The Granular Phase Diagram

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Abstract

The kinetic energy distribution function satisfying the Boltzmann equation is studied analytically and numerically for a system of inelastic hard spheres in the case of binary collisions. Analytically, this function is shown to have a similarity form in the simple cases of uniform or steady-state flows. This determines the region of validity of hydrodynamic description. The latter is used to construct the phase diagram of granular systems, and discriminate between clustering instability and inelastic collapse. The molecular dynamics results support analytical results, but also exhibit a novel fluctuational breakdown of mean-field descriptions.

Key words: granular Dynamics

Granular media such as sand provide an attractive opportunity to revisit a number of topics in classical physics, and contribute new angles. In this paper we describe the granular phase diagram which we hope will be helpful to a broad community given the raising interest in granular systems [1]. The researchers interested in diluted granular gases, such as in astrophysical applications [2], and researchers who study, say, compaction of sand [3] use different approaches. The phase diagram may represent a ground for communication.

The phase of granular system depends on inelasticity of collisions, $r$, (restitution coefficient approximation, [4, 5, 6]), particle density, $\rho$, particle size, $a$, system size,
and the observation time, \( t \). The external influence of shaking, gravity, boundaries enters through the above parameters. Granular temperature, or any other characteristic of the rate of bulk-averaged motion doesn’t appear on the phase diagram since there is no characteristic energy scale for hard-core interaction assumed here.

As we shall see below, a combination, \( \rho L a^{d-1} \), where \( d \) is the system dimension, is playing the key role. It represents the average number of particles inside an imaginary tube of length \( L \) and cross-section \( a^{d-1} \). Presently we think that the phase diagram consists of at least three regions, Fig. 1. In the region \( 1 - r \ll (\rho L a^{d-1})^{-2} \) the system is gas-like. In the region \( (\rho L a^{d-1})^{-2} \ll 1 - r \ll (\rho L a^{d-1})^{-1} \) the system is condensed but doesn’t collapse, and for \( (\rho L a^{d-1})^{-1} \ll 1 - r \) the system contains inelastically collapsed chains of particles, mixed with non-collapsing regions (see [7] for the description of collapse). Fig. 2 is a snapshot from an event-driven Molecular Dynamics (MD) with 5000 particles in a circle the wall of which is maintained at a constant temperature. Particles undergo inelastic collisions with a constant coefficient of restitution (see Eqs. (3,4) below). When a particle hits the wall it experiences diffusive angular scattering, while the distribution of the scattered velocity amplitude is Maxwellian, with the temperature being that of the wall. After a period of equilibration the system finds a state drawn in Fig. 2, which is in many respects a steady state. In Fig. 2 the gray scale linearly codes the relative number of collisions experienced by each particle during the previous \( 10^5 \) time steps, where black means high collision rate. One can clearly distinguish three different regions: (i) close to the wall there is a gas-like phase with low density where the mean free path is large. (ii) The region between the center of the bulk and the wall consists of closely packed particles. We consider this region as condensed phase. It is well separated from the gas like phase. (iii) Close to the center of the container we find a region which isn’t distinguishable from the second region by “naked eye”. With the help of gray scale coding one observes collision chains, i.e. almost linearly arranged chains of particles which made major contribution to the previous \( 10^5 \) collisions, in comparison to their neighbors. Quantitatively, about 5 percent of the particles participated in about 96 percent of the collisions. Sometimes we observed that very few (\( \approx 5 \ldots 10 \)) particles
take almost all of the collisions in a certain time interval (not shown). The lengths
of the appearing chains as well as their life times vary irregularly, their statistics will
be discussed in detail elsewhere [8]. This figures serves as an illustration that in MD
simulation one encounters all three regions of the phase diagram given in Fig. 1.

The role of observation time is not included here, and will be discussed separately
below. Notwithstanding this classification, the condensed phases may be ordered in
space (a crystal) or disordered (a glass). This different structure-based classification
exists in the elastic limit. We now present the arguments used for constructing Fig. 1
beginning from the gas-like phase.

As the parameter \( \rho La^{d-1} \) increases the first condensation or clustering transition
of granular gas occurs. The above given criterion has been obtained by Goldhirsch
and Zanetti [8] with the usage of granular hydrodynamics [4, 5]. This description is
based on the assumption of molecular chaos [9] and Maxwellian distribution functions
for particle velocities [8] which have been questioned and related to the concept of
inelastic collapse [10]. In this paper we present the limits of validity of granular
hydrodynamics, and then use the hydrodynamical clustering instability to argue that
it occurs before the condition of inelastic collapse is satisfied. The latter is the line sep-
rarating collapsing and non-collapsing condensed phases, \( 1 - r \sim (\rho La^{d-1})^{-1} \), [7, 10].

