Local equation of state and velocity distributions of a driven granular gas

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We present event-driven simulations of a granular gas of inelastic hard disks with incomplete normal restitution in two dimensions between vibrating walls (without gravity). We measure hydrodynamic quantities such as the stress tensor, density and temperature profiles, as well as velocity distributions. Relating the local pressure to the local temperature and local density, we construct a local constitutive equation. For strong inelasticities the local constitutive relation depends on global system parameters, like the volume fraction and the aspect ratio. For moderate inelasticities the constitutive relation of state, even though the system is highly inhomogeneous with heterogeneous temperature and density profiles arising as a consequence of energy injection. With respect to local velocity distributions we find that they do not scale with the square root of the local granular temperature. Moreover the high-velocity tails are different for the distribution of the *x* and the *y* components of the velocity, and even depend on the position in the sample, the global volume fraction, and the coefficient of restitution.

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I. INTRODUCTION

The physics of vibrofluidized granular materials is far from being fully understood. In particular, the applicability of hydrodynamics [1-5] is still an object of debate [6-8]. Despite similarities to the hydrodynamics of elastic hardsphere systems, with respect to, e.g., the appearance of instabilities [9–12], a main difference from ordinary fluids is the fact that continuous energy injection is vital to maintain a stationary state. Otherwise a gas of inelastic spheres would collapse, even in the absence of gravity. The nontrivial nature of this stationary state has been elucidated in experiments of vibrating grains. Among the most striking features found are non-Gaussian velocity distributions [13-16] and cluster formation [17-19]. Furthermore, several attempts have been made to extract either an equation of state or more generally a scaling relation for the thermodynamic variables [20,21].

Specification of the driving mechanism is a crucial ingredient for any model that describes vibrated granular fluids. A rather simple approach, although less appealing from an experimental point of view, utilizes stochastic bulk heating by uncorrelated random forces, which act on every particle at every instant of time [22-25]. In Ref. [26] random restitution coefficients were considered with a probability distribution allowing values both smaller and greater than one. However, this yields nonuniversal properties depending on the specific form of this distribution. The multiplicative bulk driving defined in terms of stochastic collision rules in Ref. [27] is similar in spirit. Compared to bulk driving, a driving mechanism which acts only at a boundary of the system is much closer to experiments. In Ref. [2] the energy influx at the boundary has been modeled by a heuristically motivated ansatz for the heat current, while Refs. [28,29] assume heating at the boundary through thermal walls. In the present paper the driving mechanism consists of incrementing a particle's velocity by a constant amount after every collision with a wall. This way of driving was used before in, e.g., Refs. [30,31], and with slight modifications in Refs. [32-34].

One theoretical approach by which to study granular systems at fluid densities rests on kinetic theory, see, e.g., Refs. [5,35–37], and references therein. Particular applications to driven granular gases can be found in Refs. [23,26–28,38]. Much of this work has been based on the Boltzmann or Boltzmann-Enskog equation, modified for inelastic collisions. One method by which to solve this nonlinear equation is based on the local equilibrium distribution [39], which is only known for the elastic case. For systems with strongly inelastic collisions, the stationary state is not known so a systematic discussion of transport properties within kinetic theory is severely hampered.

Hydrodynamic studies [2,11,12,30] have been motivated partly by the search for understanding of temperature and density profiles [2], but also by experiments on hydrodynamic-like instabilities [9]. In the hydrodynamic approach the question of how to relate the pressure to the density and temperature arises. Several equations of state for the *global* quantities have been proposed, which go beyond the ideal-gas limit. In Ref. [2] an interpolation between the low and high density limits was constructed. On the other hand, in Refs. [31,39] the Boltzmann-Enskog equation was used to determine the collisional contribution to the pressure. It is not clear whether an equation of state also holds for the *local* hydrodynamic fields in a strongly driven, nonequilibrium system, where the fields are strongly *inhomogeneous*.

Most simulations of vibrated granular gases have been based either on event-driven molecular dynamics [31,40,41] or the direct-simulation Monte Carlo method [28,30,42,43]. If the system is driven through the boundaries, inhomogeneous density and temperature profiles are measured. For low densities the computed temperature profiles agree well with hydrodynamic theory [2,30], whereas for moderate or high densities the profiles are not well understood with the exception of almost elastically colliding particles. The full stress tensor, including potential contributions, has only been computed for freely cooling systems [44]. In simulations of driven granular gases [28] the collisional part of the stress tensor has not been measured, and in Ref. [31] the stress tensor is not measured directly but instead computed from a local equation of state.

Inspired by experiment [16], a lot of emphasis has been put on the tails of the velocity distribution functions, which were found to be overpopulated compared to a Maxwellian. In fact, all intermediate types of decay between a Gaussian and an exponential were observed [31,42,45,46]. In addition, mixtures [31,47,48] and rough spheres [49] have been investigated as well as hydrodynamic instabilities such as convection [50,51] and pattern formation [52].

