

Viscous potential flow analysis of peripheral heavy ion collisions

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The conditions for the development of a Kelvin-Helmholtz instability (KHI) for the quark-gluon plasma (QGP) flow in a peripheral heavy ion collision is investigated. The projectile and target side particles are separated by an energetically motivated hypothetical surface, characterized with a phenomenological surface tension. In such a view, a classical potential flow approximation is considered and the onset of the KHI is studied. The growth rate of the instability is computed as a function of phenomenological parameters characteristic for the QGP fluid: viscosity, surface tension, and flow layer thickness.

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I. INTRODUCTION

The first models of high-energy heavy ion collisions in the 1970s were successful by assuming highly idealized shock fronts where the matter was heated up and compressed in a front (having a discontinuity in perfect fluid flow [1,2]). This led to high-pressured, shock-compressed domains that collectively deflected the incoming nuclear fluid. The observation of this directed flow (side splash of bounce-off) was the first proof of the collective fluid dynamical behavior of nuclear matter [3]. Recent theoretical developments and experimental observation of high multiplicity fluctuations indicate that the quark-gluon plasma (QGP) is a low-viscosity fluid, which makes turbulent phenomena possible [4–6].

Here we adopt again a fluid dynamical picture and discuss the strong shear flow arising in the initial states of peripheral heavy ion collisions at ultrarelativistic energies, which may lead to KHI under favorable conditions, as discussed recently [4].

A simple analytic study showing the development of the KHI in a highly idealized situation is discussed. The investigated phenomenon has some resemblance to the initial state of a peripheral heavy ion collision. In these collisions the collective flow should be a “shear flow” because the top participant layers move nearly with projectile velocity while the bottom layers with the target velocity.

In the reaction plane, the height of the participant profile is $L = 2R - b$, where b is the impact parameter, and the half height is $l = L/2 = (2R - b)/2$. In the following we denote by n the nuclear matter density and by η its phenomenological viscosity. The matter coming from different sources is marked by subscript/superscript t (top) and b (bottom); see Fig. 1.

For simplicity, we consider a two-dimensional nonrelativistic dynamics in the reaction plane. The position vector is $\mathbf{x} = (x, z)$, and the velocity vector is $\mathbf{v} = (v_x, v_z)$.

The nearly perfect QGP provides a possibility of a strong idealization in this situation. The velocity profiles for v_z presented on Fig. 2 illustrates (the dotted line) that the KHI develops if we have a strong shear flow at the $x = 0$ plane, which leads to large vorticity and circulation. With decreasing viscosity we could idealize this configuration in a way that the vorticity is constrained into a narrow layer around the $x = 0$ plane, while the circulation remains constant. In a limiting

case if we constrain the vorticity and shear to the dividing, $x = 0$, plane, this plane would represent an infinite vorticity, providing the same circulation for a trajectory surrounding the dividing plane.¹ In this limiting case the velocity profile is idealized to the one indicated by the full line in Fig. 2. Thus, in the top and bottom domain the flow has no shear and can be described as potential flow, which is an important simplification and idealization. This makes the analytic study of the KHI possible.

II. THE ANALYTIC MODEL

Thus, the idealized dividing layer represents a discontinuity of the flow velocity (i.e., unconstrained slip). At the same time a small viscosity would contribute the transverse momentum transfer to a small distance across the dividing front. The particles in this narrow layer, scattering over from the other side of the dividing plane, would have a high relative velocity, so this layer would exhibit an extra energy increase compared to the general fluid body on the top or the bottom side of our system. This can be taken into account as an effective surface energy of the dividing layer. This surface energy can be estimated both from a microscopic kinetic theory approach and from a rough energy balance calculation. In a microscopic approach one would assume that the this extra energy depends on the (viscosity-dependent) thickness of the layer and the (temperature-dependent) rate of transverse flow crossing the dividing plane. A quantitative estimate of this surface energy in this idealized situation is not feasible, but its existence and a rather qualitative estimate can be made. In contrast with this, a phenomenological energy balance calculation is easier to perform. The advantage of such an approach would be that one does not rely on further estimates for the involved physical parameters. Here, we use this later method to approximate the surface tension of the dividing layer.

