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Particle-In-Cell simulation of laser irradiated two-component microspheres in 2 and 3 dimensions

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ABSTRACT

We examine proton acceleration from spherical carbon-hydrogen targets irradiated by a relativistic laser pulse. Particle-In-Cell (PIC) simulations are carried out in 2 and 3 dimensions (2D and 3D) to compare fast proton spectra. We find very different final kinetic energies in 2D and 3D simulations. We show that they are caused by the different Coulomb fields in 2D and 3D. We propose a correction scheme for the proton energies to test this hypothesis. In the case of sub-focus diameter targets comparison of corrected 2D energies with 3D results show good agreement. This demonstrates that caution is required when modeling experiments with simulations of reduced dimensionality.

1. Overview

1.1. Motivation

Recently many simulations and experiments have been carried out to investigate the properties of fast particles emitted from plasmas that are produced when mass limited targets are irradiated by relativistic pico- or femtosecond laser pulses (comp. [6,10,8,14,11]). The long-term goal of these investigation is to develop new accelerating technologies and radiation sources. The cooperative effort of experiment and simulation in this area is fruitful for both fields, as comparing the results can help to improve the methods and models used by both. The simulations presented here are carried out in the context off recent experiments using levitating targets by the group of Jörg Schreiber at LMU (Garching). We make use of the PSC [7,1,5,4] developed at the chair for Computational Physics at LMU to simulate the interaction of isolated spherical two-component targets with relativistic laser pulses lasting 55 fs. We ask whether a two dimensional representation can yield satisfying results by comparing these to 3D simulations.

1.2. Set-up

The target in the associated experiment is a polystyrene ball of varying diameter consisting of a 1:1 mixture of hydrogen and carbon. It is modeled by a spherical plasma (fully ionized ab initio) with the corresponding amount of electrons (1 per hydrogen, 6 per carbon ion). We scan the properties of the resulting proton spectrum for different target sizes, ranging in diameter from 500 nm to 20 μm.

A 55 fs FWHM Gaussian laser pulse ($\lambda = 1$ μm, 8.3 μm FHWM transversal Gaussian profile) is seeded into the simulation domain.

A rectangular simulation domains of 2048×4096 cells in 2D and $2688 \times 1536 \times 1536$ cells for 3D are used corresponding to a spatial resolution of 64 cells per micrometer. For 2D simulations, boundaries are open at the slice edges and periodic on the slice planes (y direction) for fields and particles. For 3D simulations all boundaries are open. To avoid the danger of unstable plasma wave modes due to the finite resolution (as described in [15]), we use one half of the solid density n_s of polystyrene corresponding to $n_s/2 = 1.7 \cdot 10^{23}/\mathrm{cm}^3$, unless otherwise specified. For the same reason particles of the target are initiated with a Maxwellian momentum distribution corresponding to a plasma temperature in the keV range. Unless specified otherwise the initial temperature was 10 keV in the simulations.

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2. 2 Dimensional simulations

Since the computational load (machine run time and data output) is much smaller in 2D as compared to 3D, a reduced dimension model is often employed (for example [2,8,3,11]). We try to investigate whether simulations with reduced dimensions can still lead to reliable results for the scenario of a short laser pulse interacting with polystyrene microspheres as described in this paper.

The 2D geometry implies that the linear polarized laser pulse interacts with a cylinder rather than a sphere. This means that the laser can be polarized either along the cylinder axis (s-polarization, 2D-s) or perpendicular to it (p-polarization, 2D-p). We expect that the s and p cases lead to different results for otherwise same parameters.

2.1. Difference between polarization directions

The electron motion is governed by the Lorentz force leading to ponderomotive potential effects. For p-polarization significant numbers of electrons can move along the electric field and be extracted from the target, forming ultrashort bunches separated by one laser wavelength (see Fig. 1). This process is described in detail in [9]. For the s-polarization case displacement along the direction of the electric field is impossible. However, electrons still experience the $\vec{j} \times \vec{B}$-force, but do not form bunches to the same extent as in the p case. Since the $\vec{j} \times \vec{B}$-force changes sign 4 times during a laser cycle, density modulations with the separation of half a wavelength are observed. In addition the longitudinal ponderomotive force is overestimated as electrons only see the maximum field strength in the middle of the pulse (see Fig. 1). We also note that for a 2 μm target average electron temperature is 0.3 MeV for p but only 0.12 MeV for s-polarization, in both cases exceeding the initial temperature substantially.

We observe that ion kinetic energies are significantly higher in the p case than in the s case. The energy mismatch is largest for the smallest targets as is seen in Fig. 2. We attribute this mismatch in energy to a difference in electron removal, which in the p case is about 77% for 0.5 μm and 12% for a 2 μm target, but much lower in the s case (40% for 0.5 μm and 3% for 2 μm). Here we specify the percentage of electrons still present in a 16×32 μm rectangle around the target 500 fs after pulse maximum. The suppressed electron depletion in the s-case obviously implies, that a 2D-s set-up cannot model the interaction scenarios in 3D.

