

TIF330
Computational Continuum
physics

Problem Set 1

Date: 2023-04-18

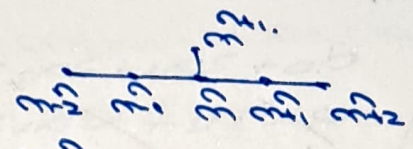
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Problem set 1:

Problem 1: Expand f(z) with 5-point stencil for the heat equation.

Heat equation: $\frac{\partial u}{\partial t} - \alpha \frac{\partial^2 u}{\partial x^2} = 0$

5 point stencil



$$\frac{\partial u}{\partial t} = \frac{u^n_{j-2} - 2u^n_j + u^n_{j+2}}{2\Delta x^2}$$

$$\frac{\partial^2 u}{\partial x^2} = \frac{\alpha_2 u^n_{j-2} + \alpha_1 u^n_{j-1} + \alpha_0 u^n_j + \alpha_1 u^n_{j+1} + \alpha_2 u^n_{j+2}}{\Delta x^2}$$

If we consider symmetry: $\alpha_2 = \alpha_2$ and $\alpha_1 = \alpha_1$

$$u^n_j = \frac{u^n_{j-2} - 2u^n_j + u^n_{j+2}}{2} - \frac{\alpha_2 u^n_{j-2} + \alpha_1 u^n_{j-1} + \alpha_0 u^n_j + \alpha_1 u^n_{j+1} + \alpha_2 u^n_{j+2}}{2}$$

We can perform Taylor expansion:

~~$$u^n_j = u^n_j + \Delta x^2 \frac{\partial^2 u^n_j}{\partial x^2} + \frac{\Delta x^4}{24} \frac{\partial^4 u^n_j}{\partial x^4} + \dots$$~~

~~$$\text{I: } \alpha_2 (u^n_{j-2} - 2u^n_j + u^n_{j+2}) + \frac{\Delta x^2}{2} \frac{\partial^2 u^n_j}{\partial x^2} + \frac{\Delta x^4}{24} \frac{\partial^4 u^n_j}{\partial x^4} + \dots$$~~

~~$$\text{II: } \alpha_1 (u^n_{j-1} - u^n_{j+1}) + \frac{\Delta x^2}{2} \frac{\partial^2 u^n_j}{\partial x^2} + \frac{\Delta x^4}{24} \frac{\partial^4 u^n_j}{\partial x^4} + \dots$$~~

~~$$\text{III: } \alpha_0 (u^n_j - u^n_j) + \frac{\Delta x^2}{2} \frac{\partial^2 u^n_j}{\partial x^2} + \frac{\Delta x^4}{24} \frac{\partial^4 u^n_j}{\partial x^4} + \dots$$~~

~~$$\text{IV: } \alpha_1 (u^n_{j+1} - u^n_{j-1}) + \frac{\Delta x^2}{2} \frac{\partial^2 u^n_j}{\partial x^2} + \frac{\Delta x^4}{24} \frac{\partial^4 u^n_j}{\partial x^4} + \dots$$~~

~~$$\text{V: } \alpha_2 (u^n_{j+2} - 2u^n_j + u^n_{j-2}) + \frac{\Delta x^2}{2} \frac{\partial^2 u^n_j}{\partial x^2} + \frac{\Delta x^4}{24} \frac{\partial^4 u^n_j}{\partial x^4} + \dots$$~~

Summing up:

$$u^n_j = u^n_j + \Delta x^2 \frac{\partial^2 u^n_j}{\partial x^2} + \frac{\Delta x^4}{24} \frac{\partial^4 u^n_j}{\partial x^4} + \dots$$

$$\begin{aligned} \frac{\partial^2 \mathcal{L}}{\partial x^2} = & \frac{\partial}{\partial x} \left(\frac{\partial \mathcal{L}}{\partial x} + \alpha_1 - \frac{\partial}{\partial x} \left(\frac{\partial \mathcal{L}}{\partial x} + 2\alpha_1 + 2\alpha_2 \right) \frac{\partial \mathcal{L}}{\partial x} + \right. \\ & + \left(\frac{\partial^2 \mathcal{L}}{\partial x^2} \alpha_2 + \frac{\partial^2 \mathcal{L}}{\partial x^2} \alpha_1 \right) \frac{\partial^4 \mathcal{L}}{\partial x^4} + \\ & \left. + \left(\frac{\partial^2 \mathcal{L}}{\partial x^2} \frac{\partial^4 \mathcal{L}}{\partial x^4} \alpha_2 + \frac{\partial^2 \mathcal{L}}{\partial x^2} \frac{\partial^4 \mathcal{L}}{\partial x^4} \alpha_1 \right) \frac{\partial^6 \mathcal{L}}{\partial x^6} + \dots \right) \end{aligned}$$