¹The conservation of circulation occurs in classical, barotropic flow. In QGP the temperature dominates the pressure change, so the circulation is not conserved but decreases during the expansion of the system [9].

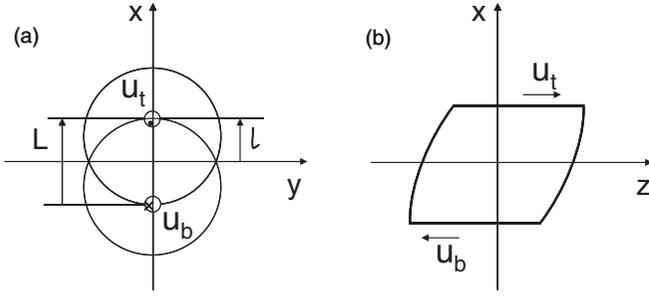


FIG. 1. Sketch of a collision. Panel (a) is a view in the transverse, $[x, y]$ plane; panel (b) is an illustration in the reaction, $[x, z]$ plane. The almond shape in the middle of figure (a) is the participant zone of the event. Right after the collision, streaks are formed and the top streaks move along the z direction while bottom ones move along the $-z$ direction. Owing to this velocity shear, an instability wave will appear on the interface plane between the top and the bottom sheets.

Following Ref. [7] we idealize the problem and assume an initial state where the shear is localized at the dividing plane between the top (t) half and the bottom (b) half of the fluids, in order that we can use the potential flow description in the top and bottom parts of the fluid; see Fig. 3. We assume that the fluid in the top and bottom parts are allowed to slip at the top and bottom boundaries, as well as at the dividing surface between them. We reference these as unconstrained slip conditions. The initial flow velocity is assumed to be uniform in the two layers, so that for the top layer $\mathbf{v}^t = (0, U_t)$ for $0 < x < l$ and for the bottom layer $\mathbf{v}^b = (0, U_b)$ for $-l < x < 0$ initially. This means that initially the amplitude of the wavelike instability is extremely small, and we are looking for the conditions to have a growing amplitude for this instability. For the sake of the analytic model we assume that the density is constant. Numerical studies [4] show that this constraint can be relaxed.

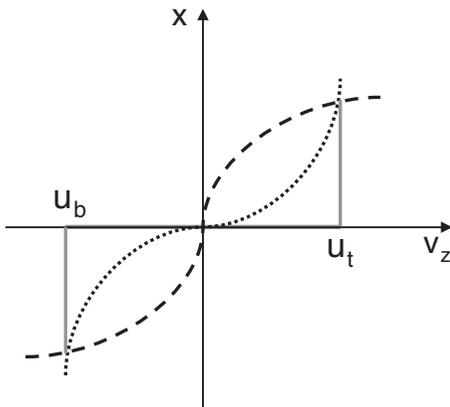


FIG. 2. The velocity along the z axis is represented by the dotted curve, calculated in our computational fluid dynamics (CFD) model and presented in Ref. [4]. The full line with two singularities in its derivative is the case used in Ref. [7]. Here we have to mention that the velocity profile of the dotted curve will induce the KHI effect, while the velocity profile illustrated with the dashed curve will not; see Chapter 8 of Ref. [8].

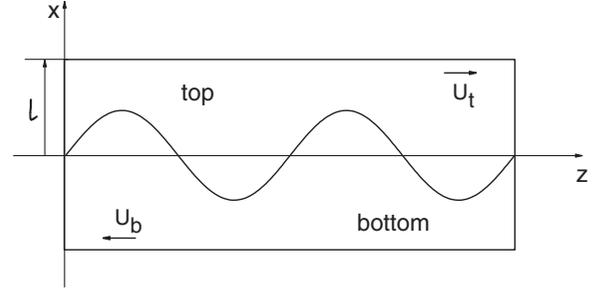


FIG. 3. The profile of the top and bottom fluid layers with a dividing surface wave on it. The top external fluid moves with velocity U_t along the z direction, while the bottom fluid moves with velocity U_b along the $-z$ direction. The slip is unconstrained at the dividing surface.