The rising maximal energies for smaller target radii in the p-case do not agree with experimental findings (e.g. [13,12]). Noticeably the protons continue to considerably gain kinetic energy until they reach the simulation domain boundaries. This is caused by the Coulomb field of the positive charge surplus at the target. To retrieve reliable kinetic energy spectra we conduct 3 dimensional simulations of the same set-up.

3. 3 Dimensional simulations

The same set-up was then simulated in 3D using the same resolution and pulse parameters but a shorter box ($2688 \times 1536 \times 1536$ cells) to reduce memory demand.

When comparing the p-pol 2D set-up with a slice of the 3D simulation domain along the laser electric field polarization

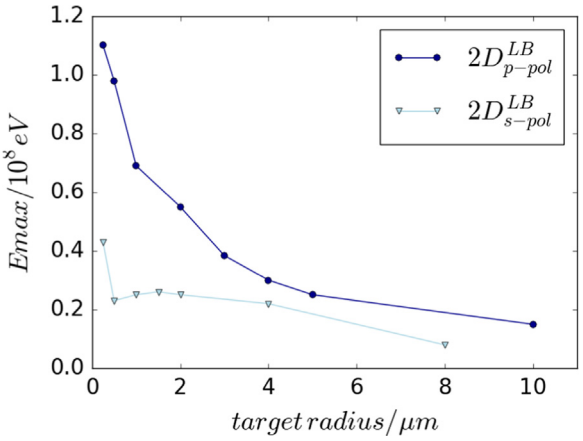

Fig. 2. Maximum proton energies for p and s polarized laser field using a long box (2048×4096, corresponding to 32 μm and 64 μm), an initial temperature of 50 keV, $n_0 = 0.25 \cdot n_s$ and $a_0 = 15$.

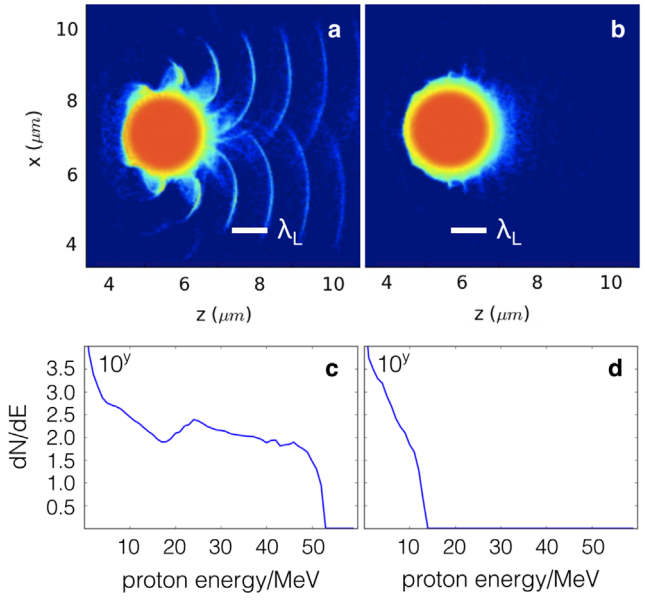

Fig. 1. 2D simulation of a 2 μm target, $a_0 = 12$, laser wavelength $\lambda = 1$ μm. Above: Qualitative comparison of electron dynamics in the p (a) and s (b) polarization case (electron density). We observe the different frequency of the density waves. Below: Proton energy spectrum 200 fs after pulse maximum (power of 10 of the number of quasi-particles, p-pol (c) and s-pol (d)).

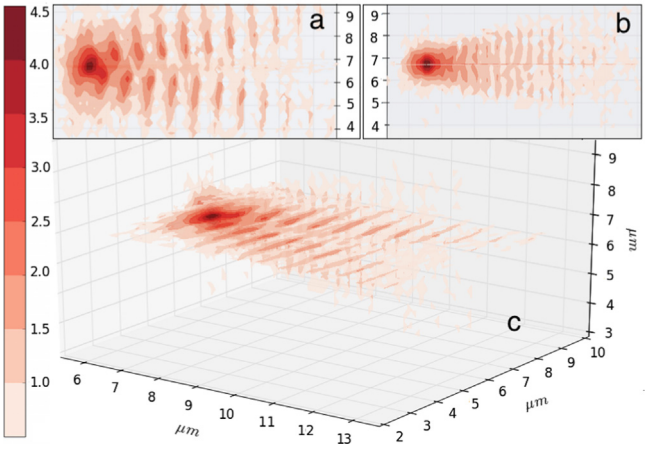

Fig. 3. 500 nm target in 3D, $a_0 = 12$. Contour plot of electron density in 50 nm slices through the center of the 3D simulation domain: (a) in polarization plane, (b) perpendicular to polarization plane and (c) both. The ultrashort electron bunches appear as in the 2D-p case. Density scale shows the powers of 10 of the number of quasi-particles in a slice volume element $10\Delta x \cdot 10\Delta x \cdot \Delta x$.