Now, in order to achieve the right order accuracy, we require that:

$$\begin{cases} \beta = 0 \\ \frac{\partial \mathcal{L}}{\partial x} = 0 \\ \delta = 0 \end{cases} \Rightarrow \begin{cases} \alpha_0 + 2\alpha_1 + 2\alpha_2 = 0 \\ \frac{\partial^2 \mathcal{L}}{\partial x^2} \alpha_2 + \frac{\partial^2 \mathcal{L}}{\partial x^2} \alpha_1 = 0 \\ \frac{\partial^2 \mathcal{L}}{\partial x^2} \frac{\partial^4 \mathcal{L}}{\partial x^4} \alpha_2 + \frac{\partial^2 \mathcal{L}}{\partial x^2} \frac{\partial^4 \mathcal{L}}{\partial x^4} \alpha_1 = 0 \end{cases} \Rightarrow \begin{cases} \alpha_1 = 1 - 4\alpha_2 \\ 10\alpha_2 + 1 - 2\alpha_2 = 0 \\ 12\alpha_2 = 2 \end{cases}$$

$$\Rightarrow \begin{cases} \alpha_1 = 1 - 4\alpha_2 \\ \alpha_2 = \frac{1}{12} \end{cases} \Rightarrow \begin{cases} \alpha_1 = \frac{11}{12} \\ \alpha_2 = \frac{1}{12} \end{cases}$$

+ This would imply:

$$\begin{aligned} \mathcal{L} &= \mathcal{L}_0 + \alpha_1 \mathcal{L}_1 - \frac{\partial}{\partial x} \left(\mathcal{L}_0 + 12 \frac{\partial^4 \mathcal{L}}{\partial x^4} + 0 + 0 \right) \\ &= \mathcal{L}_0 - \alpha \frac{\partial^4 \mathcal{L}}{\partial x^4} + 0 \mathcal{L}_1 + \alpha \mathcal{L}_2 \\ &= 0 + 0 \mathcal{L}_1 + \alpha \mathcal{L}_2 = \frac{1}{12} \mathcal{L}_2 \end{aligned}$$

So, the FDR scheme would be:

$$\frac{\partial^2 \mathcal{L}}{\partial x^2} = \frac{-\mathcal{L}_0 + 2\mathcal{L}_1 + 10\mathcal{L}_2 - 30\mathcal{L}_3 + 10\mathcal{L}_4 - \mathcal{L}_5}{12h^2}$$

Suppose we wanted to (try to) go to even higher order (FDR):

for this, our Taylor expansion would have to go to even higher order. Namely the R.H.S. of the equation would, explicitly, be:

$$\begin{aligned} \mathcal{L}_2 &= \alpha_2 \cdot 2 \cdot \frac{\partial^4 \mathcal{L}}{\partial x^4} + \alpha_1 \cdot 2 \cdot \frac{\partial^4 \mathcal{L}}{\partial x^4} + \dots \\ &= \frac{1}{300} \left(84\alpha_2 + \alpha_1 \right) \frac{\partial^4 \mathcal{L}}{\partial x^4} + \dots \end{aligned}$$

going to higher order would require:

$$\begin{aligned} 84\alpha_2 + \alpha_1 &= 0 \Rightarrow \alpha_1 = -84\alpha_2 \\ &\Rightarrow \frac{1}{12} + \frac{84}{12} \alpha_2 = 0, \text{ which is false...} \end{aligned}$$

We have arrived at a contradiction. So we cannot reach higher order approximations than $\mathcal{O}(h^4)$.

Now, we do Stability Analysis (on numerical):

FDM: $\frac{u_m^{n+1} - u_m^n}{\tau} = \frac{\partial}{\partial x^2} (-u_{m-2}^n + 10u_{m-1}^n - 30u_m^n + 10u_{m+1}^n - u_{m+2}^n)$

Assume wave-type solution: $u_m^n = \lambda^n e^{ikh_m} \Rightarrow$

$$\Rightarrow \frac{\lambda - 1}{\tau} = \frac{\partial}{\partial x^2} \left(-\frac{e^{-2ikh}}{2} + 10\frac{e^{-ikh}}{2} - 30 + 10\frac{e^{ikh}}{2} - \frac{e^{2ikh}}{2} \right) \Rightarrow$$

$$\Rightarrow \frac{\lambda - 1}{\tau} = \frac{\partial}{\partial x^2} (-30 + 32 \cos kh - 2 \cos 2kh)$$

Using: $\cos kh = 1 - 2 \sin^2(\frac{kh}{2})$, we get:

$$\lambda = 1 + \frac{\partial \tau}{\partial x^2} (32 - 64 \sin^2(\frac{kh}{2}) - 2 + 4 \sin^2(\frac{kh}{2}) - 30) \Rightarrow$$

$$\Rightarrow \lambda = 1 + \frac{\partial \tau}{\partial x^2} (\sin^2 kh - 10 \sin^2(\frac{kh}{2}))$$

Stability requires that: $|\lambda| \leq 1$:

$$-1 \leq 1 + \frac{\partial \tau}{\partial x^2} (\sin^2 kh - 10 \sin^2(\frac{kh}{2})) \leq 1 \Rightarrow$$

$$\Rightarrow -2 \leq \frac{\partial \tau}{\partial x^2} (\sin^2 kh - 10 \sin^2(\frac{kh}{2})) \leq 0 \Rightarrow$$

$$\Rightarrow -\frac{2 \cdot 3h^2}{\partial^2} \leq \sin^2 kh - 10 \sin^2(\frac{kh}{2}) \leq 0$$

remember that $\frac{\partial^2}{\partial x^2} < 0$.