In this paper, we present results from event-driven simulations of inelastic spheres in two dimensions confined between two vibrating walls without gravity. Our focus is on the stationary state, which is reached when dissipation by particle collisions is equal to the energy injection due to the vibrating walls. In the stationary state, density and temperature profiles are shown to be strongly inhomogeneous due to the driving walls—even in a range of parameters where clustering is only a minor effect. This has led us to:

(a) derive a constitutive equation by relating the measured hydrodynamic fields, granular temperature T(x), volume fraction $\phi(x)$, and pressure p(x), at each point *x*, and

(b) check whether this constitutive equation is universal or depends on the global system parameters of the model, like the aspect ratio of the cell, the overall volume fraction, or the coefficient of restitution of the disks. For moderately inelastic systems (coefficient of restitution α =0.9) the constitutive equation is (almost) independent of the remaining global system parameters. In that case the constitutive equation can be interpreted as a local equation of state for a driven granular gas in the stationary state, even though the system is highly inhomogeneous with heterogeneous temperature and density profiles. In contrast, for strongly inelastic systems (α =0.5) the constitutive equation depends significantly on the global system parameters so the concept of a local equation of state cannot be sustained in this case.

We furthermore discuss the one-particle distribution function in the stationary state and show that:

(c) the local distribution $f_x(x, v_x)$ of v_x , the velocity in the direction of driving, is not a function of the rescaled variable $v_x/\sqrt{T_x(x)}$ alone. Similarly, curves of $f_y(x, v_y)$ cannot be mapped onto a master curve for different x, when plotted against $v_y/\sqrt{T_y(x)}$. Here $T_i(x)$ denotes the local granular temperature associated with the translational motion in the *i* direction. We find deviations from scaling at small and large arguments.

(d) The local velocity distributions f_x and f_y have high-velocity tails whose decay ranges from stretched exponential to almost Gaussian. The particular type of decay depends on the position in the sample, the overall particle density, and the coefficient of restitution. Furthermore, the decay of f_x for large velocities is generally different from that of f_y .

Taken together, one has to acknowledge that the stationary state of a driven granular gas is by no means universal. It shows peculiar features that depend on the precise values of the system parameters in contrast to the corresponding elastic system.



FIG. 1. (Color online) Model of N disks, driven in the x direction with periodic boundary conditions in the y direction.

The outline of the paper is as follows. In Sec. II we introduce the model, specify the driving, and define the observables. We briefly discuss balance of energy input through the walls and energy dissipation in binary collisions in Sec. III. Subsequently we present data for the profiles of the density, the temperature, and the components of the stress tensor (Sec. IV). In Sec. V we relate the local density, pressure, and temperature to derive "experimentally" an equation of state and check its universality. Finally, in Sec. VI we discuss velocity distribution functions and their scaling behavior and present conclusions in Sec. VII.

II. MODEL AND OBSERVABLES

We investigate a driven granular gas in two dimensions consisting of N identical inelastic smooth hard disks of diameter a and mass m which are confined to a rectangular box with edges of length L_x and L_y . The gas is driven through the walls perpendicular to the x direction, which vibrate in an idealized sawtooth manner (see below), while periodic boundary conditions are imposed in the y direction. Figure 1 shows a typical snapshot of this. The gas evolves over time through ballistic center-of-mass motion, binary inelastic collisions, and particle–wall collisions.

A. Binary collisions

The inelastic nature of interparticle collisions is the most important characteristic of granular media. As is often done we assume that it can be taken into account by a constant coefficient of normal restitution and briefly recall the collision rules for binary collisions, see, e.g., Refs. [53,54].

The unit vector from the center of sphere two to the center of sphere one is denoted by $\hat{\boldsymbol{n}} = (\boldsymbol{r}_1 - \boldsymbol{r}_2)/|\boldsymbol{r}_1 - \boldsymbol{r}_2|$ where \boldsymbol{r}_i is the position vector of the center of mass of particle *i*. The center-of-mass velocities before a collision are denoted by \boldsymbol{v}_1 and \boldsymbol{v}_2 , and the relative velocity by $\boldsymbol{v}_{12} = \boldsymbol{v}_1 - \boldsymbol{v}_2$. A prime denotes postcollisional quantities. The relative velocity after a collision is assumed to obey

$$\hat{\boldsymbol{n}} \cdot \boldsymbol{v}_{12}' = -\alpha \hat{\boldsymbol{n}} \cdot \boldsymbol{v}_{12}, \tag{1}$$

where $\alpha \in [0, 1]$ is the (constant) coefficient of normal restitution. The value $\alpha = 1$ describes elastic collisions with energy conservation, while for $\alpha < 1$ energy will not be conserved but instead decreased in each collision.