Inelastic collapse, when a chain of particles experiences an infinite number of
collisions in finite time as discussed by McNamara and Young [7], is realized inside
dense clusters, specified by granular hydrodynamics. Inside such condensed regions we
have numerically observed collision chains. The relation of hydrodynamic instabilities
of collision chains to their collapse and the importance of the upper line in Fig. 1 will
be published elsewhere [8].

We begin with diluted phase. The study of granular gases is greatly simplified by
the fact that the binary collisions dominate and goes back to the works of Boltzmann
and others [11, 12]. It is unclear a priori, whether inelasticity represents a regular
or singular perturbation. The first part of this problem is the reduction of Boltz-
mann equation to hydrodynamics, it can be studied independently from the known
complications associated with hydrodynamical description, i.e. divergence of trans-
The Boltzmann equation for inelastically colliding identical particles reads \[ \frac{\partial f}{\partial t} + \vec{v} \frac{\partial f}{\partial x} = \hat{C}(f, f) \] (1)

where \( f(\vec{v}, x, t) \) is the velocity distribution function, and \( \hat{C} \) is the bilinear collision operator \[ \hat{C}(f, f) = \int \left\{ \frac{1}{r^2} f_1 f_1' - ff \right\} |\vec{v} - \vec{v}_1| d\sigma d\vec{v}_1. \] (2)

The differential cross-section of two spheres of radius \( a \) is \( d\sigma = \pi a^2 \sin^2 \theta d\theta / 2 \), where \( \theta \) is the angle between the vectors \( \vec{q} \) and \( \vec{v} - \vec{v}_1 \). An incoming collision event to the state \( \{\vec{v}, \vec{v}_1\} \) occurs between particles with velocities \( \vec{v}' \) and \( \vec{v}_1' \),

\[
\begin{align*}
\vec{v}' &= \vec{v} - \frac{1 + r}{2r} \vec{q}(\vec{v} - \vec{v}_1) \\
\vec{v}_1' &= \vec{v}_1 + \frac{1 + r}{2r} \vec{q}(\vec{v} - \vec{v}_1).
\end{align*}
\] (3)

The problem is (hopefully!) isotropic and homogeneous and the distribution function, \( f_0 \), must depend only on the absolute value of particle velocity, \( v \), i.e., on kinetic energy, \( E = \frac{mv^2}{2} \). When the initial distribution is forgotten one may search for a similarity solution \( f_0(E, t) = \frac{n}{T(t)} \phi \left( \frac{E}{T(t)} \right) \), (5)

where \( T(t) \) is a single scale of the kinetic energy. The normalization condition to the number density of particles \( n \) is \( \int dz \phi(z) = 1 \). Eqs. (1, 2) result in an integro-
differential equation for function $\phi(z)$

$$-A(r) \left( \frac{d\phi}{dz} + z \frac{d\phi}{dz} \right) = \sqrt{2} \int d\sigma d\alpha \sqrt{g(z)}$$

$$\times \left\{ \frac{1}{r^2} \phi' - \phi \right\} \left( z - z_1 - \sqrt{zz_1} \cos \alpha \right)^{1/2},$$

and in an equation for $T(t)$,

$$dT/dt = -A(r)T^{3/2}. \quad (7)$$

Here $A(r)$ is the constant introduced by separation of variables, $\alpha$ is the angle between vectors $\vec{v}$ and $\vec{v}_1$, and $g(z)$ is the density of states. Examination of the collision integral allows to get asymptotic form of $\phi(z)$ at large $z$. Namely, it can be shown that the incoming term contains extra factor $O(z^{-1/2})$ [this is contributed by the events $(\vec{q}, \vec{v}) \propto z^{-1/2}$] and is small as compared to the outgoing term. The asymptotic form of Eq. (6) is $-A(r) z (d\phi/dz) \sim \phi z^{1/2}$, up to a number. Therefore,

$$\ln \phi(z) \sim \sqrt{z}/A(r). \quad (8)$$

The moments of this function converge, and one can restrict oneself to the hydrodynamical description (7) of the granular temperature which gives

$$T(t) = T_0 \left[ 1 + \frac{t}{2} A(r) \sqrt{T_0} \right]^{-2}, \quad (9)$$

and $T(t) \propto t^{-2}$ at large $t$, (3). One can get this power law from dimensional arguments. $A(r)$ remains undetermined, it can only be found from the full solution of Eq. (3). Asymptotically, $A(r) \propto (1 - r)$ when $r$ is close to 1. The enhanced population of large energies, (8), dominates at $z \gg 1/A^2(r)$. For lower energies the distribution is Maxwellian.

