

# Pushing charged particles around

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The goal is to simulate the motion of charged particles in a given electric field  $\mathbf{E}(\mathbf{x}, t)$ , and magnetic field  $\mathbf{B}(\mathbf{x}, t)$ . This is the test-particle approximation.

For each particle  $p$  of charge<sup>1</sup>  $q_p$  and mass  $m_p$  the equations to be solved (i.e. integrated in time) are

$$\frac{d\mathbf{x}_p}{dt} = \mathbf{v}_p \quad (1)$$

$$m \frac{d\mathbf{v}_p}{dt} = q_p [\mathbf{E}(\mathbf{x}_p, t) + \mathbf{v}_p \times \mathbf{B}(\mathbf{x}_p, t)] \quad (2)$$

where  $\mathbf{v}_p(t)$  is the velocity and  $\mathbf{x}_p(t)$  is its position. For simplicity we shall drop the subscript  $p$ . Possibly, the simplest discretization is that of Euler:

$$\frac{\mathbf{x}_{n+1} - \mathbf{x}_n}{\Delta t} = \mathbf{v}_n \quad (3)$$

$$\frac{\mathbf{v}_{n+1} - \mathbf{v}_n}{\Delta t} = \frac{q}{m} [\mathbf{E}_n(\mathbf{x}_n) + \mathbf{v}_n \times \mathbf{B}_n(\mathbf{x}_n)] \quad (4)$$

where the electric and magnetic fields are evaluated at the particle position  $\mathbf{x}_{n+\frac{1}{2}}$ . This scheme is however only first order in  $\Delta t$  and not very useful, but for better understanding the next method.

A much better numerical scheme is achieved by staggering in time the position and velocity (leapfrog), so that

$$\frac{\mathbf{x}_{n+\frac{1}{2}} - \mathbf{x}_{n-\frac{1}{2}}}{\Delta t} = \mathbf{v}_n \quad (5)$$

$$\frac{\mathbf{v}_{n+1} - \mathbf{v}_n}{\Delta t} = \frac{q}{m} \left[ \mathbf{E}_{n+\frac{1}{2}} + \frac{\mathbf{v}_n + \mathbf{v}_{n+1}}{2} \times \mathbf{B}_{n+\frac{1}{2}}(\mathbf{x}_{n+\frac{1}{2}}) \right] \quad (6)$$

The electric and magnetic fields are evaluated at the particle position  $\mathbf{x}_{n+\frac{1}{2}}$ . In these equations, time has been discretized as  $t = n\Delta t$  where the subscript  $n$  is an integer. The solution is then found at times  $(1, 2, 3, \dots)\Delta t$  for the velocity and at times  $(-\frac{1}{2}, \frac{1}{2}, \frac{3}{2}, \dots)\Delta t$  for the position.

One of the most widely used, for its accuracy and simplicity, is the so-called Boris pusher<sup>2</sup>. The method is based on defining two new variables for the velocity

$$\mathbf{v}^- = \mathbf{v}_n + \frac{q}{m} \mathbf{E}_{n+\frac{1}{2}} \frac{\Delta t}{2} \quad (7)$$

$$\mathbf{v}^+ = \mathbf{v}_{n+1} - \frac{q}{m} \mathbf{E}_{n+\frac{1}{2}} \frac{\Delta t}{2} \quad (8)$$

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<sup>1</sup> $q = -e$  for electrons and  $Ze$  for ions, where  $Z$  is the ion charge

<sup>2</sup>J. Boris, in Proceedings of the Fourth Conference on Numerical Simulation of Plasmas (Naval Research Laboratory, Washington DC, 1970), p. 3.

which when substituted into eq. 6 gives

$$\frac{\mathbf{v}^+ - \mathbf{v}^-}{\Delta t} = \frac{q}{m} \frac{\mathbf{v}^- + \mathbf{v}^+}{2} \times \mathbf{B}_{n+\frac{1}{2}} \quad (9)$$

The Boris algorithm results in the “separation” of the electric field and magnetic field forces acting on the particle.

The steps to update the velocity from  $\mathbf{v}_n$  to  $\mathbf{v}_{n+1}$  are: given by:

1. Add half of the electric impulse to  $\mathbf{v}_n$  using eq.7 to obtain  $\mathbf{v}^-$
2. Perform a rotation using eq. 9 to obtain  $\mathbf{v}^+$ . This is done by decomposing the rotation as

$$\mathbf{v}' = \mathbf{v}^- + \mathbf{v}^- \times \mathbf{t} \quad (10)$$

$$\mathbf{v}^+ = \mathbf{v}^- + \mathbf{v}' \times \mathbf{s} \quad (11)$$

where

$$\mathbf{t} = \frac{q\Delta t}{2m} \mathbf{B}_{n+\frac{1}{2}} \quad (12)$$

$$\mathbf{s} = \frac{2}{1+t^2} \mathbf{t} \quad (13)$$

3. Add the remaining electric impulse using eq. 8 to obtain the updated velocity  $\mathbf{v}_{n+1}$

$$\mathbf{v}_{n+1} = \mathbf{v}^+ + \frac{q}{m} \mathbf{E}_{n+\frac{1}{2}} \frac{\Delta t}{2} \quad (14)$$

## Today's work

Write a code to solve the motion of a charged particle using the Boris particle pusher described above. Particles can be ions ( $q = Ze$ ), electrons ( $q = -e$ ) or positrons ( $q = e$ ). For numerical stability and accuracy use  $\omega_0 \Delta t = 0.1$  where  $\omega_o = \frac{|q|B_0}{m}$ . **Think carefully about initial conditions and the first time-step. Ideally the code should be able to deal with N particles.**

Like all numerical models, we **must** test the code against known analytical solutions. The tests **must** be quantitative. Examples of tests that you should consider are: simple gyro-motion:  $\mathbf{B} = (0, 0, B_0)$  and  $\mathbf{E} = (0, 0, 0)$ ;  $\mathbf{E} \times \mathbf{B}$  drift ( $\mathbf{E}$  and  $\mathbf{B}$  are constant in time and do not vary in space); energy conservation when  $\mathbf{E} = 0$ ; when possible, you should compare to the known analytical solutions.