Worst case scenario: $\frac{kh}{2} = \frac{\pi}{2} \Rightarrow$

$$\Rightarrow -\frac{9h^2}{\partial^2} \leq 0 - 10 \leq 0 \Rightarrow$$

$$\Rightarrow \frac{h^2}{\tau^2} \geq \frac{100}{9} \Rightarrow \frac{h^2}{\tau^2} \geq \frac{80}{9} \Rightarrow \left\{ \frac{\tau}{h^2} \leq \frac{3}{80} \right\} \Rightarrow$$

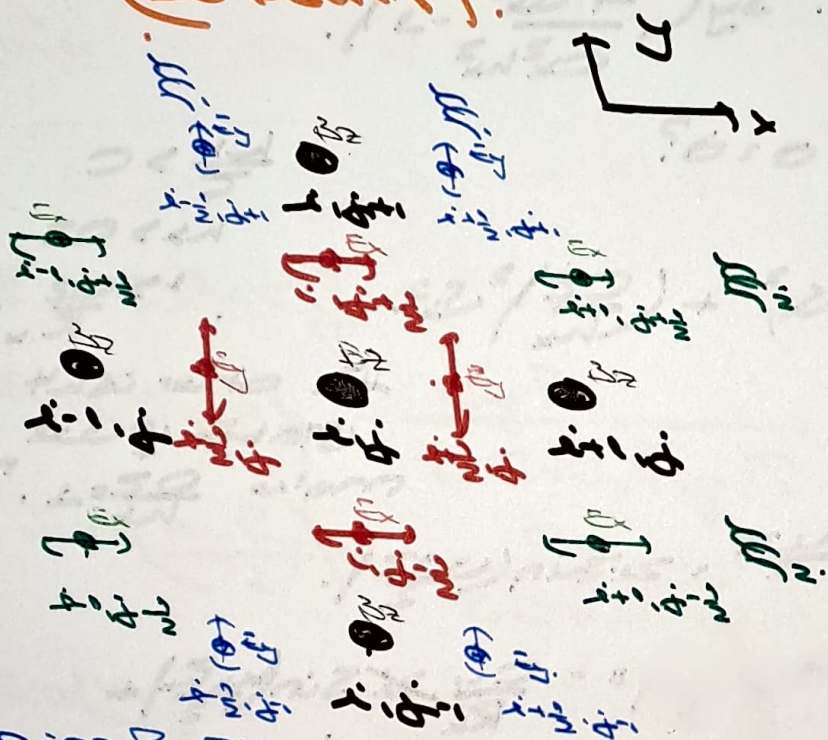
$$\Rightarrow \left\{ \tau \leq \frac{3}{80} \right\}, \tau = \frac{\tau}{h^2} - \text{parabolic Courant number.}$$

Stability condition.

Note: note that $\sin^2 kh - 10 \sin^2(\frac{kh}{2}) \leq 0$ is always true regardless of the value of $k \in \mathbb{R}$ (or \mathbb{Z}).

Problem 2: FDTD modification without numerical dispersion along one of the grid axes (20 points):

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• blue and green
points on the
y-axis are
added to the
original strain.

$\frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}$

$$\Delta = \frac{1}{2} \left(\frac{c_1^2 + c_2^2 - c_3^2}{2} - \frac{c_1^2 + c_2^2 - c_3^2}{2} \right)$$

Introducing the blue and green points, we get:

(Introduce the blue points):

$$\left\{ \begin{aligned} & \frac{C(i+2, j+1) - C(i+1, j+1)}{A_x} + \frac{C(i+1, j+1) - C(i, j+1)}{A_x} + \frac{C(i, j+1) - C(i, j)}{A_x} \\ & + (1-2) \frac{C(i, j+1) - C(i, j)}{A_x} \end{aligned} \right\} \text{Blue points}$$

$$\left\{ \begin{aligned} & \frac{C(i+1, j+2) - C(i+1, j+1)}{A_y} + \frac{C(i+1, j+1) - C(i+1, j)}{A_y} + \frac{C(i+1, j) - C(i, j)}{A_y} \\ & + (1-2) \frac{C(i+1, j) - C(i, j)}{A_y} \end{aligned} \right\} \text{Green points}$$

+ 2D Laplace FD scheme becomes:

$$\frac{C(i+2, j) - 2C(i+1, j) + C(i, j))}{A_x^2} + \frac{C(i, j+2) - 2C(i, j+1) + C(i, j))}{A_y^2} + (1-2) \frac{C(i, j+1) - C(i, j)}{A_y} + (1-2) \frac{C(i+1, j) - C(i, j)}{A_x}$$

(1) (2) } these two equations stay the same as in the original scheme (see previous page).

