Boris Push with Spatial Stepping

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- Boris particle push
- Spatial Boris Push Equations
- advantages of Boris push
- Stability and Accuracy
- Comparison with Runge-Kutta



Boris Push



The Boris push is commonly used in plasma physics simulations because of its speed and stability.

- It uses a leapfrog scheme, where the particle is moved, then half of the energy changed is applied.
- After this, the momentum is rotated by the B field, and the rest of the energy change applied.
- The rotation is automatically energy conserving, and the algorithm is symmetric to time reversal, which improves the performance.Second order accurate, but requires only one field evaluation per time step.
- The Boris push is rarely used for accelerator simulations, because it is more convenient to propagate particles in **z**.



Spatial Boris Push



In collab. with P. Stoltz, J. Cary, at Tech-X (& U. Colorado, Boulder)

The spatial Boris scheme exchanges U for \boldsymbol{p}_z and t for z

First, we replace the equation for pz with the equation for U:

$$\frac{dp_x}{dt} = q(E_x + v_y B_z - v_z B_y)$$
$$\frac{dp_y}{dt} = q(E_y + v_z B_x - v_x B_z)$$
$$\frac{d(U/c)}{dt} = q(E_x v_x + E_y v_y + E_z v_z)$$

Replacing t with z, the governing equations of the spatial Boris scheme are:

$$\begin{aligned} \frac{dp_x}{dz} &= \frac{1}{v_z} \frac{dp_x}{dt} = q \left(\frac{E_x}{v_z} + \frac{v_y B_z}{v_z} - B_y \right) \\ \frac{dp_y}{dz} &= \frac{1}{v_z} \frac{dp_y}{dt} = q \left(\frac{E_y}{v_z} - \frac{v_x B_z}{v_z} + B_y \right) \\ \frac{dU/c}{dz} &= \frac{1}{v_z} \frac{dU/c}{dt} = q \left(\frac{v_x E_x}{v_z c} + \frac{v_y E_y}{v_z c} + \frac{E_z}{c} \right) \end{aligned}$$



Decomposition (1)



For spatial Boris push, the equations separate into terms that directly change p_z and terms that don't

 In the temporal Boris scheme, the separation is into one piece that changes U (E-fields) and one that doesn't (B-fields)

The terms that directly change p_z are E_z , B_x , and B_y

$$\frac{d}{dz} \begin{pmatrix} p_x \\ p_y \\ U/c \end{pmatrix} = \frac{q}{p_z} \begin{pmatrix} 0 & B_z & E_x/c \\ -B_z & 0 & E_y/c \\ E_x/c & E_y/c & 0 \end{pmatrix} \begin{pmatrix} p_x \\ p_y \\ U/c \end{pmatrix} + q \begin{pmatrix} -B_y \\ B_x \\ E_z/c \end{pmatrix}$$

For simplicity, rewrite:

$$\frac{dw}{dz} = Mw + b$$



Decomposition (2)



The Boris scheme integrates vector and matrix terms separately

The Boris scheme says first push the vector term one-half step:

$$w^- = w^n + \frac{\Delta z}{2}b$$

This step is implicit and requires some

Then push the matrix term a full step:

more massaging

$$w^{+} - w^{-} = M\left(\frac{w^{+} + w^{-}}{2}\right)\Delta z$$

Finally, push the vector term the final half step:

$$w^{n+1} = w^+ + \frac{\Delta z}{2}b$$

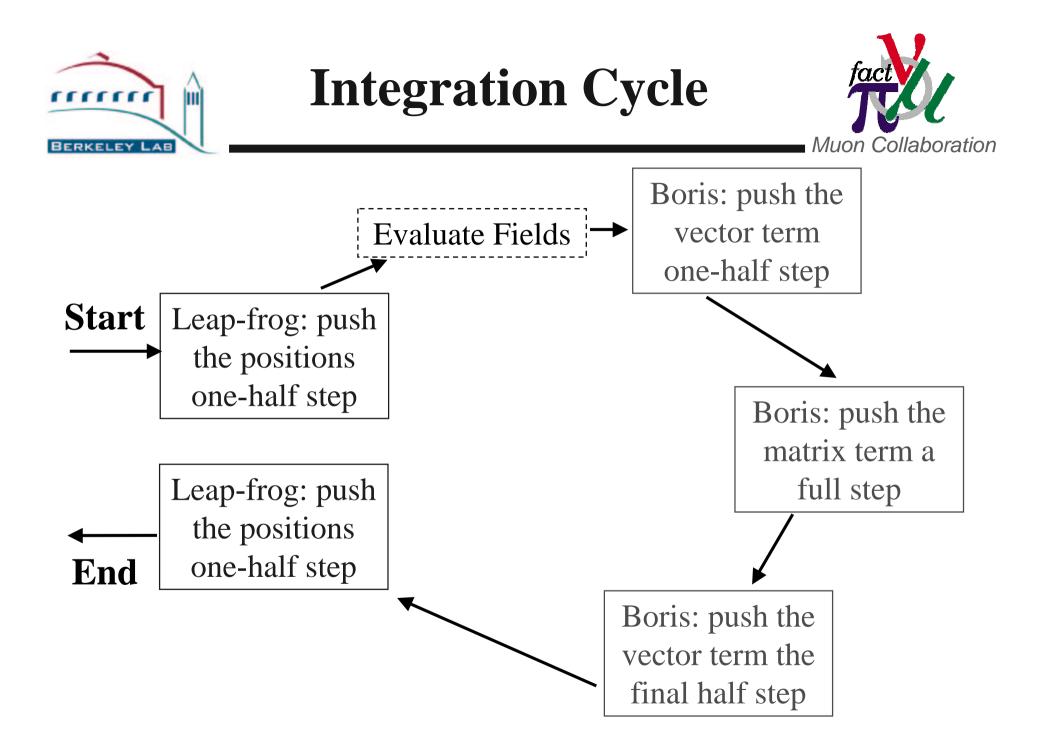


Explicit Expression for Boris Push



Step-centered push of matrix term is 2nd-order accurate

Because M is constant, $w^+ - w^- = M\left(\frac{w^+ + w^-}{2}\right)\Delta z$ a step-centered scheme will be 2nd-order accurate Solving for w^+ gives: $w^+ = \left(I - M\frac{\Delta z}{2}\right)^{-1} \left(I + M\frac{\Delta z}{2}\right) w^ = (I + R) w^{-},$ $R = \frac{q\Delta z/p_z}{1 + \frac{q^2\Delta z^2}{4p_z^2} \left(B_z^2 - \frac{E_x^2 + E_y^2}{c^2}\right)} \begin{pmatrix} \frac{q\Delta z}{2p_z} \left(\frac{E_x^2}{c^2} - B_z^2\right) & B_z + \frac{q\Delta z}{2p_z} \frac{E_x E_y}{c^2} & \frac{E_x}{c} + \frac{q\Delta z}{2p_z} \frac{B_z E_y}{c} \\ -B_z + \frac{q\Delta z}{2p_z} \frac{E_x E_y}{c^2} & \frac{q\Delta z}{2p_z} \left(\frac{E_y^2}{c^2} - B_z^2\right) & \frac{E_y}{c} - \frac{q\Delta z}{2p_z} \frac{B_z E_x}{c} \\ \frac{E_x}{c} + \frac{q\Delta z}{2p_z} \frac{B_z E_y}{c} & \frac{E_y}{c} + \frac{q\Delta z}{2p_z} \frac{B_z E_x}{c} & \frac{q\Delta z}{2p_z} \left(\frac{E_x^2}{c^2} + \frac{E_y^2}{c^2} + \frac{E_y^2}{c^2}\right) \end{pmatrix}$





Boris Push speeds up particle tracking



Spatial motion is calculated a 1/2 step off from momentum/energy evolution:

- typically, use leap-frog
- in ICOOL, split into two half-steps, before and after evolution of *w*.

