Two-dimensional quantum hydrodynamic model for the heating of a solid target using a Gaussian cluster

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Abstract

This paper presents numerical simulations to study the heating of a two-dimensional (2D) solid target under an ion cluster interaction. 2D quantum hydrodynamic (QHD) model is employed for the heating of solid target to warm dense matter on a picosecond time scale. A Gaussian cluster is used to uniformly heat the solid target to a temperature of several eV. The density and temperature of the target are calculated by a full self-consistent treatment of the QHD formalisms and the Poisson's equation. The technique described in this paper provides a method for creating uniformly heated strongly coupled plasma states.

Keywords: Cluster; Uniform heating; Warm dense matter

1. INTRODUCTION

Interaction processes of ion beam or cluster with matter are important for the applications in basic and applied physics (Patel et al., 2003; Brambrink et al., 2006; Tahir et al., 2009a), laboratory accelerator (Tahir et al., 2009b), inertial confinement fusion (ICF) (Deutsch, 1986; Deutsch et al., 1989; Hora, 2007; Hoffmann, 2008), astrophysical processes (Nettelmann et al., 2008), heavy ion beam research (Hoffmann et al., 2010; Stöckl et al., 1996; Tahir et al., 2007), and high energy density physics (HEDP) (Tahir et al., 2005; Nellis, 2006; Hoffmann et al., 2009). HEDP is understood to be the thermodynamic regime where the energy density exceeds 10^{11} J/m³ or equivalent pressure of 1 Mbar and above. Pioneering experiments of beam-matter interactions (Deutsch et al., 1989) have been conducted to investigate the stopping power for hydrogen plasma, which have demonstrated that the stopping power can increase with the increasing ionization of plasma due to the free electrons interacting with beam ions. It will be shown very soon that intense clusters may be of considerable potential interest for driving an ICF (Deutsch, 1992). Using clusters as a ICF driver, can break through the space-charge limit of ion beam, which is helpful to achieve high heating effect. Recently, several attempts have been made to investigate the beam-matter

interactions involving heavy ion (Zhao *et al.*, 2009; Hoffmann *et al.*, 2010), proton (Patel *et al.*, 2003; Tahir *et al.*, 2009a) and laser beams (Zhao *et al.*, 2012). In particular, ICF is a promising scheme, where a target matter can be irradiated by intense ion beams or clusters, and be transformed into a plasma of high density and high temperature which can initiate fusion reactions (Deutsch, 1992; Tahir *et al.*, 2005).

The basic properties of interactions between ions and matter, which play an essential role in the above scenario, have been encouraged to be explored thoroughly. The evaluations of target temperature and density are an ideal tool to investigate these interactions for ion beams or clusters immersed in a target. It is expected in such studies that target properties would be influenced significantly by the modulation of beam or cluster parameters such as particle energy, total number of particles per beam, and focal spot size. These beam parameters will be available at the new powerful heavy ion synchrotron, SIS100, at GSI Darmstadt, which will deliver strongly bunched intense beams and clusters of all stable ion species from protons up to uranium with particle energies a few MeV-GeV/u (Tahir et al., 2009a, 2011). These intense ion beams and clusters present a promising source for volumetric heating and bring up unique capabilities for HEDP. In this field, one of the special interests is warm dense matter (WDM) research, since the accelerator technology has improved sufficiently to provide intense ion beams and clusters to heat samples of target and reach the

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regime of WDM. WDM is the region in temperature (several eV) and density that can be described as strongly coupled plasma. The importance of WDM arises from its wide occurrence, namely planetary science, cold star physics, and all plasma production devices, which start from cold dense matter.

In a strong coupling plasma considering nonlinear effects, the interaction processes of ions with matter have been simulated by employing various methods, such as particle-in-cell (PIC) and molecular-dynamics (MD) (Zwicknagel et al., 1999; Hu et al., 2011, 2012). In particular, quantum hydrodynamic (QHD) model has been developed by solving the nonlinear Schördinger-Poisson or the Wigner-Poisson kinetic models, which can be devoted to studying the properties of strongly coupled plasmas. Large number of calculations have been carried out by using the QHD model since the use of this theory was introduced (Chen et al., 1995; Haas et al., 2000; Manfredi and Haas, 2001), due to the advantages that the QHD equations are quite simple for numerical studies and have a straightforward interpretation in terms of fluid quantities that are employed in classical physics. In our previous work (Zhang et al., 2011b), we have simulated the wake potential and energy loss of a test charge propagating in a one-dimensional (1D) solid target, by using the 1D QHD theory. However, the QHD theory has not been applied to examine an ion cluster interacting with a two-dimensional (2D) solid target so far.