Now, in order to avoid dispersion, we introduce wave-type solutions:

$$\begin{aligned} & u(x, y, t) = A e^{i(k_x x + k_y y - \omega t)} \\ & \text{where } \omega = \omega(k_x, k_y) \end{aligned}$$

... notation

$$\frac{1}{2} \cdot \frac{20}{20} \cdot \frac{20}{20} = 1$$

$\frac{1}{2} \cdot \frac{20}{20} \cdot \frac{20}{20} = 1$
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 $\frac{1}{2} \cdot \frac{20}{20} \cdot \frac{20}{20} = 1$

$$P = \begin{pmatrix} \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} \end{pmatrix} \begin{pmatrix} \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} \end{pmatrix} = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

$$\Rightarrow S_x^2 = \left(\frac{C^2}{R_x^2}\right)^2 S_x^2 (1 - 4 \beta^2 S_x^2) + \left(\frac{C^2}{R_x^2}\right)^2 S_x^2 (1 - 4 \beta^2 S_x^2)$$

numerical dispersion relation.

Now, ~~for instance~~, we consider a wave traveling along $x=0$ ~~at~~ $y=0 \Rightarrow S_y = \sin(k_y R_y), \sin(0)=0$.

$$\Rightarrow S_x^2 = \left(\frac{C^2}{R_x^2}\right)^2 S_x^2$$

In order to have no dispersion right, we add the condition $\left\{\frac{C^2}{R_x^2} = 1\right\}$ + this condition.

$$S_x^2 = \left(\frac{C^2}{R_x^2}\right)^2 S_x^2 \Rightarrow S_x^2 = S_x^2 \Rightarrow \sin^2\left(\frac{k_x R_x}{2}\right), \sin^2\left(\frac{k_y R_y}{2}\right) = 1$$

$$\Rightarrow \omega^2 = k_x R_x$$

$$\Rightarrow \omega = k_x R_x = \sqrt{\omega^2} = \sqrt{k_x R_x}$$

Let's:

$$\frac{C^2}{R_x^2} = 1 \Rightarrow C = R_x$$

relationship of x & y no dispersion!

Some of the dispersion relation!

Additionally, it's important to over the stability of this method.

+ this [see numerical] analysis

follows: pretty much directly from the numerical dispersion relation.

We have stability if $\omega \leq \omega_{max} \Rightarrow$ max in our solution doesn't explode as $t \rightarrow \infty$. + this condition implies that:

$$S_x^2 = \left(\frac{C^2}{R_x^2}\right)^2 S_x^2 (1 - 4 \beta^2 S_x^2) + \left(\frac{C^2}{R_x^2}\right)^2 S_x^2 (1 - 4 \beta^2 S_x^2) \leq 1$$

$$0 \leq \beta^2 \leq 1$$

We can check if this is verified for the maximal first dispersion: $k_y=0$ and $\left\{\frac{C^2}{R_x^2} = 1\right\} \Rightarrow$

$$0 \leq \beta^2 = \left(\frac{1}{2} S_x^2 (1 - 0) + 0\right) = \frac{1}{2} S_x^2 \leq 1$$

which means that:

method is stable under this condition for any order of time space derivatives. + this is true for any order of time space derivatives.

Problem 3

In this problem, we are asked to develop a C++ implementation of the 1D FDTD method for numerical solution of Maxwell's equations in vacuum:

$$\begin{cases} \frac{\partial E_y}{\partial t} = -\frac{\partial B_z}{\partial x} \\ \frac{\partial B_z}{\partial t} = -\frac{\partial E_y}{\partial x} \end{cases} \quad (1)$$

Moreover, we consider the propagation of a linearly polarized electromagnetic pulse, for $x \in [-1, 1]$, towards the positive x direction, and for a time $t \in [0, 2.5]$, starting from the following initial conditions:

$$E_y(x) = \begin{cases} \sin(20\pi x), & |x| < 0.1 \\ 0 & |x| \geq 0.1 \end{cases} \quad (2)$$

Since the value of $B_z(x, t = 0)$ is not specified, it is assumed to be zero everywhere.

Periodic boundary conditions

In this case, we are asked to plot $E_y(x)$ and $B_z(x)$ at $t = 2.5$. My code will generate a gif with an animation of the propagation of the electromagnetic fields (as well as many images, used to build that gif). The requested plot will be the last image. For convinence, I reproduce it here.