Just as in Ref. [4] we use an average energy and mass density for our estimates. Because the effective particle density is constant the continuity equation of the flow velocity, \mathbf{v} , will become

$$\nabla \cdot \mathbf{v} = 0. \quad (1)$$

For the top and bottom parts of the fluid we assume small velocities and neglect velocity gradients and we assume that the rotation of the flow velocity is $\nabla \times \mathbf{v} = 0$. Under such conditions we can describe the flow as a potential flow, i.e., $\mathbf{v} = \nabla \phi$, where ϕ is the velocity potential, $\phi \equiv \phi(t, x, z)$. The continuity equation is $\Delta \phi \equiv \nabla^2 \phi = 0$; applying this for the top and bottom layers gives

$$\nabla^2 \phi_t = 0 \quad \text{for } 0 < x < l, \quad (2)$$

$$\nabla^2 \phi_b = 0 \quad \text{for } -l < x < 0. \quad (3)$$

We assume that, owing to the raising instability, the initially plane interface will experience a perturbation and will deviate from the $x = 0$ plane. The height of the deviation from the $x = 0$ plane is denoted by $h = h(t, z)$, and it is taken as wavelike perturbation in the z direction with wave number k . We also allow the amplitude at a given coordinate to change in time. The most general form for such a wavelike interface would be

$$h(t, z) = A_0 e^{(\sigma t + i k z)}, \quad (4)$$

where, σ is the complex growth rate and A_0 is a complex amplitude.

Consequently, the fluid on the top and bottom sides next to the dividing surface will have vertical velocity components,

$$v_x^{t,b} = \frac{dh}{dt} = \frac{\partial h}{\partial t} + U_{t,b} \frac{\partial h}{\partial z}, \quad (5)$$

where U_t and U_b are the flow velocities of the fluid at the top and bottom of the bounding surfaces, respectively. These are assumed to be the average velocity of the fluids at the surface neglecting the horizontal velocity fluctuations arising from wave formation. Initially, these velocity fluctuations are small, and our aim is to study the initial development of KHI. At the $h(t, z)$ dividing surface we also assume unconstrained slip conditions of the two inversely flowing fluid slabs.

The boundary conditions for the external border of the profile are

$$v_x^{t,b} = \frac{\partial \phi_{t,b}}{\partial x} = 0 \quad \text{at } x = \pm l. \quad (6)$$

Initially (at time $t = 0$), one would also have to satisfy

$$v_z^{t,b}(t = 0, x = \pm l) = U_{t,b}. \quad (7)$$

At each time moment the potential ϕ_t and ϕ_b is satisfying Eqs. (2), (3), (5), and (6) for the top and bottom sides, respectively:

$$\begin{aligned} \frac{\partial^2 \phi_{t,b}}{\partial x^2} + \frac{\partial^2 \phi_{t,b}}{\partial z^2} &= 0, \\ v_x^{t,b} &= \frac{dh}{dt} = \frac{\partial h}{\partial t} + U_{t,b} \frac{\partial h}{\partial z} \quad \text{at } x = 0, \\ v_x^{t,b} &= \frac{\partial \phi_{t,b}}{\partial x} = 0 \quad \text{at } x = \pm l. \end{aligned} \quad (8)$$

Assuming for the interface, $h(t, z)$, in Eq. (4), a wavelike perturbation, which is symmetric in $\pm x$ and exponentially decreasing away from the surface, the solution can be searched in the form

$$\phi_{t,b} = A_{t,b} \cosh[k(x-l)]e^{(\sigma t + ikz)} + zU_{t,b}, \quad (9)$$

where A_t , A_b , and A_0 are the complex amplitudes, σ is the growth rate, and k is the wave number. From the kinematic conditions on the dividing layer Eq. (5), we get the following equations at the dividing surface:

$$(\sigma + ikU_{t,b})A_0 = \mp kA_{t,b} \sinh(kl). \quad (10)$$

The pressure (p), viscosity (η), and surface tension (γ) balance at the interface writes as

$$-p_t + 2\eta \frac{\partial v_x^t}{\partial x} - \left(-p_b + 2\eta \frac{\partial v_x^b}{\partial x} \right) = -\gamma \frac{\partial^2 h}{\partial z^2}. \quad (11)$$

