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# **3D PARTICLE SIMULATION OF BEAMS USING THE WARP CODE: TRANSPORT AROUND BENDS**<sup>†</sup>

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WARP is a discrete-particle simulation program which was developed for studies of space-chargedominated ion beams. It combines features of an accelerator code and a particle-in-cell plasma simulation. The code architecture, and techniques employed to enhance efficiency, are briefly described. Current applications are reviewed. In this paper we emphasize the physics of transport of three-dimensional beams around bends. We present a simple bent-beam PIC algorithm. Using this model, we have followed a long, thin beam around a bend in a simple racetrack system (assuming straight-pipe self-fields). Results on beam dynamics are presented; no transverse emittance growth (at mid-pulse) is observed.

# 1 INTRODUCTION

In a Heavy Ion inertial confinement Fusion (HIF) driver, space-charge-dominated beams are to be accelerated and transported over large distances. WARP<sup>1,2</sup> is a discrete-particle simulation program developed specifically for the study of emittance growth resulting from the nonlinear self-fields of such beams.

In the sections below, the code's architecture is reviewed, and its numerical methods (which have been described previously) are summarized. Some of the applications of the code are then reviewed. The new bent-beam algorithm currently in use is described, and some early results presented. Although somewhat preliminary in nature, these simulations suggest that it is possible to transport a space-charge-dominated beam around a fairly sharp bend without inducing unacceptable emittance growth. Finally, some directions of our ongoing research are indicated.

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# 2 CODE OVERVIEW

The WARP code contains a number of distinct parts, including: a 3d particle-in-cell (PIC) package, WARP6, which uses a "warped Cartesian" mesh to describe bends; an r,z (axisymmetric) PIC package, WARPRZ; an envelope-equation solver (used for particle loading and lattice testing); and facilities for initialization, diagnostics, etc. In this paper we focus exclusively on WARP6.

The code has been built up from independent modules; each can be loaded and used as a code by itself. This is made possible through the facilities of BASIS,<sup>3</sup> which provides a development system and an interactive user interface. WARP runs on Cray computers and on Sun workstations.

The simulation takes place in the laboratory frame. The computational mesh fills a moving window and is laid down anew at each timestep. The self-field is electrostatic in the beam frame. At present (slow beams), we use these fields directly. For faster beams we will obtain the lab-frame self- $\mathbf{E}$  and  $\mathbf{B}$  via a Lorentz transformation.

Self-field boundary conditions are those of a square metal pipe in x,y; a round pipe is an option. Periodicity of the electrostatic potential  $\phi$  in z is assumed; this is reasonable provided the beam does not approach the ends of the mesh, or is itself assumed periodic. The algorithm for the self-field conserves axial momentum algebraically; energy is not identically conserved, but is instead used as a check on the accuracy of the calculation.

The envelope solver is used for particle loading, since it is important to be able to initialize a beam which is approximately "matched"; it is also useful for the initial testing of an accelerator lattice specification.

It is necessary to follow beams over long distances, and 3d simulations are computationally challenging. Therefore, efficiency was a critical element of the code's design and has been obtained via a number of means, summarized here.

In leapfrog motion, if a particle were to land within a sharp-edged focusing or bending element on four steps while its neighbor did so on only three, they would receive dramatically different impulses. Thus, the advance is modified to incorporate "residence corrections" for element forces; these take into account the fraction of the velocity advance step actually spent within the element. This allows much bigger steps than otherwise would be possible.

No mesh arrays for the electric field components are used; instead,  $\phi$  is gathered from cells in the neighborhood of each particle, and then differenced on a particleby-particle basis. This saves the space of three 3d arrays.

The FFT field-solver is fully vectorized and uses no scratch space. The particle advance is vectorized. Deposition of the charge density  $\rho$  is vectorized with length 8, over cells overlapped by each particle.

# **3 SUMMARY OF APPLICATIONS**

In addition to the bent-beam modeling described herein, current studies include the following areas; some of these are described elsewhere, and are merely noted:

#### 3.1 Drift Compression.

Drift compression is current enhancement resulting from a head-to-tail velocity gradient or "tilt". We are examining this process, as well as the effects of quadrupole magnet misalignments on it. Relatively small misalignments can, depending on how they are phased with respect to each other, lead to significant off-axis displacements. Image forces and fringing fields can then induce emittance growth.<sup>4</sup> A goal of this work is to gain an understanding of how rapidly the beam may be compressed without unacceptable emittance degradation. The early stages of this work have been described previously.<sup>2</sup>

#### 3.2 Equilibration.

We are examining processes that lead to the transfer of thermal energy between transverse and longitudinal motions. WARP simulations indicate that, for certain ranges of physical parameters (phase advance, etc.), a beam initialized colder in z than in x,y will heat rapidly in z until  $T_z$  is a large fraction of  $T_{x,y}$ . In one such test, described previously,<sup>2</sup> a periodic beam was followed for 150 full focusing-defocusing (FODO) lattice periods (9000 steps). The beam had a phase advance per cell of  $\sigma_0 = 60^\circ$ , depressed to  $\sigma = 20^\circ$  by space charge. It was initially cold longitudinally. A rapid conversion of transverse thermal energy into longitudinal thermal energy at the beginning of the run (growth rate on the order of 1/10 of the plasma frequency,  $\omega_n$ ) was followed by a slower collisional phase, during which  $T_z$  and  $T_{x,y}$  rise together.