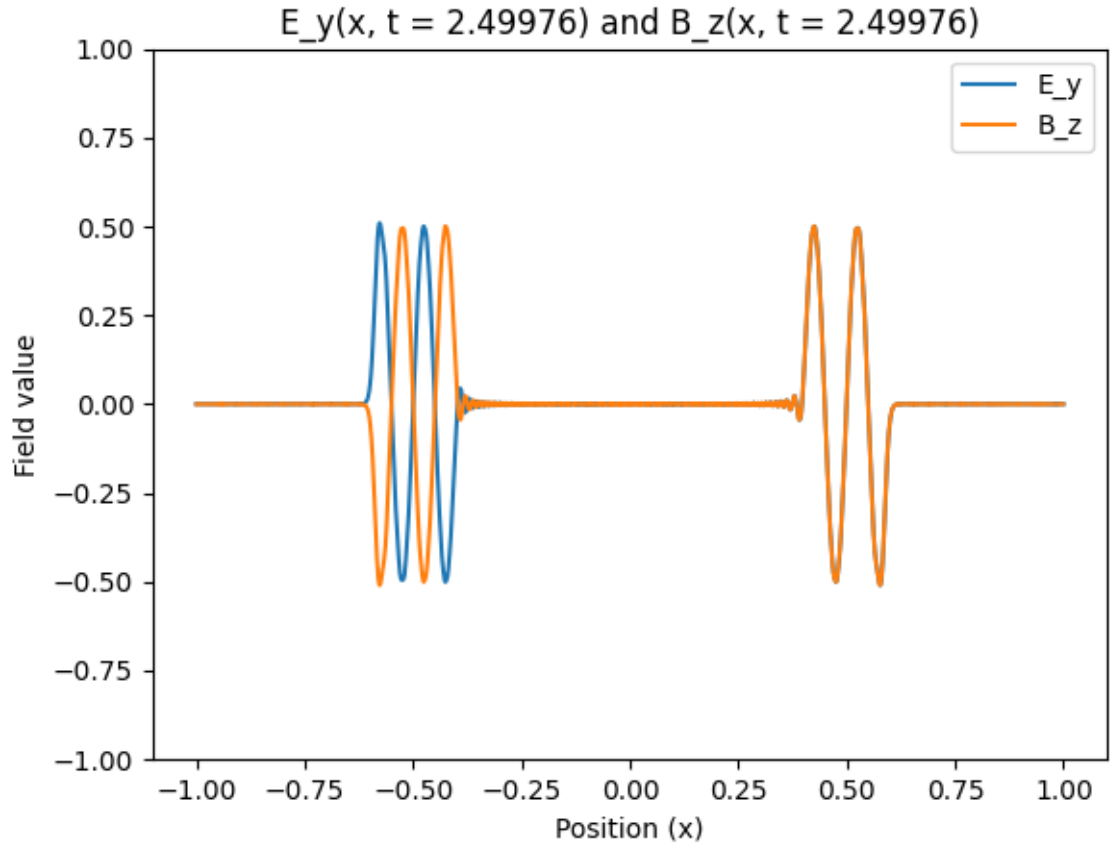


Figure 1: Electromagnetic fields at $t = 2.5$, in the periodic boundary condition case. The onset of numerical dispersion is clearly visible!

In order to produce the aforementioned `gifs` and images, first compile and run the `C++` code, followed by the `Python` script. The first will generate a bunch of `.txt` files with the values of the E and B fields (for all the many time-steps), resultant from the FDTD calculations, for each of the three boundary condition cases. The later will generate the corresponding images and `gifs`, to help in the visualization of the results. Import aspects to mention about the developed code:

1. My space grid always starts at an E field point and, likewise, always ends at an E field point. This is just a choice that I made. As a consequence, we have one less B field point than E field points. By doing things this way, I also only need to implement boundary conditions for the (extreme) E field points (**not** for the B field), since there is no B field point at the grid boundary;
2. The time-step is chosen to be equal to the distance between adjacent E and B field points. This will be important, later, for the implementation of the perfect boundary condition. Also, note that (Courant) stability is assured so long as: $2h/\tau > c_0$ ¹ (I assume propagation in vacuum. Otherwise, we'd need to include n [refractive index] here as well). This condition, in our system of units, is written as $2h/\tau > 1$, since c_0 is taken to be 1. Thus, we have chosen $2h/\tau = 2$ (Stable!), which means that a certain value of the $E(B)$ field will take two time-steps (2τ) to travel to the next position (index) in its corresponding vector, i.e., to go from `ez[0]` to `ez[1]`², if we take Fig. 2 as reference.

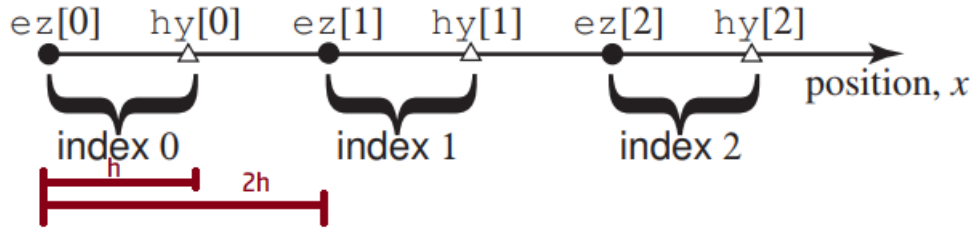


Figure 2: Considered FDM grid and relevant distances. Taken from Ref. [1]. Slight modifications to include distances h and $2h$.

3. Finally, my code computes all the three boundary condition cases all at once. So, it needs only to be run once, followed by the `Python` script (also once). **Notes:** run the `C++` and `Python` codes from the same directory, and keep in mind that the `Python` code requires `numpy`, `matplotlib` (plots) and `imageio` (for making the gif).