The aim of this paper is to extend our previous work (Zhang et al., 2011b) to a 2D case with an ion cluster interaction, moreover, the energy-balance equation is considered in this work. The 2D QHD approach will be employed to analyze the isochoric heating (i.e., heating at constant volume) of aluminum solid target under an ion cluster interaction, which can heat solid density material to warm dense states on a picosecond time scale. We shall show how the temperature and density of the target are affected by the modulation of cluster properties and how the WDM is generated from a cold target. The nonlinear QHD equations will be solved by employing the flux-corrected-transport (FCT) method (Boris et al., 1993), which has proven to be an accurate and easy to use algorithm to solve nonlinear timedependent conservative equations of hydrodynamics. This paper is organized as follows. In Section 2, the nonlinear 2D QHD formalisms coupled with the Poisson's equation, are presented. In Section 3, we show numerical results of the temperature and density of the target. Finally, a short summary is given in Section 4. Gauss units will be adopted throughout the paper except specific definition.

2. QUANTUM HYDRODYNAMIC MODEL

We consider a Gaussian proton cluster of density profile n_c $(z, r, t) = n_{c0} exp(-r^2/r_c^2 - (z - V_c t)^2/l_c^2)$ with cluster radius r_c and cluster half-length l_c , propagating with constant velocity $V_c = \sqrt{2E_{cp}/m_{cp}}$ through a solid aluminum target along the z axis. Here, m_{cp} is the proton mass and E_{cp} is the proton energy. The total number of protons N_{cp} within cluster is $N_{cp} = 4\pi \int_0^t \int_0^{r_c} n_c (z, r, t) r dr dz$, functions of n_{c0} , r_c , and l_c . Note that the total number and Gaussian distribution of protons within cluster are unvaried with time *t* evaluation. We treat the valence electrons in the target as free electrons with an equilibrium density $n_0 = 7.829 \times 10^{+23}$ cm⁻³, which is aluminum solid density, immersed in a uniform background of motionless ions. The electron dynamics can be analyzed by a collective-stimulated interaction model, that is, the transmission of the cluster ions can cause dynamics polarization effect in the target.

The simulation technique is the 2D QHD model with the electron velocity field $\mathbf{u}_e(z, r, t)$, density $n_e(z, r, t)$, and energy density $W_e(z, r, t)$ by the continuity equation

$$\frac{\partial n_e}{\partial t} + \nabla \cdot (n_e \mathbf{u}_e) = 0, \tag{1}$$

the momentum-balance equation

$$\frac{\partial \mathbf{u}_e}{\partial t} + (\mathbf{u}_e \cdot \nabla) \mathbf{u}_e = \frac{e}{m_e} \nabla \Phi - \frac{\nabla P_e}{m_e n_e} + \frac{\hbar^2}{2m_e^2} \nabla \left(\frac{1}{\sqrt{n_e}} \nabla^2 \sqrt{n_e}\right) - \gamma \mathbf{u}_e,$$
⁽²⁾

and the energy-balance equation (Chen et al., 1995)

$$\frac{\partial W_e}{\partial t} + \nabla \cdot (W_e \mathbf{u}_e) = e n_e \mathbf{u}_e \cdot \nabla \Phi - \nabla \cdot (\mathbf{u}_e P_e) - \nabla \cdot \mathbf{q}_e - \frac{W_e - \frac{3}{2} n_e T_e}{\tau_w},$$
(3)

with $W_e = \frac{3}{2}n_eT_e + \frac{1}{2}m_en_eu_e^2 - \frac{\hbar^2 n_e}{24m_e}\nabla^2 \log(n_e) + O(\hbar^4)$. The above equations are closed by adding Poisson's equation

$$\nabla^2 \Phi = 4\pi e (n_e - n_0 - Z_1 n_c).$$
(4)