The surface energy and consequently the surface tension of the dividing layer is approximated later. As it was already emphasized in the introductory paragraphs, although the top (t) and bottom (b) sides are of the same nuclear matter, the velocity jump or the sharp velocity change contribute to additional surface energy owing to the large shear at the interface exhibiting extra energy or to a smaller extent by the momentum dependance of nuclear interaction potential.

Because we have unconstrained slip conditions on the dividing surface between the top and bottom layer, p_t and p_b can be written by the classical equation of motion without the viscous term as

$$\rho \left(\frac{\partial v_z^{t,b}}{\partial t} + U_{t,b} \frac{\partial v_z^{t,b}}{\partial z} \right) = -\frac{\partial p_{t,b}}{\partial z}. \quad (12)$$

Then, first we apply ∇_z on both sides of the equation and substitute the equation of continuity, $\partial_z v_z = -\partial_x v_x$, into it:

$$\rho \left(\frac{\partial^2 v_x^{t,b}}{\partial t \partial x} + U_{t,b} \frac{\partial^2 v_x^{t,b}}{\partial x \partial z} \right) = \frac{\partial^2 p_{t,b}}{\partial z^2}. \quad (13)$$

Here ρ is the effective mass density of the QGP, we use $\rho = 10 \text{ GeV/fm}^3 c^2$ [4] in our work. To substitute the above equations into Eq. (11), we consider the second-order derivative of Eq. (11) as a function of z and substitute Eq. (13) into

it. Thus, the pressure, viscosity, and surface tension balance will be written in the following form:

$$\begin{aligned} -\rho \left(\frac{\partial^2 v_x^t}{\partial t \partial x} + U_t \frac{\partial^2 v_x^t}{\partial x \partial z} \right) + 2\eta \frac{\partial^3 v_x^t}{\partial x \partial z^2} \\ + \rho \left(\frac{\partial^2 v_x^b}{\partial t \partial x} + U_b \frac{\partial^2 v_x^b}{\partial x \partial z} \right) - 2\eta \frac{\partial^3 v_x^b}{\partial x \partial z^2} = -\gamma \frac{\partial^4 h}{\partial z^4}. \end{aligned} \quad (14)$$

By inserting the velocity derived from Eq. (9) and the considered interface profile, Eq. (4), into the above equation and expressing the top and bottom amplitudes, $A_{t,b}$ from Eq. (10), after simplifying all over with A_0 and putting the condition $x = 0$, we obtain an equation for σ and k :

$$\begin{aligned} [\rho(\sigma + ikU_t)^2 + 2\eta k^2(\sigma + ikU_t)] \coth(kl) \\ + [\rho(\sigma + ikU_b)^2 + 2\eta k^2(\sigma + ikU_b)] \coth(kl) + \gamma k^3 = 0. \end{aligned} \quad (15)$$

Considering this as an equation for σ , one can write it in a simplified form as

$$A\sigma^2 + 2B\sigma + C = 0, \quad (16)$$

where the coefficients, A , B , C are defined as

$$\begin{aligned} A &= 2\rho \coth(kl), \\ B &= 2k^2\eta \coth(kl) + ik\rho(U_b + U_t) \coth(kl) = B_R + iB_I, \\ C &= -k^2\rho \coth(kl)(U_t^2 + U_b^2) + \gamma k^3 \\ &\quad + 2ik^3\eta \coth(kl)(U_t + U_b) = C_R + iC_I. \end{aligned} \quad (17)$$

