Spinodal Amplification of Density Fluctuations in Fluid-Dynamical Simulations of Relativistic Nuclear Collisions

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Extending a previously developed two-phase equation of state, we simulate head-on relativistic leadlead collisions with fluid dynamics, augmented with a finite-range term, and study the effects of the phase structure on the evolution of the baryon density. For collision energies that bring the bulk of the system into the mechanically unstable spinodal region of the phase diagram, the density irregularities are being amplified significantly. The resulting density clumping may be exploited as a signal of the phase transition, possibly through an enhanced production of composite particles.

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Strongly interacting matter is expected to possess a rich phase structure. In particular, compressed baryonic matter may exhibit a first-order phase transition that persists up to a certain critical temperature [[1](#page-4-0)], and experimental efforts are under way to search for evidence of this phase transition and the associated critical end point $[2-4]$ $[2-4]$ $[2-4]$.

For this endeavor to be successful, it is important to identify observable effects that may serve as signals of the phase structure. This is a challenging task, because the colliding system is relatively small, nonuniform, far from global equilibrium, and rapidly evolving, features that obscure the connection between experimental observables and the idealized uniform equilibrium matter described by the equation of state. Therefore, to understand how the presence of a phase transition may manifest itself in the experimental observables, it is necessary to carry out dynamical simulations of the collisions with suitable transport models.

A first-order phase transition introduces spinodal instabilities [\[5](#page-4-3)–[8\]](#page-4-4) and the associated nonequilibrium dynamics may generate observable fluctuations in the chiral order parameter $[9,10]$ $[9,10]$ and the baryon density $[11-14]$ $[11-14]$ $[11-14]$ $[11-14]$ $[11-14]$. In order to explore this latter prospect, we simulate nuclear collisions with finite-density fluid dynamics, using a previously developed two-phase equation of state. Fluid dynamics has the important advantage that the equation of state appears explicitly, contrary to most microscopic transport models where it is often unknown or is cumbersome to determine. Because it is essential to incorporate finite-range effects when seeking to describe the spinodal phase decomposition [\[5](#page-4-3)[,6,](#page-4-9)[8](#page-4-4)[,15](#page-4-10)], we introduce a gradient term into the expression for the local pressure in a nonuniform medium. This term ensures that two coexisting bulk phases will develop a diffuse interface, and we calculate the associated temperature-dependent tension. The gradient term also causes the dispersion relation for the collective modes in the unstable phase region to exhibit a maximum, as is a characteristic feature of spinodal decomposition [[5\]](#page-4-3). Thus we obtain a transport model that has an explicitly known two-phase equation of state and that treats the associated physical instabilities in a numerically reliable manner. This is the first time a transport model with these key characteristics has been developed for high-energy nuclear collisions.

As a first application of this novel tool, we simulate head-on collisions of lead nuclei at various energies. For a certain window of collision energies, several GeV per nucleon, the bulk of the system will reside within the spinodal region of the phase diagram for a sufficient time to allow the associated instabilities to enhance the initial density irregularities, a phenomenon that was exploited previously to obtain experimental evidence of the nuclear liquid-gas phase transition $[5,16]$ $[5,16]$ $[5,16]$. As a quantitative measure of the effect, we extract the moments of the density distribution and compare the degree of enhancement with what would result in the absence of the phase transition.

In order to obtain a suitable equation of state, we employ the method developed in Ref. [\[8\]](#page-4-4). Thus we work (at first) in the canonical framework, and, for a given T , we obtain the free energy density $f_T(\rho)$ in the phase coexistence region by performing a suitable spline between two idealized systems (either a gas of pions and interacting nucleons or a bag of gluons and quarks) held at that temperature. In Ref. [\[8](#page-4-4)], the focus was restricted to subcritical temperatures $T < T_{\text{crit}}$, so for each T the spline points were adjusted such that the resulting $f_T(\rho)$ would exhibit a concave anomaly; i.e., there would be two densities, $\rho_1(T)$ and $\rho_2(T)$, for which the tangent of $f_T(\rho)$ would be common. This ensures phase coexistence, because then the chemical potentials $\mu_T = \partial_{\rho} f_T(\rho)$ match, $\mu_T(\rho_1) =$ $\mu_T(\rho_2)$, as do the pressures $p_T(\rho_1) = p_T(\rho_2)$. In the present work, we have extended the equation of state to $T>T_{\text{crit}}$ by using splines that have no concavity, as is characteristic of single-phase systems. After having thus constructed $f_T(\rho)$ for a sufficient range of T and ρ , we may obtain the pressure $p_T(\rho) = \rho \partial_\rho f_T(\rho) - f_T(\rho)$ and the energy density $\varepsilon_T(\rho) = f_T(\rho) - T \partial_T f_T(\rho)$ by suitable interpolation and then tabulate the equation of state $p_0(\varepsilon, \rho)$, on a convenient Cartesian lattice.