Here, $\nabla = \frac{\partial}{\partial r} \mathbf{e}_r + \frac{\partial}{\partial z} \mathbf{e}_z$, m_e is the electron mass, e is the elementary charge, Φ is the total electric potential, \mathbf{q}_e , T_e , and P_e are the electron heat flux, temperature and pressure, respectively, γ and τ_w are the frictional coefficients with $\tau_w = 2\gamma$. In particular, the electron pressure P_e can be calculated by finite temperature Thomas-Fermi theory (More *et al.*, 1988) and the heat flux is specified by Fourier's law $\mathbf{q}_e = -\kappa \nabla T_e$ (Chen *et al.*, 1995) with thermal conductivity κ . In the right-hand side, the second and third terms of Eq. (2) are regarded as the quantum effects, and the last terms of Eqs. (2) and (3) are the frictional forces with γ and τ_w obtained from our previous work (*Zhang et al.*, 2011*a*).

The above Eqs. (1)–(4) involve a self-consistent determination of the electron density, velocity, and temperature, and the total electric field $\mathbf{E} = -\nabla \Phi$ with the components E_r and E_z . The FCT method is adopted to numerically solve Eqs. (1)–(3) by split-time integration from the initial time t = 0 when the values of all quantities are known. The Poisson's equation is solved by the successive over relaxation (SOR) method.

In the following section, we obtain the numerical values of the density and temperature of the target under the effect of a Gaussian proton cluster. In the simulation, $n_{c0} = n_0 \times 10^{-9}$ and the cluster is projected from (z = 0, r = 0) at t = 0. Our numerical results are obtained for solid aluminum with initial solid density n_0 and initial temperature $T_e = 0$ eV.

3. NUMERICAL RESULTS

Figure 1 show the density as functions of r and z at 5 ps Figure 1a and 10 ps Figure 1b for $E_{cp} = 20$ MeV, and at 5 ps Figure 1c and 10 ps Figure 1d for $E_{cp} = 10$ MeV. Here, $r_c = 3 \mu m$ and $l_c = 15 \mu m$. The density demonstrates a clear time and space variations and wake effects are generated behind the cluster. A comparison between Figures 1a and 1b shows that the maximum density is almost unchanged with the cluster penetration time. As the cluster penetrates deeper into the target, the perturbed density region becomes larger. When $E_{cp} = 20$ MeV in Figures 1a–b, the range of the cluster is longer, while the maximum density is smaller, in comparison with that for $E_{cp} = 10$ MeV in Figures 1c–1d. The perturbed density decreases and tends to reach an initial equilibrium value with the increasing distance behind the cluster.

Temperature corresponding to Figures 1a–1d is shown in Figures 2a–2d. Cluster protons impinging on the target deposit their energy in the target that leads to an increase in the temperature. Thus the spatial profile of the temperature is similar to the Gaussian distribution of the cluster. In the

cluster heated region, temperature of $0.6 \le T_e \le 1$ eV is obtained in Figures 2a–2b and that of $1 \le T_e \le 1.5$ eV is obtained in Figures 2c–2d, where WDM is generated. The WDM temperature range is roughly $T_e = 1.0-20$ eV for aluminum. As the cluster transmission time increases, a larger volume of the target has been heated. When $E_{cp} =$ 20 MeV in Figures 2a–2b, the range of the cluster is longer and the heated region is larger, while the maximum temperature is smaller, in comparison with those for $E_{cp} = 10$ MeV in Figures 2c–2d. Moreover, more uniform temperature is obtained in Figs. 2(a-b), as compared to that shown in Figures 2c–2d. Note that nearly uniform temperature of several eV is achieved in the cluster heated region and the target can be converted into WDM on a picosecond time scale, which is important for the study of HEDP and ICF research.

To further examine the influence of the cluster parameters on the density, we present in Figures 3a–3c the density as functions of r and z with $r_c = 1 \mu m$, $l_c = 10 \mu m$ Figure 3a, $r_c = 3 \mu m$, $l_c = 15 \mu m$ Figure 3b and $r_c = 6 \mu m$, $l_c = 20 \mu m$ Figure 3c. Here, $E_{cp} = 20$ MeV and t = 10 ps. In Figures 3a–3b the maximum density decreases with decreasing r_c and l_c due to reduction in n_c that has a direct ratio relation with r_c and l_c .