The solution is

$$\begin{aligned} \sigma &= -\frac{B}{A} \pm \sqrt{\frac{B^2}{A^2} - \frac{C}{A}} \\ &\rightarrow \sigma_R + i\sigma_I = -\frac{B_R + iB_I}{A} \pm \frac{\sqrt{D}}{A}, \end{aligned} \quad (18)$$

where $D = D_R + iD_I$ and

$$\begin{aligned} D_R &= k^2\rho^2 \coth^2(kl)(U_t - U_b)^2 \\ &\quad + 4\eta^2 k^4 \coth^2(kl) - 2\rho \coth(kl)\gamma k^3, \\ D_I &= 0; \end{aligned} \quad (19)$$

thus, the real part and the imaginary part can be expressed as

$$\sigma_R = \frac{-B_R \pm \sqrt{D_R}}{A}, \quad \sigma_I = -\frac{B_I}{A}. \quad (20)$$

In heavy ion collisions, the matter will expand after the collision, and, in fact, there is no external boundary (top and bottom) of the fluid shown in Fig. 3. If we assume $l \rightarrow \infty$, the above equations can be simplified as

$$\sigma_R = -\frac{k^2\eta}{\rho} \pm \sqrt{\frac{k^4\eta^2}{\rho^2} + \frac{k^2(U_t - U_b)^2}{4} - \frac{\gamma k^3}{2\rho}}, \quad (21)$$

$$\sigma_I = -\frac{k(U_t + U_b)}{2}. \quad (22)$$

In Eq. (21), for the typical parameters of a peripheral heavy ion collision, the real part of the growth rate, σ_R , is dominantly dependent on the viscosity η , namely the first term and the first term in the square root. In our expanding system the dominant wave number of KHI is changing with time.

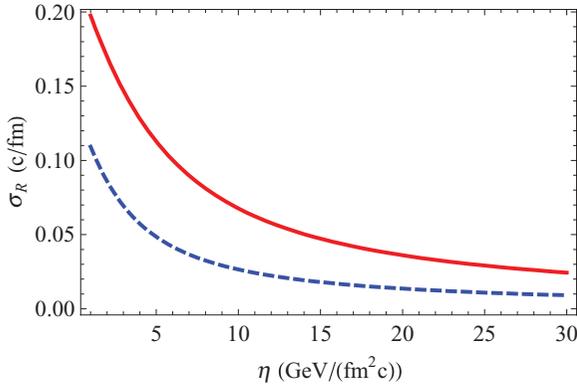


FIG. 4. (Color online) The real part of the growth rate, σ_R , is shown as function of the viscosity η . The full (red) line is for the surface tension $\gamma = 0.4$ GeV/fm² and the dashed (blue) line is for $\gamma = 3.5$ GeV/fm². The wave number, k , is taken to be $k = 0.6$ fm⁻¹ and the effective mass density is $\rho = 10$ GeV/fm² c². The growth rate decreases when the viscosity increases, suggesting that the KHI grows weaker for a more viscous fluid.

III. RESULTS

According to the CFD observations [4], initially we have a small wave formation with $k \approx 1$ fm⁻¹, but with time and expansion, the possible largest wavelength takes over with $k \approx 0.6$ fm⁻¹, which decreases further with the expansion of the system. By assuming $|U_t - U_b| = 0.8c$, we can obtain the growth-rate dependence of the viscosity, η , and the surface tension, γ , which are shown in Figs. 4 and 5.

Similarly to the viscosity, the effective surface energy also influences the growth rate of KHI. As expected, larger surface tension or surface energy damps the growth of KHI. Beyond a critical surface energy (in our model at $\gamma_{\text{crit}} \approx 5.3$ GeV/fm²) the surface tension will lead to a decrease in the KHI. Interestingly, this threshold value is independent of the

