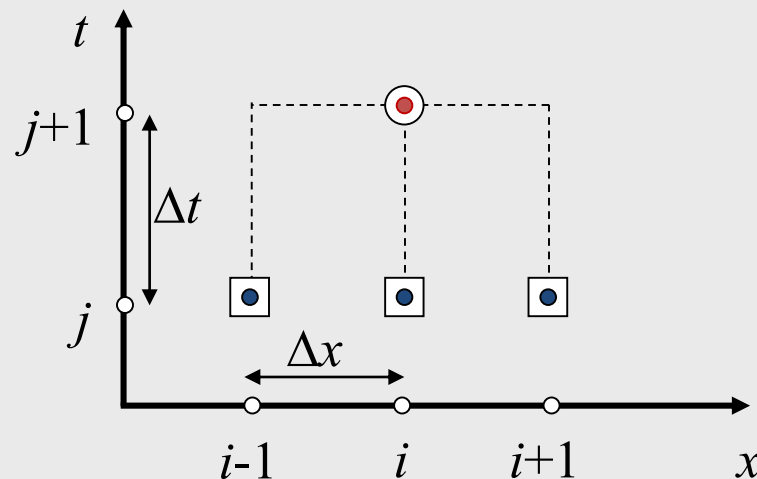
After re-ordering the equation, we obtain:

$$\Phi(i, j+1) = r \Phi(i+1, j) + (1-2r) \Phi(i, j) + r \Phi(i-1, j) \quad (\text{with } r = \frac{\Delta t}{k(\Delta x)^2}).$$

This expression represents an **explicit formula** to determine $\Phi(x, t+\Delta t)$ in terms of $\Phi(x, t)$.

COMPUTATIONAL SCHEME

- ⊙ value to determine
- ◻ known values of Φ



This explicit formula is **stable** only for certain values of r .

$$0 < r \leq \frac{1}{2}$$

STABILITY CONDITION

As r depends on the ratio between Δt and Δx , the stability condition sets the maximum time step for a given space grid.

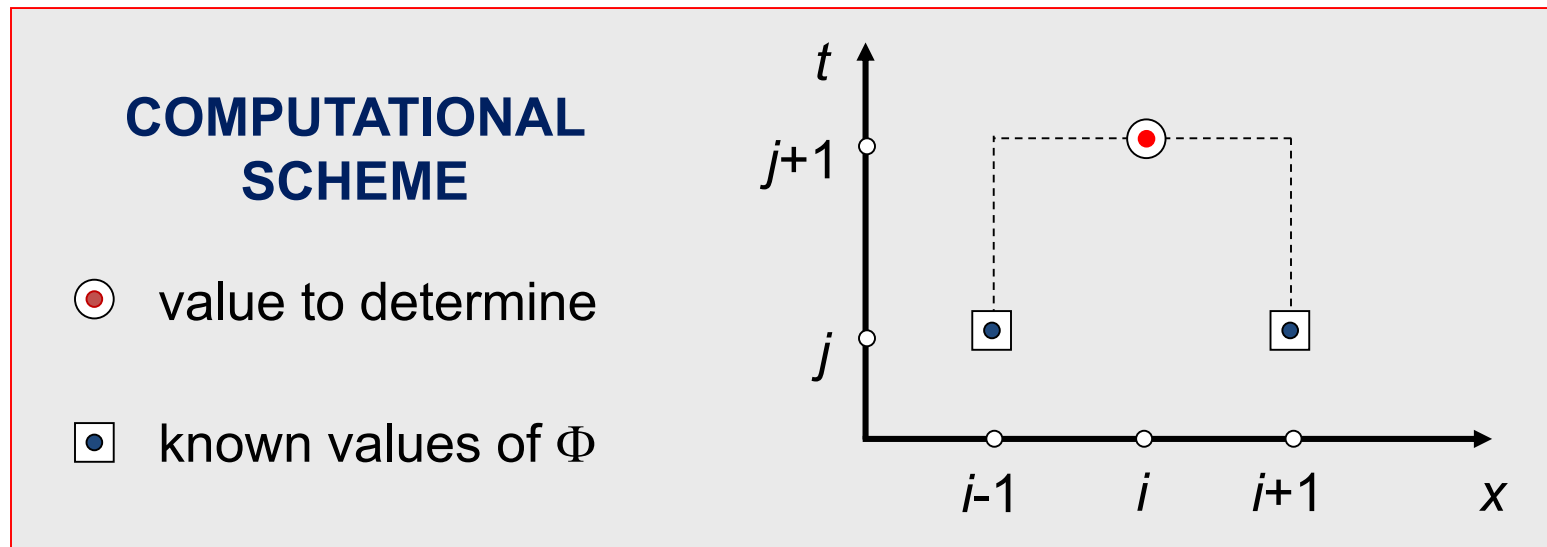
$$0 < \Delta t \leq \frac{1}{2} k(\Delta x)^2$$