Two sets of MD results by an event-driven code are obtained with 5000 (20000) particles of unit radius in a circle with radius 130 (260). In both cases the surface fraction covered by particles is 50/169. Units of mass, length and time are arbitrary, initial velocity distribution is taken to be uniform in velocity in the square $-1 \leq v_x, v_y \leq 1$. The employed model of collision is the same as given above, $r = 0.999$. Fig. 3 shows the temporal evolution of mean kinetic energy and the number of collisions occurred.
The fit to the energy decay over the entire range is achieved with the help of Eq. (9).

The distribution function is self-similar over four orders of energy decay as is seen in Figs. 3, 4. The latter figure shows the energy distribution function with the argument rescaled by \( \langle E(t) \rangle = T \) for different times. One can see that the hydrodynamical description is quite precise. The system remains uniform and gas-like.

However, at times \( t \approx 10^4 - 10^5 \) we observed appearance of a rarefied space in the center of the circle followed by “condensation” of granular gas on the wall. This is a novel type of transition, which is different from clustering [6] or collapse [7]. One can understand the growth of fluctuations using the following argument. The mean free path \( l \) is given by \( l = 1/\eta = a/c \) with the volume fraction \( c = na^d \) in a \( d \) dimensional system with concentration \( n \). The diffusivity of particles decreases in time like \( D(t) \sim (\delta v) l \sim l \sqrt{T} \), where \( T \) is given in energy units. At large \( t \) (Eq. (9)),

\[
T = \frac{4}{(tA)^2} \sim \left[ \frac{l}{t(1-r)} \right]^2 , \tag{10}
\]

Therefore,

\[
D \sim \frac{l^2}{t(1-r)} = \frac{(a/c)^2}{t(1-r)} , \tag{11}
\]

and the diffusion length of particles in time \( t \) is

\[
l_D \sim \sqrt{\int D(t) \, dt} \sim \frac{(a/c)^2}{1-r} \ln^{1/2} \left( \frac{t \, v_0}{a} \right) , \tag{12}
\]

where \( v_0 \) is the typical initial velocity. The number of particles inside the sphere of radius \( l_D \) is \( N \sim nt_D^d = c (l_D / a)^d \). These particles preserve a fluctuational collective velocity \( v_0 \sqrt{c (l_D / a)^d} \). Equating this velocity to thermal velocity \( \delta v \sim \sqrt{T(t)} \sim (a/c) / (t(1-r)) \) yields the time when the collective velocity exceeds the fluctuational individual velocity \( \delta v \)

\[
t_c \sim \frac{a}{v_0} \frac{c}{(1-r)^{d/4}} \ln^{-d/4} \left( \frac{c}{1-r} \right) , \tag{13}
\]

given with logarithmic precision. Granular flow becomes fluctuationally supersonic at \( t \gg t_c \) and molecular chaos assumption along with the Boltzmann equation and granular hydrodynamics are no longer applicable. The r.h.s. of Eq(13) can be evaluated, and the corresponding time is \( 5.0 \times 10^4 \). Given that the numerical prefactor
in the estimate (13) isn’t known, one finds a very satisfactory agreement between the estimate (13) and the time where hydrodynamical prediction deviates from the MD time-trace of kinetic energy density in Fig. 3. At time $t_c$ groups of particles of the size $l_c \sim \left[ \int_{t_c}^t dt D(t) \right]^{1/2}$ become special. Their mutual encounters lead to structures containing travelling shock waves. The difference between the systems with 5000 and 20000 particles as seen in Fig. 3 is due to the fact that the regions of the size $l_c$ occupy a different fraction of the area, and in the second case it takes longer for inhomogeneities to evolve from the size $l_c$ up to the system size. The observed fluctuational transition is the reason why observation time may explicitly appear in the phase diagram.

We now add a constant energy supply from the border, and allow the system to reach a steady state. The system is no longer uniform in space, and we assume that the corresponding length scales exceed the mean free path, $l$, everywhere. The similarity Ansatz Eq. (5) with a non-uniform temperature $T(x)$ can be used again. A study of the asymptotic form of the collision term shows that in the regions of the phase space where the detailed balance is absent the incoming term is again small in $O(z^{-1/2})$ times as compared to the outgoing term, and no global balance is possible. If $1 - r \ll 1$, the detailed balance approximately holds for the energies, $z \ll (1 - r)^{-1}$, above this range there are effectively no particles. Therefore, we find that the distribution function is almost Maxwellian for the energies less than $T/(1-r)$ and small otherwise,

$$\phi(z) = \begin{cases} 
\exp(-z) & z \ll (1 - r)^{-1} \\
0 & z \gg (1 - r)^{-1} 
\end{cases} \tag{14}$$