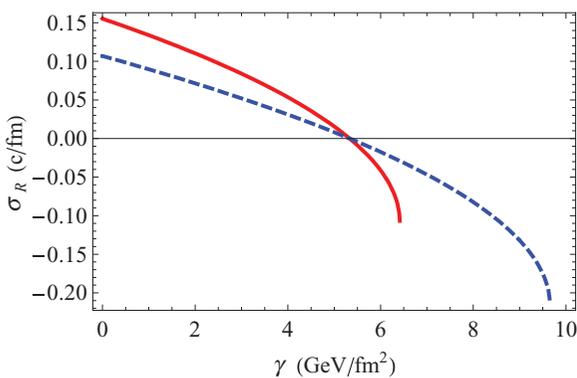


FIG. 5. (Color online) The real part of the growth rate, σ_R , as a function of the surface tension γ with different values of viscosity η . The full (red) line represents $\eta = 3$ GeV/fm² and the dashed (blue) line represents $\eta = 6$ GeV/fm². The wave number k is taken as 0.6 fm⁻¹ and the effective mass density ρ is 10 GeV/fm² c². As we can see in the figure, the two curves cross each other at $\sigma_R = 0$, which is around $\gamma = 5.3$ GeV/fm², and then the growth rate becomes negative. With bigger surface tension the KHI effect is less probable to appear.

viscosity. This is part of the general feature that the behavior of the zero-growth ($\sigma_R = 0$) curve is independent of the value of the viscosity in this model. The growth rate and damping rate are, of course, dependent on the viscosity.

The condition to have a growing instability is to have a solution with $\sigma_R > 0$. Taking into account that D is a real number ($D_I = 0$, $D_R > 0$), and $B_R > 0$, from Eq. (20) it follows that to have a positive growth rate ($\sigma_R > 0$) one has to satisfy the condition

$$\sqrt{D_R} > B_R. \quad (23)$$

Thus, using Eqs. (17) and (19) we get the condition for positive growth:

$$V^2 > \frac{2\gamma k}{\rho \coth(kl)}, \quad (24)$$

where $V \equiv U_t - U_b$.

The above condition will limit the region of the (V, k) parameter space where the KHI can evolve. One should also keep in mind the results obtained in Ref. [4], regarding the acceptable wave numbers, k , for the considered wavelike instability. Definitely there is a lower cutoff (k_{min}) governed by the beam-directed longitudinal length of the flow, l_z :

$$k_{\text{min}} = \frac{2\pi}{l_z}. \quad (25)$$

For the $b = 0.5b_{\text{max}}$ and $b = 0.7b_{\text{max}}$ impact parameter values the calculations in Ref. [4] leads to $k_{\text{min}} = 0.598$ fm⁻¹ and $k_{\text{min}} = 0.479$ fm⁻¹ values, respectively. There is also an upper limit for the wave numbers, k_{max} governed by the Kolmogorov length scale, λ_K :

$$k_{\text{max}} = \frac{2\pi}{\lambda_K}. \quad (26)$$

According to Ref. [4], this characteristic length scale is estimated for the above-given impact parameters as $\lambda_K \approx 3.5$ fm and $\lambda_K \approx 2.5$ fm, leading to $k_{\text{max}} = 1.79$ fm⁻¹ and $k_{\text{max}} = 2.51$ fm⁻¹ values, respectively.

For the peripheral Pb + Pb collisions, the radius of Pb is $R = 7$ fm; thus, $b_{\text{max}} = 14$ fm. To get the parameter space where the KHI will grow, let us estimate now the value of the surface tension. As it has been discussed in the introductory part, this surface energy comes from the energy excess of the unbalanced energy flow in the two layers. Although a theory based on kinetic considerations would capture more from the involved physics, here we just consider a simple approach based on the energy balance. The reason for doing this is that fewer phenomenological parameters are needed.

The flow assumed in the present work has a perpendicular velocity profile illustrated in Fig. 6(a). This means that a smooth velocity profile [Fig. 6(b)], characterizing a stable and balanced viscous flow, is not formed. In the case illustrated in Fig. 6(a) one would assume that there are two distinct layers flowing with velocities U_t and U_b . For the balanced flow illustrated in Fig. 6(b), one would observe a smooth flow velocity transition from U_t to U_b . It is obvious that in the laboratory frame, this later flow has less kinetic energy in the z direction than the previous one. The difference between the two kinetic energies can be accounted as the energy surplus