At each iteration, all the "middle" points are taken care of using "regular" 1D FDTD update equations. In my code, these would be:

$$\begin{aligned} E_y[mm] &= E_y[mm] - (\tau/(2*h))*(B_z[mm] - B_z[mm - 1]); \\ B_z[mm] &= B_z[mm] - (\tau/(2*h))*(E_y[mm + 1] - E_y[mm]); \end{aligned} \quad (3)$$

¹The reason why I write $2h$ instead of just h is because of the way I defined the grid in my code. There, the distance between adjacent $E(B)$ points is $2h$, whereas the distance between adjacent E and B points is just h (Alternating E and B points). Check Fig. 2.

²Also, note that Fig. 2 does not describe my code. It was taken from Ref. [1], where they have a different convention. In my case, I compute E_y and B_z , not E_z and B_y . This figure is merely used to illustrate the distances involved (in the $(2)h$ discussion from above).

The only place where we can have "problems" is at the boundaries, where we would require the value of the magnetic field (and in my case only the magnetic field, since the grid's extreme points are E at both ends) outside of the grid. In the case of periodic boundary conditions, we simply allow our grid to loop back around. This means that, for the last E point, we utilize the B point before it, as well as the first B point of the grid. The same goes for the first E point of the grid. In my code, this reads:

$$\begin{aligned} E_y[0] &= E_y[0] - (\tau/h)*(B_z[0] - B_z[SIZE - 2]); \\ E_y[SIZE - 1] &= E_y[SIZE - 1] - (\tau/h)*(B_z[0] - B_z[SIZE - 2]); \end{aligned} \quad (4)$$

The results of implementing this type of boundary condition are very easily understood if we look at the produce gif, where we see the waves propagating towards the edge of the simulation grid, only to then reappear at the other end of the grid. (**Note:** All the [full] developed codes are presented in the appendix, below.)

Perfect magnetic conductor boundary conditions

In order to simulate perfect magnetic conductor (PMC) boundaries, we merely have to set the magnetic field outside the grid to zero. We no longer allow it to loop back around. On external boundaries, a perfect magnetic conductor can be interpreted as a "high surface impedance"³, which doesn't allow current to flow into it (surface current density is zero along this boundary surface). Due to the way I defined my grid (E points on both extremes), we have a PMC on both ends. Had I defined it with a B point in either of its edges, setting the E field outside the grid to zero would be equivalent to having a perfect electric conductor-type boundary, instead. Either way, given how I implemented my grid, the B field will be inverted when it reaches the end of the grid, just like waves on a rope, reflecting from a fixed end. The electric field, on the other hand, will reflect symmetrically, not inverting its sign. This can be very readily observed in the gif that my program produces. Explicitly, this type of boundary condition is implemented as such:

$$\begin{aligned} E_y[0] &= E_y[0] - (\tau/(2*h))*(B_z[0] - 0); \\ E_y[SIZE - 1] &= E_y[SIZE - 1] - (\tau/(2*h))*(0 - B_z[SIZE - 2]); \end{aligned} \quad (5)$$

Perfect boundary conditions

Finally, in order to implement what I call "perfect boundary conditions", we come back to the definition of τ as $\tau = h$. As mentioned previously, this makes it so that after **two** time-steps, the E value that we had at position $e_y[n]$ will now be at position $e_y[n + 1]$ (Same thing for B). In my code, the perfect boundary conditions are implemented as follows:

$$\begin{aligned} E_y[0] &= E_y[0] - (\tau/(2*h))*(B_z[0] - Res_B[(qTime + 1) - 2][0]); \\ E_y[SIZE - 1] &= E_y[SIZE - 1] - \\ &(\tau/(2*h))*(Res_B[(qTime + 1) - 2][SIZE - 2] - B_z[SIZE - 2]); \end{aligned} \quad (6)$$

The reason why this works is the following. Suppose we are at an extreme E point. In order to compute its next value, we require the value of the B field **one index** outside the grid, as if there was no boundary and the grid just continued infinitely. Now, since

³More detailed definition in Ref. [2].

we know that after two time-steps, our fields will have advanced by one index, we can recognize that the value we are looking for is merely the value of the B field two time-steps ago (which we stored previously as part of our solution!). The description of this perfect boundary condition is described in more detail in Ref. [3], where a more involved discussion of the topic is presented.

With all this being said, I can now easily simulate a part of infinite space, as requested. More specifically, my code demonstrates this by simulating the continuous emission of an antenna placed inside the simulation region. The produced gif makes this type of boundary condition very easy to visualize, as it appears the waves simply wander off to infinity.