All the moments converge, and one can justify the hydrodynamical reduction, which reads

$$\nabla P(\rho, T) = 0 \tag{15}$$

$$\nabla q - B(r)T/\tau_c(\rho, T) = 0 \tag{16}$$

where $P$ is the granular pressure, and $q$ is the thermal flux. This system can be analyzed. To understand what happens as one increases inelasticity (or the number of particles) it is useful to consider the dilute limit approximation when $q = \lambda(T)\nabla T$,
\[ P = \rho T, \]  
where \( \lambda(T) = \gamma_1 T^{1/2} \) is the thermal conductivity, \( B(r) \propto (1 - r) \) is the energy fraction lost per collision. \( \tau_c^{-1}(n, T) = \gamma_2 \rho T^{1/2} \) is the time between collisions. We found that the hydrodynamical solution exists only below a threshold,

\[
a^{d-1}B^{1/2}(r) \int_0^L \rho \, dx < \xi_d(\gamma_1/\gamma_2) \ln(L/a), \tag{17}
\]

\( \xi_d \) is a pure number. Our hydrodynamical analysis is valid when \( \text{(17)} \) is fulfilled; in the opposite case a particle condensate appears, and system has more than one phase. The left-hand side of \( \text{(17)} \) is the generalization of the parameter \( \rho La^{d-1} \) for non-uniform case. Its region of applicability is more general than the derivation given above. Eq \( \text{(17)} \) describes with logarithmic precision the same clustering transition as discussed by Goldhirsch and Zanetti \[6\].

Returning to the question of validity of hydrodynamical reduction of the Boltzmann equation we note that the complete system of equations for granular hydrodynamics \[17, 4, 16\] for density, linear and angular velocity density and granular temperature(s) has a shortcoming. It follows from Eqs. \( \text{(8,14)} \) that in sufficiently inelastic complex flows where changes occur in space and time the granular temperature cannot be introduced. Indeed, if we restrict ourselves with the first moment of the energy distribution in situations when this distribution changes its functional form, we wouldn’t be able to close the hydrodynamical reduction. Therefore, in general, the full system of hydrodynamical equations \[17, 4\] cannot serve as a quantitative description. The situation is a bit easier, though, than in the kinetics of phonons: only kinetic coefficients are not exact, and the uncertainty is in numerical prefactors, which depend on local distribution function. To give an example about difficulties with phonons it is sufficient to recall that the heat transfer may be nonlocal \[18\].

Sufficiently inelastic problems require a mixed kinetic-hydrodynamical description. In this case, similar to kinetics of phonons \[18\], one finds a reduced kinetic equation for the isotropic part, \( f_0 \), of the distribution function

\[
\partial_t f_0 - \nabla \vec{J}[f_0] = \hat{C}(f_0, f_0), \tag{18}
\]
where diffusion-like flux $\vec{J}$ is defined as

$$\vec{J}(f) = \frac{4E}{9m} \left[ \frac{\delta \hat{C}}{\delta f} \right]^{-1} \nabla f. \quad (19)$$

The inverse integral operator $\left[ \frac{\delta \hat{C}}{\delta f} \right]^{-1}$ is fixed by particle number constraint, i.e. the result of applying this operator to $f_0$ must have vanishing norm. Similar expressions arise for all dissipative coefficients. Despite the cumbersome appearance, Eq. (18) together with the remaining hydrodynamic equations for density and momenta conservations offers a reduction from nine to five-dimensional space of arguments.

Our study of inelastic gas with binary collisions provided an opportunity to justify granular hydrodynamics for simple flows and for all flows in dilute and sufficiently elastic systems. This allows to discriminate between clustering instability discussed by Goldhirsch and Zanetti \cite{6} and inelastic collapse \cite{7}. In complex inelastic flows the predictions of granular hydrodynamics are valid on the order of magnitude almost everywhere, and such precision is comparable to that in Fig. \cite{6}. On the basis of this study we constructed the granular phase diagram and identified a novel type of supersonic fluctuational phase transition, leading to shock waves, which is different from clustering and collapse.

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Figure 1: The granular phase diagram. See text for details.
Figure 2: A snapshot from an event-driven Molecular Dynamics simulation of 5000 hard spheres. The wall is kept at a fixed temperature. The gray scale codes the relative number of collisions experienced by particles during the previous $10^5$ time steps. One distinguishes three regions: a gas-like state with low density, and two high density regions, with and without collision chains (c.f. Fig. [3]).
Figure 3: The kinetic energy $E$ per particle over time $t$ from MD simulations. Energy decay curves for systems with 5000 and 20000 particles are identical for a very long time (solid lines) The dashed line shows the prediction of Eq. [8] with $T_0 = 0.333$ and $A = 0.01$. The dash-dotted lines display the cumulative number of collisions per particle scaled by factors $10^{-8}$ (5000) and $4 \times 10^{-8}$ (20000).
Figure 4: Evolution of the scaled energy distribution function in MD simulations. Different symbols used for different times as shown by insets. Statistically, the distribution is indistinguishable from Maxwellian within the given range.