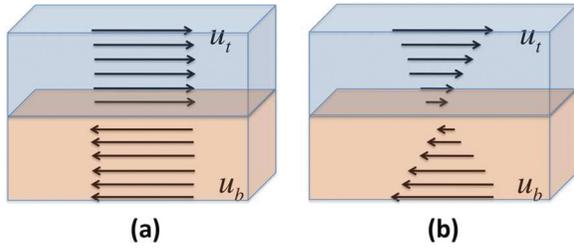


FIG. 6. (Color online) Two different velocity profiles. Panel (a) is the profile used in our present work; it has two distinct layers with two constant velocity U_t and U_b . Panel (b) has a flow transition from the two layers, and at the dividing surface the velocity is smallest.

of the dividing layer. If we denote the contact surface between the flows in the top and bottom layers by S , the surface tension could be estimated as

$$\gamma = \frac{E_{kz}^a - E_{kz}^b}{S}, \quad (27)$$

where $E_{kz}^{a,b}$ denotes the kinetic energy of the flow in the z direction for the profile illustrated in Figs. 6(a) and 6(b), respectively. The total relativistic kinetic energy of the system, E_k , in the laboratory frame is

$$E_k = 2M_{Pb}c^2 \left(\frac{1}{\sqrt{1 - \frac{V^2}{c^2}}} - 1 \right), \quad (28)$$

where $V = U_t - U_b$ is the relative speed of the two projectiles, and M_{Pb} is the mass of the collided Pb ions. Assuming that the participating zone in the collision has a surface $q \times \pi R^2$ (the overlapping regions are only a q part of the possible ones) and the kinetic energy of the participating particles in this zone is distributed equally in all the directions of the space, a rough approximation for E_{kz}^a would be $E_{kz}^a = q \frac{E_k}{3}$. However, for the flow illustrated in Fig. 6(b), owing to the balanced velocity profile, a part of this kinetic energy has to be dissipated, and assuming a linear velocity profile, one gets $E_{kz}^b = 1/2 E_{kz}^a$. The above arguments lead us to a first approximation of the surface tension value:

$$\gamma = \frac{q}{3} \frac{M_{Pb}c^2}{S} \left(\frac{1}{\sqrt{1 - \frac{V^2}{c^2}}} - 1 \right). \quad (29)$$

Assuming $q \approx 0.5$ and estimating the surface of the dividing layer, S , from Ref. [4], one gets the values of γ for different impact parameter values.

The surface tension is estimated to be $\gamma = 0.4 \text{ GeV/fm}^2$ from Eq. (29). This value is used in the following examples. The critical velocity Eq. (24) for different impact parameters is shown in Fig. 7. These curves show the border of instability of the growth rate, σ_R . The curves divide the space into two areas; the upper side above the curve is the region where the instability grows and the area below the critical velocity curve is where the instability does not grow. The KHI development region is also limited by the k_{\min} and k_{\max} values as drawn in figure Fig. 7.

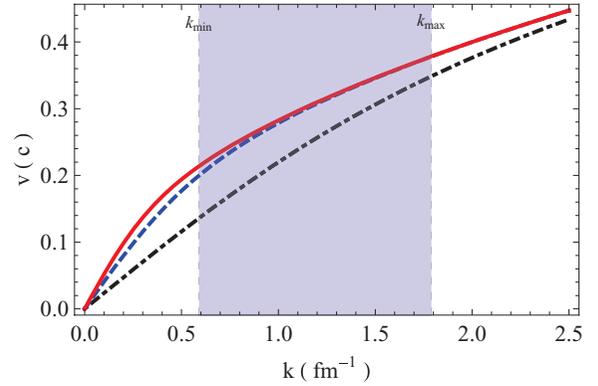


FIG. 7. (Color online) The critical velocity $V \equiv U_t - U_b$ as function of the wave number, at the condition of vanishing growth rate, $\sigma_R = 0$. The red full line, blue dashed line, and black dot-dashed line are for impact parameters $b = 0.5b_{\max}$, $0.7b_{\max}$, $0.9b_{\max}$, respectively. On the graph we also illustrated the two natural boundaries k_{\min} and k_{\max} for $b = 0.5b_{\max}$. The KHI will evolve thus above the critical velocity curves and between these two limits. For increasing impact parameters, the instability is less able to grow and the system tends to be stable.