References

- [1] Chapter 3: Introduction to the Finite-Difference Time-Domain Method: FDTD in 1D, Lecture notes by John Schneider. Available at: <https://eecs.wsu.edu/~schneidj/ufttd/chap3.pdf>;
- [2] Perfect Magnetic Conductor. Available at: https://doc.comsol.com/5.5/doc/com.comsol.help.rf/rf_ug_radio_frequency.07.12.html;
- [3] Lecture 6 (FDTD) - Implementation of 1D FDTD. Available at: <https://www.youtube.com/watch?v=hNN7EtZsJuU>;
- [4] Chapter 8: Two-Dimensional FDTD Simulations, Lecture notes by John Schneider. Available at: <https://eecs.wsu.edu/~schneidj/ufttd/chap8.pdf>;
- [5] Perfect Electric Conductor. Available at: https://doc.comsol.com/5.5/doc/com.comsol.help.rf/rf_ug_radio_frequency.07.11.html;

Developed code

C++ code: P3_Final.cpp

```
#include <iostream>
#include <cmath>
#include <vector>
#include <fstream>
#include <string>

#define SIZE 2048

using namespace std;

/* Auxiliary functions */
void Write_to_file(vector<vector<double>> &Res, string filename)
{
    cout << "Entering \"Write_to_file\"... " << endl;
    cout << "Matrix dimensions: " << size(Res) << " " << size(Res[0])
    << endl;
    fstream w_File;
    w_File.open(filename, fstream::out);
    for (int i = 0; i < size(Res); i++)
    {
        for (int j = 0; j < size(Res[0]); j++)
        {
            if(j < (size(Res[0]) - 1)) w_File << Res[i][j] << ",";
            else w_File << Res[i][j] << endl;
        }
    }
    w_File.close();
}

void Write_Vec_to_file(vector<double> &Vec, string filename)
{
    cout << "Entering \"Write_Vec_to_file\"... " << endl;
    fstream w_File;
    w_File.open(filename, fstream::out);
    for (int i = 0; i < size(Vec); i++)
    {
        if(i < (size(Vec) - 1)) w_File << Vec[i] << ",";
        else w_File << Vec[i] << endl;
    }
}

void Add_to_Res(double* E_y, double* B_z, vector<vector<double>> &Res_E
, vector<vector<double>> &Res_B)
{
    vector<double> E_y_, B_z_;
    E_y_.assign(E_y, E_y + SIZE);
```



```

    B_z_.assign(B_z, B_z + (SIZE - 1));
    Res_E.push_back(E_y_);
    Res_B.push_back(B_z_);
    E_y_.clear();
    B_z_.clear();
}

/* Main program */
int main(int argc, const char** argv)
{
    /* Boundary condition type (different cases):
    1 - Periodic;
    2 - Perfect Magnetic Couductor;
    3 - Perfect (Infinity). */
    for (int BC_type = 1; BC_type < 4; BC_type++)
    {
        /* Variables */
        vector<double> t_vec;
        vector<vector<double>> Res_E, Res_B;
        int qTime, mm;
        double E_y[SIZE] = {0.}, B_z[SIZE - 1] = {0.};
        double h = 2/(2 * double(SIZE) - 2);
        double tau = h;
        long int maxTime = 2.5/tau;
        double x[2 * SIZE - 1] = {0.};
        bool delay = false;
        vector<string> names = {"Periodic_BC", "PMC_BC", "
Perfect_BC"};

        /* These ended up not being needed */
        int numberOfImages = 500; /* We only save 500 images */
        int stride = int(double(maxTime)/double(
numberOfImages));

        /* Printing for debugging/reference */
        printf("Parameters: \n");
        printf("h = %f\n", h);
        printf("gridSize = %d\n", SIZE);
        printf("tau = %f\n", tau);
        printf("2h/tau = %f > c = 1 (Stable)\n", (2*h)/tau);
        ;
        printf("numberOfIterations = %ld\n", maxTime);
        printf("numberOfImages = %d\n", numberOfImages);
        printf("stride = %d\n", stride);

        /* Filling up space grid vector */
        for(mm = 0; mm < (2 * SIZE - 1); mm++) x[mm] = -1 + (mm * h);

        if(BC_type < 3)

```

```

{
    /* Initial condition */
    for(mm = 0; mm < SIZE; mm++)
    {
        if(abs(x[2 * mm]) < 0.1) E_y[mm] = sin(20 * M_PI * x[2 * mm
]);
        else E_y[mm] = 0;
    }
}

/* 'Pushing back' intial condition */
t_vec.push_back(0.);
Add_to_Res(E_y, B_z, Res_E, Res_B);

/* Do time stepping */
for (qTime = 0; qTime < maxTime; qTime++)
{

    /* Update magnetic field. Was missing factors of (tau/h) here (
and in the electric field equation(s))! */
    for (mm = 0; mm < (SIZE - 1); mm++) B_z[mm] = B_z[mm] - (tau
/(2*h))*(E_y[mm + 1] - E_y[mm]);

    /* Update electric field */
    switch (BC_type)
    {
        case 1:
            /* Boundary conditions - Periodic! First and last points,
respectively */
            E_y[0] = E_y[0] - (tau/h)*(B_z[0] - B_z[SIZE
- 2]);
            E_y[SIZE - 1] = E_y[SIZE - 1] - (tau/h)*(B_z[0] - B_z[SIZE
- 2]);
            break;
        case 2:
            /* Perfect magnetic conductor at edges */
            E_y[0] = E_y[0] - (tau/(2*h))*(B_z[0] - 0);
            E_y[SIZE - 1] = E_y[SIZE - 1] - (tau/(2*h))*(0 - B_z[SIZE -
2]);
            break;
        case 3:
            /* Perfect boundary conditions */
            if(qTime > 0)
            {
                E_y[0] = E_y[0] - (tau/(2*h))*(B_z[0] -
Res_B[(qTime + 1) - 2][0]);
                E_y[SIZE - 1] = E_y[SIZE - 1] - (tau/(2*h))*(Res_B[(
qTime + 1) - 2][SIZE - 2] - B_z[SIZE - 2]);
            }
    }
}