Fields are only evaluated once, where momentum kick is applied

- compared to 4 field evaluations for RK
- local effect on particle, so almost indep. of coordinates
- track as if no field, then replace Δz with separation between planes (now a function of transverse co-ords)

As in RK, assumes small energy loss per step.



Boris Push is Space-Symmetric



- except ionization energy loss, scatter, are applied at end of step

Second-order conservation of energy

- also canonical momentum when applicable
- robust for large stepsizes

Errors tend to average out

• in RK scheme, errors will slowly accumulate

Both schemes work well for small phase advances, but Boris push is simpler to calculate

- especially if field calculations are expensive
- less savings for curvilinear (where even field-free is complicated)



Error Scaling



Comparison of Runge-Kutta with Boris, for conserved quantities:

Runge Kutta is approx fifth order accurate (special case, solenoid?) tends to yield slowly increasing errors.

Boris push only second order, but preserves invariants

maximum error, Θ , reached in 1 betatron oscillation (or cell period) Relevant length scale is L, step size is Δ

RK: $e \sim R_0 (\Delta_R / L)^5 z / L$ note exponent = 5, not 4

Boris: max $e \sim B_0 (\Delta_B / L)^2$

with 4 x more field calculations per step for RK,

Boris is faster when acceptable to have $P = B_0^{5/3} R_0^{-2/3} (L / z)^{2/3}$



Examples



10 T Uniform Solenoid – conserves P_{\perp}

40 cm for π phase advance: after 400 m, crossover point is e ~ 2 x 10⁻⁵

well below fluctuations for an ensemble of ~ 10^6 particles

FOFO lattice:

1 m half-period with B=2 T, P=200 MeV/c

well above cutoff momentum

after 400 meters, crossover point is roughly the same, $e \sim 3 \ge 10^{-5}$

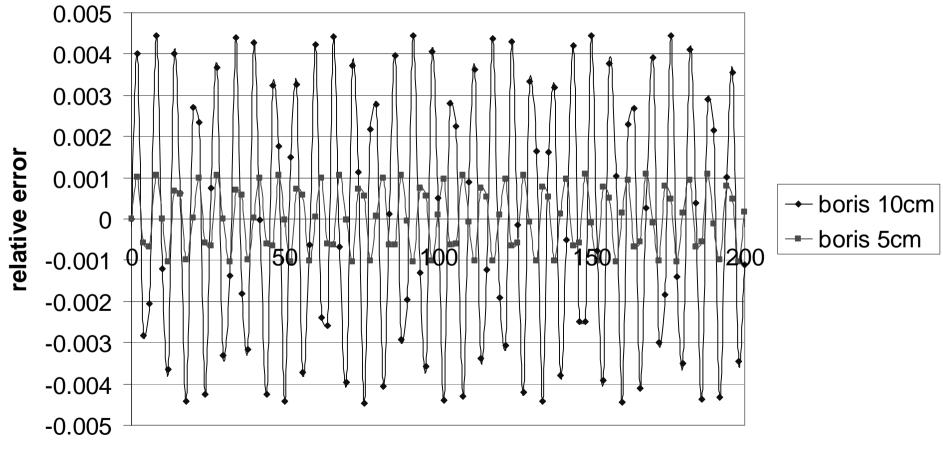
For minimum step sizes fixed by other concerns (e.g., scattering), Boris step is 4 x faster and may still be more accurate