The resulting phase diagram in the (ε, ρ) phase plane is illustrated in Fig. [1](#page-1-0). At a given net baryon density ρ , the lower bound on the energy density is given by $\varepsilon_{T=0}(\rho)$, which represents the effect of compression. The lower and upper boundaries of the phase coexistence region are traced out by $\rho_1(T)$ and $\rho_2(T)$, respectively. Uniform matter is thermodynamically unstable inside this region; while it is mechanically metastable near the phase coexistence boundaries, it loses mechanical stability when the speed of sound vanishes and density irregularities are amplified inside this spinodal boundary.

It is important to recognize that the model describes the instabilities not only in the unstable spinodal region but also those in the surrounding metastable region in which finite seeds are required for amplification to occur (yielding nucleation or bubble formation [\[13\]](#page-4-12)). Thus, density irregularities may be amplified by the metastable region as well, and any clumping generated inside the spinodal region may be further enhanced as the system expands through the nucleation region.

When the dynamical evolution is governed by ideal fluid dynamics (see below), the instability boundary is characterized by the vanishing of the isentropic sound speed, $v_s = 0$, where $v_s^2 \equiv (\rho/h)(\partial p/\partial \rho)_{s/\rho}$, with s being the entropy density and $h = p + \varepsilon$ the enthalpy density. For dissipative evolutions, the finite heat conductivity causes the instability region to widen as the boundary is then characterized by the vanishing of the isothermal sound speed, $v_T = 0$, where $v_T^2 \equiv (\rho/h)(\partial p/\partial \rho)_T \le v_s^2$. These instability boundaries are also indicated in Fig. [1](#page-1-0).

FIG. 1 (color online). The (ε, ρ) phase diagram: $\varepsilon_{T=0}(\rho)$, the minimal ε for a given ρ (black line); the phase-coexistence boundary (red line) ending at the critical point (c. p.), as well as the isothermal spinodal boundary where $v_T = 0$ (green line) and the isentropic spinodal boundary where $v_s = 0$ (blue line). Points corresponding to counterstreaming Lorentz-contracted nuclei are shown for various energies E_{lab} , indicated in A GeV (circles). The two contours (dashed line) show where the points of maximum compression are concentrated in ensembles of collisions simulated with fluid dynamics for $E_{\text{lab}} = 2$ or 5 A GeV.

In order to ascertain the dynamical effects of the firstorder phase transition, we construct a one-phase partner equation of state by replacing, for any $T < T_{crit}$, the actual $f_T(\rho)$ by the associated Maxwell construction, $f_T^M(\rho) =$ $f_T(\rho_i) + (\rho - \rho_i)\mu(\rho_i)$, through the coexistence region $\rho_1(T) < \rho < \rho_2(T)$, where $f_T^M(\rho) < f_T(\rho)$.

For our present investigation, we describe the evolution of the colliding system by ideal fluid dynamics, because dissipative effects are not expected to play a decisive role for the spinodal clumping [\[15\]](#page-4-10). The basic equation of motion expresses four-momentum conservation, $\partial_{\mu}T^{\mu} = 0$, where $T^{\mu\nu}(x) = [p(x) + \varepsilon(x)]u^{\mu}(x)u^{\nu}(x)$ – $p(x)g^{\mu\nu}$ is the stress tensor. It is supplemented by the continuity equation $\partial_{\mu}N^{\mu} = 0$ for the baryon charge current density $N^{\mu}(x) = \rho(x)u^{\mu}(x)$, where $u^{\mu}(x)$ is the fourvelocity of the fluid. These equations of motion are solved by means of the code SHASTA [\[17\]](#page-4-13) in which the propagation in the three spatial dimensions is carried out consecutively.

A proper description of spinodal decomposition requires that finite-range effects be incorporated [[5](#page-4-3),[6](#page-4-9)]. Therefore, following Refs. $[8,15]$ $[8,15]$ $[8,15]$ $[8,15]$ $[8,15]$, we write the local pressure as

$$
p(\mathbf{r}) = p_0(\varepsilon(\mathbf{r}), \rho(\mathbf{r})) - a^2 \frac{\varepsilon_s}{\rho_s^2} \rho(\mathbf{r}) \nabla^2 \rho(\mathbf{r}), \qquad (1)
$$

where we recall that $p_0(\varepsilon, \rho)$ is the equation of state, the pressure in uniform matter characterized by ε and ρ . Furthermore, $\rho_s = 0.153/\text{fm}^3$ is the nuclear saturation density, and $\varepsilon_s \approx m_N \rho_s$ is the associated energy density. The strength of the gradient term is then conveniently governed by the length parameter a.