```



```

        break;
    }

    /* All the other 'middle' points */
    for (mm = 1; mm < SIZE - 1; mm++) E_y[mm] = E_y[mm] - (tau/(2*h
    ))*(B_z[mm] - B_z[mm - 1]);

    /* Delay counter: Antenna beeps and then waits for 500
    iterations before beeping again */
    if(BC_type == 3)
    {
        if ((qTime+1) % 500 == 0) delay = !delay;
        if (!delay) E_y[SIZE/2] += 0.75 * abs(sin(
M_PI * (double(qTime)/500.)));
    }

    /* Results */
    if (qTime % 1 == 0) // 'stride' was changed to 1 for now.
    {
        cout << "Progress: " << 100*(qTime/double(maxTime)) << " %"
<< endl;
        t_vec.push_back((qTime + 1)*tau);
        Add_to_Res(E_y, B_z, Res_E, Res_B);
    }

} /* End of time-stepping */

/* Writing to text file */
switch (BC_type)
{
case 1:
    Write_to_file(Res_E, "E_Periodic_BC.txt");
    Write_to_file(Res_B, "B_Periodic_BC.txt");
    break;
case 2:
    Write_to_file(Res_E, "E_PMC_BC.txt");
    Write_to_file(Res_B, "B_PMC_BC.txt");
    break;
case 3:
    Write_to_file(Res_E, "E_Perfect_BC.txt");
    Write_to_file(Res_B, "B_Perfect_BC.txt");
    break;
}

/* Additionally saving the time vector, for easier plotting later
*/
// ofstream outFile("timeVec.txt");
// for (const auto &e : t_vec) outFile << e << "\n"; // Find out
why this doesn't work!

```

```

        Write_Vec_to_file(t_vec, "timeVec.txt");
    }

    return(0);
}

```

Listing 1: C++ code utilized in Problem 3 to resolve the FDTD scheme. For ease of copying: <https://codeshare.io/6p1Wbm>.

Python script: P3_Gif.py

```

import matplotlib.pyplot as plt
import numpy as np
import imageio.v2 as imageio
from pathlib import Path

# Definitions
h = 2/(2 * float(2048) - 2); tau = 4/(2048**2)
x_E = [-1 + (j * 2*h) for j in range(2048)]
x_B = [(-1 + h) + (j * 2*h) for j in range(2047)]
Data = {}
ids = ['E_y_1', 'B_z_1',
        'E_y_2', 'B_z_2',
        'E_y_3', 'B_z_3', 'TimeVec']
cases = ['./Periodic_BC', './PMC_BC', './Perfect_BC']
filenames = ['./E_Periodic_BC.txt', './B_Periodic_BC.txt',
              './E_PMC_BC.txt', './B_PMC_BC.txt',
              './E_Perfect_BC.txt', './B_Perfect_BC.txt', 'timeVec.txt']

]

for folder in cases:
    Path(f"{folder}").mkdir(parents=True, exist_ok=True)

for file, id in zip(filenames, ids):
    with open(file, 'r') as f:
        Data[id] = [[float(num) for num in line.split(',') for line in f]]

for ind in range(1,4):
    for i in range(0, len(Data[f'E_y_{ind}'])):
        if((i % 40 == 0) or (i == (len(Data[f'E_y_{ind}']) - 1))):
            plt.plot(x_E, Data[f'E_y_{ind}'][i], label = 'E_y')
            plt.plot(x_B, Data[f'B_z_{ind}'][i], label = 'B_z')
            plt.title(f"E_y(x, t = {Data['TimeVec'][0][i]:.5f}) and B_z(x, t = {Data['TimeVec'][0][i]:.5f})")
            plt.ylabel(f"Field value")
            plt.xlabel("Position (x)")
            plt.ylim(-1, 1)
            plt.legend()

```



```

plt.savefig(f"./{cases[(ind - 1)]}/Img{int(i/40)}.png", dpi
= 100)
plt.clf()

# Small piece of code to join all the 127 images into a gif, for
easier visualization. Takes some minutes to run.
filenames = [f'./{cases[(ind - 1)]}/Img{n}.png' for n in range(0,
int(i/40) + 1)]
with imageio.get_writer(f'./{cases[(ind - 1)]}/Movie_E+B.gif', mode
='I') as writer:
    for filename in filenames:
        image = imageio.imread(filename)
        writer.append_data(image)

```

Listing 2: Python code utilized in Problem 3 to generate the plots and gif. For ease of copying: <https://codeshare.io/OdyxwN>.