We have run many simulations to demonstrate that numerical artifacts are not a likely source of the heating. It has been observed over significant ranges of numerical and physical parameters (including runs with meshes which resolve the initial axial Debye length), and in finite-length beams. Furthermore, it disappears for certain ranges of physical parameters, in particular when  $\omega_p$  is reduced sufficiently, e.g.,  $\sigma = 45^{\circ}$ . Such axial heating might arise from a mechanism analogous to an anisotropy-driven "Harris" mode in a magnetized plasma.<sup>5</sup> Previous researchers have noted collective equilibration mechanisms (in other geometries).<sup>6,7</sup> We also have made a preliminary observation of rapid equilibration in an axisymmetric beam, using WARPRZ.

The equilibration issue is of interest because it is important to identify in advance all constraints on driver design. For present driver designs, equality of  $T_z$  and  $T_{x,y}$ is acceptable, but a significantly greater thermal spread in z might make final focus problematic. As driver parameters evolve it may be desirable to revisit this issue.

# 3.3 Axial Confinement and Nature of Equilibria.

To follow a finite-length beam for a long time, it is necessary to apply an axial confining force; without it the beam will blow apart electrostatically. Considerable work has been done on this subject, and we have had success in modeling axially confined beams that remain "quiescent" over the entire run. We have followed beams for up to 175 lattice periods without significant numerical emittance degradation. This work is described in a companion paper.<sup>8</sup>

3.4 Longitudinal Stability With Nonzero Gap Impedance.

This work is being carried out using WARPRZ, and is described in a companion paper.<sup>9</sup>

# 4 BENT-BEAM ALGORITHM

We have developed a family of techniques for modeling bends. These are based upon following a particle's position and velocity in a sequence of rotated inertial (laboratory) frames. In the "exact" method (not yet in WARP), no pseudo-forces are used. The scheme conserves phase-space area identically, and implies no large-aspect-ratio (gentle bend) expansion. We have formulated 3d and 2d (transverse) algorithms of this type.<sup>10</sup>

In this paper, we concentrate upon the "simplified" method, which is now in use. The method is inexact, but is accurate enough for our present purposes. It folds the velocity rotation part of the coordinate transformation into the physical dipole field as a "pseudo-gyrofrequency." Minor changes to the position advance are also needed. "Intermediate" methods are also possible, and in fact may be optimal for many purposes.

The radius of curvature of the reference orbit (usually the vessel centerline) is  $r_*$ . Time is the independent variable for particle orbits. The conventional (for accelerator codes) independent variable s is, in WARP, a dependent variable for orbits, as are the transverse coordinates x,y. In the straight sections,  $s \equiv z$ , while in the bends,  $s \equiv -r_*\theta$ . The "radial" coordinate is  $x \equiv r - r_*$ ; the unit vectors  $\hat{x}$  and  $\hat{s}$  evolve as a particle moves, and are different for each particle. This geometry is depicted in Figure 1(a). The axial speed is  $v_z = -r\dot{\theta}$ . The axial position (projection onto the centerline) is advanced in time using

$$ds/dt = -r_*\dot{\theta} = (r_*/r)v_z. \tag{1}$$

In this derivation of the simplified bent-beam algorithm, we show that the necessary "pseudo-forces" are equivalent to a velocity rotation. The radial coordinate increases for an unaccelerated particle due to "centrifugal force"; with q/m = 1:

$$\frac{d}{dt}v_{\mathbf{x}} = \frac{v_{z}^{2}}{r_{\star} + x} + \{\mathbf{E} + \mathbf{v} \times \mathbf{B}\}_{\mathbf{x}}.$$
(2)

For the "Coriolis force," use conservation of angular momentum:

$$\frac{d}{dt}r^{2}\dot{\theta} = r\{\mathbf{E} + \mathbf{v} \times \mathbf{B}\}_{\theta}.$$
(3)

After algebra:

$$\frac{d}{dt}v_z = -\frac{v_z v_x}{r_* + x} + \{\mathbf{E} + \mathbf{v} \times \mathbf{B}\}_z.$$
(4)



FIGURE 1(a) Geometry of the bent-beam algorithm. (b) Effect of folding the pseudo-gyromotion into the bending field.