The gradient term will cause a diffuse interface to develop when matter of two coexisting phases is brought into physical contact. The associated interface tension can readily be determined from the equation of state [\[8\]](#page-4-4)

$$
\sigma_T = a \int_{c_1(T)}^{c_2(T)} [2\varepsilon_s \Delta f_T(c)]^{1/2} dc,\tag{2}
$$

where $c \equiv \rho / \rho_s$ denotes the degree of compression and $\Delta f_T(c) \equiv f_T(c) - f_T^M(c)$ is the difference between the actual free energy density at the specified compression and that associated with the corresponding Maxwell construction. The tension σ_T decreases steadily as T is raised and finally vanishes at T_{crit} ; it is displayed in Fig. [2](#page-2-0) as obtained for various values of the range a.

Furthermore, uniform matter inside the spinodal region (where v_s^2 < 0) is mechanically unstable, and density ripples of wave number k will be amplified at a rate $\gamma_k(\rho, \varepsilon)$ that depends on the strength of the gradient term: γ_k^2 = $|v_s|^2 k^2 - a^2(\varepsilon_s/h)(\rho/\rho_s)^2 k^4$. Thus, the gradient term introduces a penalty for the development of short-range undulations, and γ_k exhibits a maximum at the favored length scale; the familiar dispersion relation for ideal fluid dynamics, $\gamma_k = |v_s|k$, is recovered in the absence of a gradient term, $a = 0$.

FIG. 2 (color online). Effect of the gradient term: For various values of the parameter a in Eq. ([1\)](#page-1-1) is shown the interface tension σ_T as a function of T (upper panel) and the growth rate γ_k as a function of the wave number k of a disturbance, as obtained with ideal fluid dynamics (lower panel).

These spinodal growth rates have been extracted by following the fluid-dynamical evolution of small harmonic perturbations in uniform matter. Figure [2](#page-2-0) shows the resulting γ_k curves extracted at the phase point (6 ρ_s , 10 ε_s) located in the central spinodal region (see Fig. [1\)](#page-1-0). Because the numerical algorithm inevitably produces some degree of dissipation [[17\]](#page-4-13), the growth rates extracted for $a = 0$ deviate from the exact ideal dispersion relation $\gamma_k = |v_s|k$. Our studies suggest that, while those results are reliable only up to $k \leq 4$ fm, the use of a finite-range $a > 0$ extends validity to significantly finer scales.

Having thus shown that our model produces both a meaningful interface tension and reasonable spinodal growth rates, we apply it to central collisions of lead nuclei bombarded onto a stationary lead target at various kinetic energies E_{lab} . For each energy, an ensemble of 200 separate evolutions is generated, each starting from a different initial state obtained by means of the ultrarelativistic quantum molecular dynamics model (in its cascade mode) [\[18](#page-4-14)[–20\]](#page-4-15) in order to take account of the nonequilibrium dynamics in the very early stage of the collision. The switch from ultrarelativistic quantum molecular dynamics to fluid dynamics occurs at a time t_0 when the two Lorentz-contracted nuclei have just passed through each other and all initial collisions have occurred. The needed fluid-dynamical quantities $\rho(r)$, $\varepsilon(r)$, and $u(r)$ are then extracted from the ultrarelativistic quantum molecular dynamics state by representing each hadron by a Gaussian of 1 fm width and mapping this information onto a Cartesian spatial lattice [\[21\]](#page-4-16). Thus the effects of stopping as well as local event-by-event fluctuations are naturally included in the ensemble of the initial states.

Figure [2](#page-2-0) brings out the intimate relationship between the interface tension σ_T and the spinodal growth rates γ_k , and the range parameter a merely presents a convenient means for linking the two properties. For the present Letter, we have adjusted a so that the zero-temperature interface tension is approximately 10 MeV/fm², the value that was also preferred in Ref. [[13\]](#page-4-12). This corresponds to the middle curves in Fig. [2](#page-2-0), obtained for $a = 0.033$ fm.

The key effect of the first-order phase transition is the spinodal amplification of spatial irregularities [\[5](#page-4-3),[6](#page-4-9),[8,](#page-4-4)[15\]](#page-4-10). In order to obtain a quantitative measure of the resulting degree of clumping in the system, we extract the moments of the (net) baryon density distribution $\rho(r)$:

$$
\langle \rho^N \rangle \equiv \frac{1}{A} \int \rho(\mathbf{r})^N \rho(\mathbf{r}) d^3 \mathbf{r},\tag{3}
$$

where $A = \int \rho(r) d^3r$ is the total (net) baryon number. Thus the corresponding normalized moment $\langle \rho^N \rangle / \langle \rho \rangle^N$ is unity for $N = 1$. These quantities have observational relevance, because they are intimately related to the relative production yield of composite baryons.