Temperature corresponding to Figures 3a–3c, is presented in Figures 4a–4c. It is clearly seen in Figures 4a–4c where the cluster heated volume becomes larger in radial direction and the maximum temperature increases with increasing r_c and l_c . The temperature exhibits a nearly uniform distribution in the cluster heated region, which has values of $T_e = 0.4 \sim 2$ eV. Moreover, the maximum temperature (2 eV) is in favor of important experiments on HEDP.



Fig. 1. (Color online) The density at 5 ps (a) and 10 ps (b) for $E_{cp} = 20$ MeV, and at 5 ps (c) and 10 ps (d) for $E_{cp} = 10$ MeV. Here, $r_c = 3 \mu m$ and $l_c = 15 \mu m$.



Fig. 2. (Color online) The temperature at 5 ps (a) and 10 ps (b) for $E_{cp} = 20$ MeV, and at 5 ps (c) and 10 ps (d) for $E_{cp} = 10$ MeV. Here, $r_c = 3 \mu m$ and $l_c = 15 \mu m$.

The density and temperature are significantly impacted by the cluster properties, as a result, such a influence is also shown in the total electric field. Corresponding to Figure 3b and Figure 3c, Figures 5a–5b and Figures 5c–5d show E_r and E_z as functions of r and z, with E_r Figure 5a and E_z Figure 5b for $r_c = 3 \ \mu m$, $l_c = 15 \ \mu m$ and E_r Figure 5c and E_z Figure 5d for $r_c = 6 \,\mu\text{m}$, $l_c = 20 \,\mu\text{m}$. Here, $E_{cp} = 20 \,\text{MeV}$ and $t = 10 \,\text{ps}$. The total electric field is mainly determined by the Gaussian cluster, so the initial values $E_r = 0$, $E_z = 0$ are changed only at and around the cluster center, as shown in Figures 5a–5d. The volume of nonzero electric fields is smaller for a lower r_c and l_c due to reduction in n_c . The spectrum



Fig. 3. (Color online) The density with $r_c = 1 \ \mu m$, $l_c = 10 \ \mu m$ (a), $r_c = 3 \ \mu m$, $l_c = 15 \ \mu m$ (b) and $r_c = 6 \ \mu m$, $l_c = 20 \ \mu m$ (c). Here, $E_{cp} = 20 \ \text{MeV}$ and $t = 10 \ \text{ps}$.



Fig. 4. (Color online) The temperature with $r_c = 1 \ \mu\text{m}$, $l_c = 10 \ \mu\text{m}$ (a), $r_c = 3 \ \mu\text{m}$, $l_c = 15 \ \mu\text{m}$ (b) and $r_c = 6 \ \mu\text{m}$, $l_c = 20 \ \mu\text{m}$ (c). Here, $E_{cp} = 20 \ \text{MeV}$ and $t = 10 \ \text{ps}$.



Fig. 5. (Color online) $E_r(\mathbf{a})$ and $E_z(\mathbf{b})$ for $r_c = 3 \mu m$, $l_c = 15 \mu m$ and $E_r(\mathbf{c})$ and $E_z(\mathbf{d})$ for $r_c = 6 \mu m$, $l_c = 20 \mu m$. Here, $E_{cp} = 20$ MeV and t = 10 ps.

of E_r displays an antisymmetry, while E_z is symmetric, with respect to the axis (r = 0).

4. CONCLUSIONS

Numerical simulations of interaction of a Gaussian proton cluster with 2D aluminum solid target are carried out. 2D QHD model is employed for the heating of solid target. A full self-consistent treatment of this problem requires coupling of the QHD formalisms with the Poisson's equation. The density and temperature of the target and the total electric field, are calculated by solving the QHD equations and the Poisson's equation. The FCT method is adopted to numerically solve the nonlinear QHD equations. The Poisson's equation is solved by the SOR method. As the cluster penetrates into the target, a uniform temperature distribution is displayed in the cluster heated region. It is interesting to note that the target is heated to a few eV and a WDM matter is generated in the cluster heated region on a picosecond time scale. Therefore, one can access the very interesting but almost unexplored strongly coupled plasma state in this work.

In the near future, this work could be modified to consider interactions of an intense ion beam pulse with solid target by using double QHD model for beam and target particles, where high energy density (HED) matter can be generated. The developments of high-quality well-focused strongly bunched intense ion beams, from a facility of antiproton and ion research (FAIR) at GSI Darmstadt, can lead to generating samples of HED matter through isochoric and uniform heating of solid targets by such intense beams.

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