Boris push errors



FOFO: relative error in angular momentum

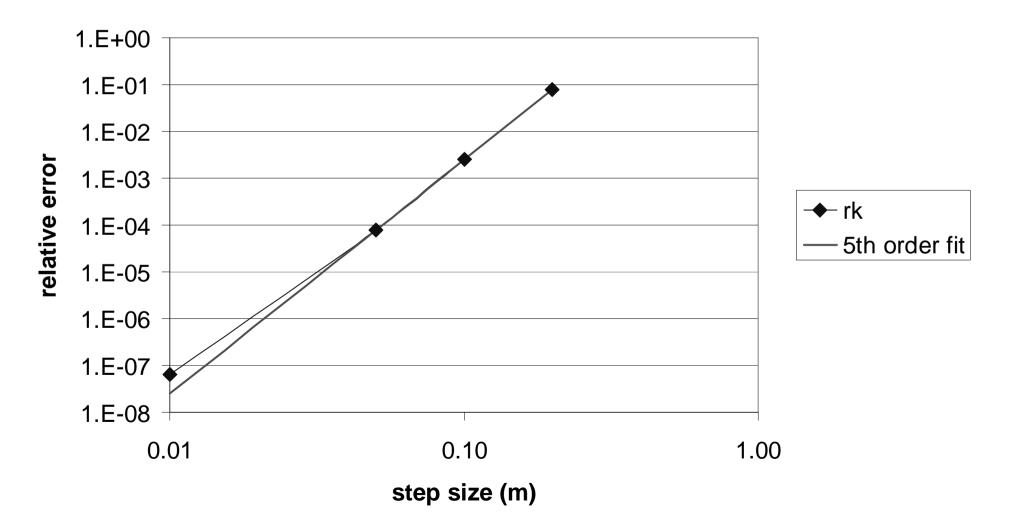




Runge Kutta errors



FOFO: relative error in angular momentum (log-log)

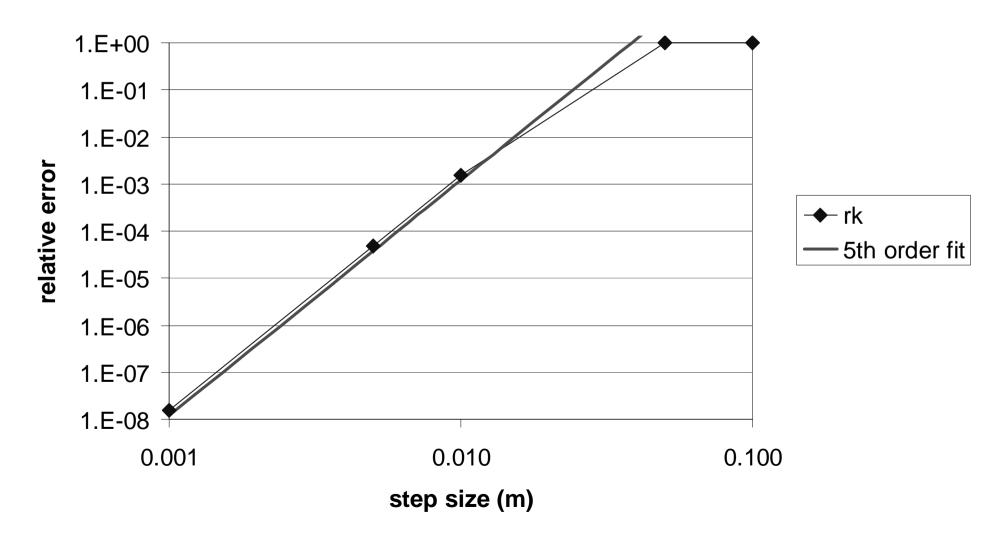




Runge Kutta errors



Solenoid: relative error in P_{\perp} (log-log)