Note there is no factor of two in the Coriolis force because the frame of reference differs from that in the classical analysis; particle velocities retain their laboratory-frame magnitudes. Considering only the "pseudo-force" terms, the rate of change of the velocity angle is consistent with the requirement  $\dot{\theta} = -v_z/r$ :

$$\frac{d}{dt}\arctan\left(\frac{v_x}{v_z}\right) = \frac{v_z}{r_* + x}.$$
(5)

The transformation is a pure rotation; one can readily verify that  $d|\mathbf{v}|/dt = 0$ .

Implementation in WARP was straightforward. We needed only to augment the dipole (bending) field with a "pseudo-gyrofrequency":

$$B_{y, \text{ dipole}} \leftarrow B_{y, \text{ dipole}} - \frac{m}{q} \frac{v_z}{r_* + x}.$$
(7)

This folds the necessary back-rotation into existing coding. The net effect is depicted in Figure 1(b). We must also advance the axial position s using Equation (1). The algorithm is inexact because  $v_z$  and x change during the step; "residence corrections" on entry to and exit from bends are necessary.

To date, we have only implemented and used "bent particle dynamics." Poisson's

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equation is still solved in a straight square pipe. The necessary changes to the field solution have been worked out but are not yet in the code.

Until recently we had only a limited, highly regular lattice; a general lattice of sharp-edged elements will soon be available.

# 5 BENT-BEAM RESULTS

A pair of runs shows the effects of a bend on a space-charge dominated beam; one run includes a bend while the other does not. Both employ a FODO lattice with period 1.2 m. The undepressed tune (phase advance per lattice period) is  $\sigma_0 = 60^\circ$ , the depressed tune is  $\sigma = 20^\circ$ ; and the ratio of space charge to emittance forces is about 6:1. The runs used a  $64 \times 64 \times 128$  mesh, with square walls at  $\pm 9$  cm. A total of 54 160 simulation particles were employed; the time-step size  $\Delta t$  corresponded to 2 cm/step. The runs used "quiet start" loading<sup>11</sup> of particles into an axially cold "cigar" beam with a "K–V" transverse distribution.<sup>12</sup> The beams corresponded to singly-charged carbon, 5.2 amperes, with energy 10<sup>7</sup> eV (speed  $1.27 \times 10^7$  m/s = 0.042 c). The bend begins at z = 1.2 m and is 3.6 m (3 FODO periods) long. The bend radius is  $r_* = 3.6/\pi = 1.146$  m, and the bend field 1.375 T. The beam ends are unconfined, and expand due to the electrostatic repulsion; the rarefaction does not have time to reach the beam center by the time the run finishes. Both runs are followed for 300 steps, to z = 6 m. Each used less than 15 minutes of Cray X/MP time.

Figure 2 shows the initial beam particle distribution; the upper half is a "top view" (z-x projection) of the beam, while the lower half is a "side view" (z-y projection). Figure 3 depicts the situation when the beam center has moved 3 m (the halfway



FIGURE 2 Initial beam particle distribution in 3-space; upper half is top view z-x, lower half is side view z-y.



FIGURE 3 Beam particle distribution midway through run for (a) straight system; (b) bent system.

point) for both "straight" and "bent" runs. In the latter, the focusing effect due to the "sector" nature of the dipole field is evident. In Figure 4 the situation at the end of the run is shown. A radially inward displacement of the beam tail (toward negative x) is evident in the bent run; this is a consequence of the tail's reduced axial velocity, which results from the axial expansion. Figure 5 shows the evolution of the mid-pulse x-x' emittance; note that these two plots use different scales for the ordinates, and both suppress the zero. The "spikes" as the beam center moves through quadrupole magnets are an artifact of the emittance diagnostic, which uses moments over a subset of particles which has finite axial length. We are encouraged by the absence of emittance growth in this run, which uses a sharp bend that was not designed to be a "matched" system. Longer runs have followed axially confined beams around as many as three bends without large emittance growth; this work will be described in a future publication.



FIGURE 4 Beam particle distribution at end of run for (a) straight system; (b) bent system.



FIGURE 5 Mid-pulse x-x' emittance vs. time for (a) straight system; (b) bent system.

## **6** FUTURE DIRECTIONS

Recent improvements to WARP include a round-pipe model (which applies a capacity matrix<sup>8</sup> to each axial Fourier mode in turn) and an axially confining force (which replaces the "ears" incorporated in real accelerating pulses).<sup>8</sup> WARP's r,z model is being used to study longitudinal stability, compression, and equilibration.<sup>9</sup>

Current and planned projects include: studying bends and recirculation with realistic lattices and bent-field corrections; modeling the Lawrence Berkeley Laboratory multiple-beam experiment MBE-4, including the effects of acceleration and compression, with both constant and intermittent "ears;" and 3d modeling to understand the effects of aberrations in final focus (wherein the beams are not truly long and thin). We plan to model beams generated by a "real" injector, instead of loading with an ideal distribution function. Many code improvements for efficiency, generality, and convenience will be folded in as the code develops.

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