NOTE: a larger value of r allows adopting a larger time step Δt .

When adopting $r=1/2$ (largest time step) it results:

$$\Phi(i, j+1) = \frac{1}{2} [\Phi(i+1, j) + \Phi(i-1, j)]$$

The value of Φ at time step $j+1$ depends only on two values of Φ at the previous time step j .



The stability issue can be solved by adopting the **implicit formula**, proposed by Crank and Nicholson in 1974, which is **unconditionally stable**, for any value of r .

In the implicit formula, the space derivative is replaced by the average of the central difference at time step j e $j+1$:

$$k \frac{\Phi(i, j+1) - \Phi(i, j)}{\Delta t} = \frac{1}{2} \left[\frac{\Phi(i+1, j) - 2\Phi(i, j) + \Phi(i-1, j)}{(\Delta x)^2} + \frac{\Phi(i+1, j+1) - 2\Phi(i, j+1) + \Phi(i-1, j+1)}{(\Delta x)^2} \right]$$

After re-ordering, we obtain:

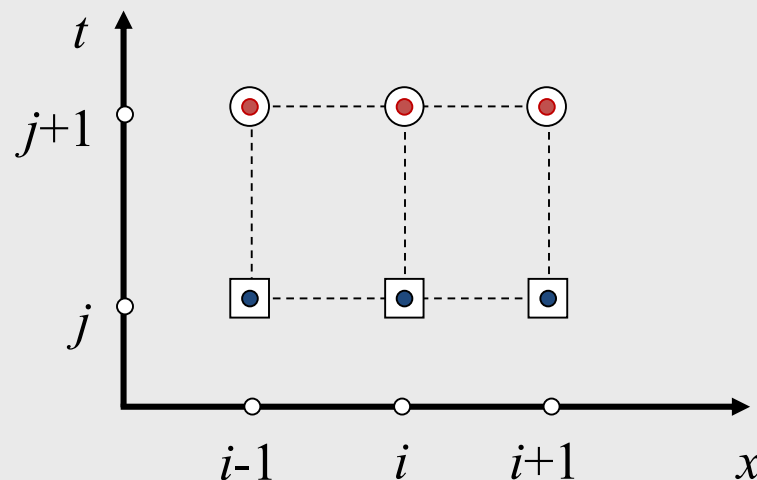
$$\begin{aligned} -r \Phi(i-1, j+1) + 2(1+r) \Phi(i, j+1) - r \Phi(i+1, j+1) = \\ = r \Phi(i-1, j) + 2(1-r) \Phi(i, j) + r \Phi(i+1, j) \end{aligned}$$

Three values of Φ at time step $j+1$ depend on three values of Φ at time step j .

COMPUTATIONAL SCHEME

○ value to determine

□ known values of Φ



The determination of the values of Φ through the implicit formula requires the set the equation in N nodes ($i=1..N$) and to solve a system of N equation with N unknowns.

The implicit formula is unconditionally stable and it allows to **adopt larger time step Δt compared to the explicit formula.**

A convenient choice is typically $r=1$, as it reduces the equation to the following:

$$\begin{aligned} -\Phi(i-1, j+1) + 4\Phi(i, j+1) - \Phi(i+1, j+1) = \\ = \Phi(i-1, j) + \Phi(i+1, j) \end{aligned}$$

The three major sources of error in the numerical solution of physical problems are:

1. **Modeling error** (approximation in the transformation of the real world problem into a model)
2. **Discretization error** (e.g., due to the transformation of the partial derivatives into finite differences)
3. **Rounding error** (due to the finite precision of digital computers)

NOTE: Accuracy increases when the grid size becomes smaller, but there is a limit set by the rounding error.