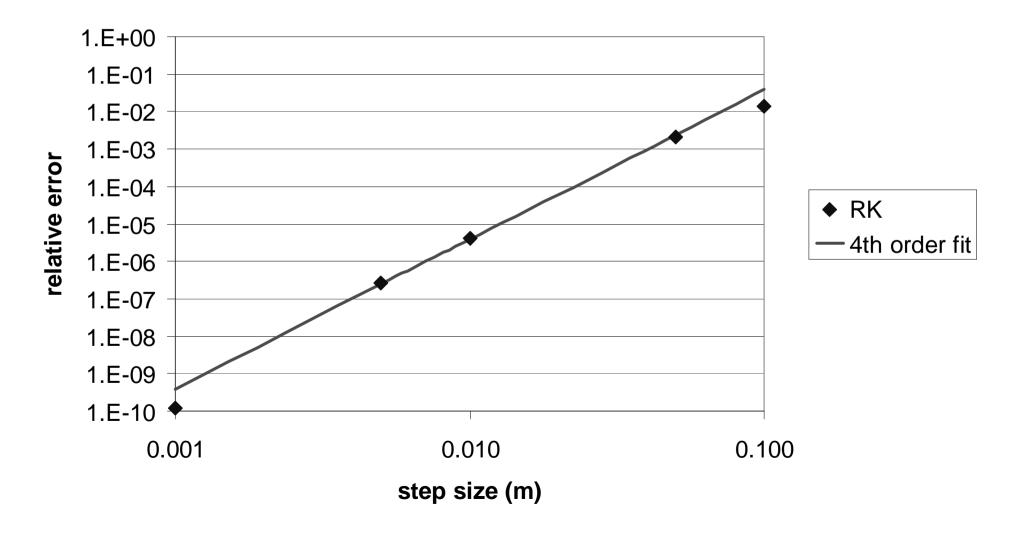




Runge Kutta Errors



Solenoid: relative error in Larmor period (log-log)

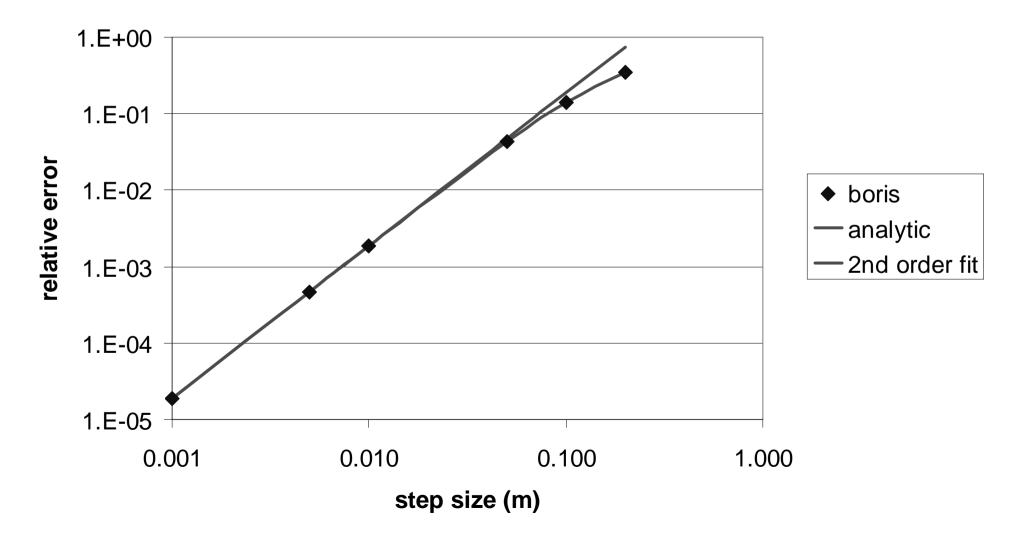




Boris errors in Solenoid



Solenoid: relative error in Larmor period (log-log)





P. H. Stoltz, G. Penn, J. Wurtele, and J. R. Cary, MC Note 229.J. Boris, *Proc. of the 4th Conf. on Numerical Simulation of Plasma* (NRL, 1970).A. Dullweber, B. Leimkuhler, and R. McLachlan, J. Chem. Phys. 15 (1997) 5840.