The time evolution of the normalized density moments is illustrated in Fig. [3.](#page-2-1) Generally, for a given density distribution $\rho(r)$, the normalized moments increase with the order N, as one would expect. At the initial time t_0 , the

FIG. 3 (color online). Mean time evolution of normalized density moments obtained for $E_{\text{lab}} = 3$ A GeV with $a = 0.033$ fm (for which $\sigma_0 \approx 10 \text{ MeV}/\text{fm}^2$), using either the two-phase equation of state (solid lines) or its one-phase partner (dashed lines).

pre-equilibrium fluctuations are transcribed into the fluiddynamical functions (see above), and they may then be amplified subsequently by the instabilities associated with the first-order phase transition. For comparison, we show also the corresponding results for the one-phase partner equation of state that has no instabilities. There is a striking difference between the two sets of results: Whereas the onephase equation of state hardly produces any amplification at all, the one having a phase transition leads to significant enhancements that increase progressively with N and amount to well over an order of magnitude for $N = 7$. This clearly demonstrates that the first-order phase transition may have a qualitative influence on the dynamical evolution of the density. Because of the variations in the initial conditions, the phase regions explored differ from one collision to the other. As a result, some of the evolutions experience amplifications that are significantly larger than the ensemble average (by up to a factor of 5 or so), whereas more evolutions are affected considerably less.

Figure [3](#page-2-1) depicts what would happen if the expansion were to continue while maintaining local equilibrium so the generated density enhancements eventually subside. However, as the hadronic gas grows more dilute, local equilibrium cannot be maintained. Consequently, if the decoupling occurs sufficiently soon after the clumps are formed, the associated phase-space correlations may survive. Studies of this issue are under way.

The degree of density clumping generated during a collision depends on how long the bulk of the matter is exposed to the spinodal instabilities. The optimal situation occurs for beam energies that produce maximum bulk compressions lying well inside the unstable phase region, because the instabilities may then act for the longest time $[8,15]$ $[8,15]$ $[8,15]$. At lower energies, an ever smaller part of the system reaches instability and the resulting enhancements are smaller. Conversely, at higher energies, the maximum compression occurs beyond the spinodal phase region and the system is exposed to the instabilities only during a relatively brief period during the subsequent expansion. For still higher energies, the spinodal region is being missed entirely.

Figure [4](#page-3-0) shows the (ensemble average) maximum enhancement achieved as a function of the beam energy for the two equations of state. The existence of an optimal collision energy is clearly brought out. While the presently employed equation of state suggests that this optimal range is $E_{\text{lab}} \approx 2-4$ A GeV, it should be recognized that others may lead to different results. Furthermore, the magnitude of the effect depends on the degree of fluctuation in the initial conditions which in turn are governed by the preequilibrium dynamics. On the other hand, our studies suggest that the optimal energy is rather insensitive to the range parameter a.

In summary, by augmenting an existing fluid-dynamical code with a gradient term, we have obtained a transport model that is suitable for simulating nuclear collisions in

FIG. 4 (color online). Mean maximum enhancement of the normalized density moments for $N = 7$ and 5 (top and bottom, respectively) as obtained for various energies by using either the two-phase equation of state (solid lines) or its one-phase Maxwell partner (dashed lines).

the presence of a first-order phase transition: It describes both the tension between coexisting phases and the dynamics of the unstable spinodal modes. Applying this novel model to lead-lead collisions, we have found that the associated instabilities may cause significant amplification of initial density irregularities. Because such clumping is expected to facilitate the formation of composite particles, such as deuterons and tritons, this effect may be experimentally observable.

However, the magnitude of the generated clumping depends considerably on the largely unknown specifics of the equation of state which determines whether the unstable region of the phase diagram is entered during the collision, for how long the system remains there, and how rapidly irregularities are amplified. It would therefore be interesting to expand the present kind of study to other equations of state and also to ascertain the importance of dissipative effects [[15](#page-4-10),[22\]](#page-4-17).

Furthermore, standard fluid dynamics propagates the system deterministically and thus ignores the effect of the ever present thermal noise. Because the initial state, in the present study, is already endowed with significant fluctuations, the thermal noise is likely to play only a minor role (see Ref. [\[5\]](#page-4-3)), but it would be interesting to investigate this quantitatively. (A study of the effect of noise in chiral fluid dynamics was recently made [[23](#page-4-18)].)

Spinodal decomposition is an inherent characteristic of first-order phase transitions, and this nonequilibrium phenomenon may therefore provide especially compelling signals. The present study demonstrates that the phase structure does affect the character of the density evolution, and we hope that these results will stimulate efforts to develop analysis techniques for extracting the related observables from experimental data.

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