The above consideration is for $\sigma_R = 0$; however, this does not show the η dependence of the growth. To see how the instability depends on the viscosity, η , we can cast Eq. (20) into the form

$$\sigma_R = \frac{k^2 \eta}{\rho} \left[-1 \pm \sqrt{1 + \frac{\rho}{\eta^2} \left(V^2 \rho - \frac{\gamma k}{\coth(kl)} \right)} \right]. \quad (30)$$

This suggests that with our characteristic parameters the dependence on the thickness of the fluid layer, l , is weak, as shown in Fig. 8.

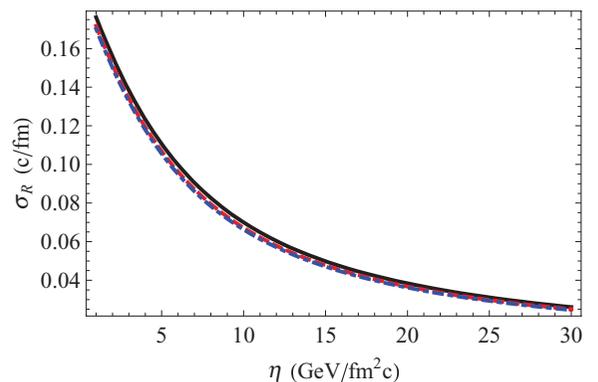


FIG. 8. (Color online) The growth rate σ_R as a function of the viscosity η at different values of l ; the black line is for $l = 0.1 \text{ fm}$, the red dashed line is for $l = 2 \text{ fm}$, and the blue dot-dashed line is for $l = \infty$. The wave number is $k = 0.6 \text{ fm}^{-1}$, the surface tension is $\gamma = 0.4 \text{ GeV/fm}^2$, the relative velocity is $V = 0.8c$, and $\rho = 10 \text{ GeV/fm}^2 c^2$. The growth rate depends weakly on l , while it depends significantly on the viscosity, increasing strongly for small viscosity values.

IV. CONCLUSIONS

In classical gravitational water waves the wave formation and wave speed depends strongly on the depth of the water, i.e., the layer thickness, l . In heavy ion collisions the role of the layer thickness is different. The material properties of the top and bottom layers are not different, these are separated from each other by the relatively thin layer of large shear. Still, the occurrence of KHI in such conditions is not uncommon, because it is frequently observed as turbulence during airplane flights, or it is even visible if the air has high humidity and condensation makes the KHI visible.

In peripheral heavy ion collisions the layer thickness is given in the initial state, but there is no solid boundary and the system expands in all directions. Thus, for this physical situation the large or infinite layer thickness is more relevant in this model, even if the initial layer thickness is finite and usually smaller than the longitudinal size of the initial state.

Large viscosity or the corresponding low Reynolds number prevents the development of turbulence and KHI, so that these phenomena appear only above a critical Reynolds number. This critical Reynolds number depends on the flow configuration, so it is separately analyzed for the KHI also; see Ref. [4]. The present study confirms that the dependence of the growth rate on the viscosity reflects the usual tendency that instability and turbulence increases with smaller viscosity.

When the KHI develops between two fluids (e.g., air/water or air/oil) the large surface tension difference at the interface damps the development of the instability; this is well known for sailors for centuries. If KHI develops inside one fluid, like in air or in quark gluon fluid, there is no surface tension in the classical sense, but the layer with large shear has extra energy, and it leads to an effective surface tension, which hinders the development of KHI.

We presented a strongly idealized analytic model for the development of the Kelvin-Helmholtz instability in ultrarelativistic heavy ion reactions. We compressed the shear zone into a central infinitesimal layer, following the idea of Ref. [7], and assumed that the remaining flow can be approximated as potential flow. The idealized dividing layer was attributed to a surface energy and unconstrained slip between the top and bottom fluid layers. It is interesting that in this model the KHI is developing under similar conditions, as in numerical high-resolution relativistic fluid dynamical calculations [4]. This model also shows that critical size KHI may occur for low-viscosity QGP.

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