Constitutive equation (1) plus momentum conservation determines the postcollisional velocities for the disks,

$$\boldsymbol{v}_1' = \boldsymbol{v}_1 + \Delta \boldsymbol{v}_{\rm pp}, \quad \boldsymbol{v}_2' = \boldsymbol{v}_2 - \Delta \boldsymbol{v}_{\rm pp}, \quad (2)$$

where

$$\Delta \boldsymbol{v}_{\rm pp} = -\frac{(1+\alpha)}{2} (\hat{\boldsymbol{n}} \cdot \boldsymbol{v}_{12}) \hat{\boldsymbol{n}}.$$
(3)

The model could easily be extended to rough spheres by also specifying the disks' moment of inertia as well as a coefficient of tangential restitution β_0 and Coulomb friction μ [55,56], but here we restrict ourselves to smooth spheres so that rotational motion is decoupled from translational motion.

B. Driving

When a particle collides with a driving wall, energy is injected into the system. This can be modeled in different ways, for example, by drawing a new velocity from a Maxwellian distribution of a given wall temperature [2] or by assuming that the wall has a coefficient of normal restitution that is greater than 1. Both of these mechanisms have no close experimental equivalents though. A more realistic model is to assume a vibrating wall moving either in a symmetric (e.g., sinusoidal) or in an asymmetric (e.g., sawtooth) way. In addition, this can be combined with a normal (and also tangential) coefficient of restitution [33,41,57]. In this paper we refrain from the latter and restrict ourselves to sawtooth driving of the walls within the limit of vanishing amplitude A and diverging frequency v such that $Av = v_{drive}/2$ is a constant. This ensures that the driving walls are always located at the same positions and leads to the following simple expression for a particle's change of velocity due to collision with the left/right wall:

where

$$\Delta \boldsymbol{v}_{\rm pw} = (-2\boldsymbol{v}_x \pm \boldsymbol{v}_{\rm drive})\boldsymbol{e}_x,\tag{4}$$

and e_x stands for the unit vector in the x direction.

C. Observables from simulations

 $\boldsymbol{v}' = \boldsymbol{v} + \Delta \boldsymbol{v}_{\mathrm{pw}},$

Following, e.g., Refs. [58,59] we performed event-driven simulations in two dimensions with periodic boundary conditions in the *y* direction and two identical idealized vibrating elastic walls in the *x* direction. We initialize the system by placing the disks on a triangular lattice with a Gaussian velocity distribution. To let the correlations of the initial state relax, the system is evolved elastically with periodic boundary conditions in the *y* direction and elastic nonvibrating walls in the *x* direction for an average of 100 collisions per particle. Then we switch on driving of the left and right walls and dissipation for particle–particle collisions. Before we start measuring the observables we let the system relax further until, at time t_0 , it has reached a stationary state, indicated by the total kinetic energy, which fluctuates around a time-independent mean value.

To measure hydrodynamic fields in the stationary state we subdivide our box into cells V_r of area $|V_r|$, centered at position $\mathbf{r} = (x, y)$. We then count the number of particles $f(\mathbf{r}, v_x, v_y, t) |V_r| dv_x dv_y$ at time t in cell V_r with an x component of the velocity between v_x and $v_x + dv_x$ and y component of the velocity between v_y and $v_y + dv_y$. This local observable fluctuates as a function of time. To eliminate these fluctuations we average over a (long) time interval of length τ and compute

$$f_{\text{stat}}(\boldsymbol{r}, \boldsymbol{v}_{x}, \boldsymbol{v}_{y}) = \frac{1}{\tau} \int_{t_{0}}^{t_{0}+\tau} dt \, f(\boldsymbol{r}, \boldsymbol{v}_{x}, \boldsymbol{v}_{y}, t).$$
(5)

Of particular interest is the particle density

$$\rho(\mathbf{r}) = \int_{\mathbb{R}} dv_x \int_{\mathbb{R}} dv_y f_{\text{stat}}(\mathbf{r}, v_x, v_y), \qquad (6)$$

or the area fraction $\phi(\mathbf{r}) = \rho(\mathbf{r}) \pi a^2/4$ and the two components i=x, y of the granular temperature,

$$T_i(\mathbf{r}) = \frac{m}{\rho(\mathbf{r})} \int_{\mathbb{R}} dv_x \int_{\mathbb{R}} dv_y f_{\text{stat}}(\mathbf{r}, v_x, v_y) (v_i - U_i(\mathbf{r}))^2, \quad (7)$$

where

$$\boldsymbol{U}(\boldsymbol{r}) = \frac{1}{\rho(\boldsymbol{r})} \int_{\mathbf{R}} d\boldsymbol{v}_x \int_{\mathbf{R}} d\boldsymbol{v}_y f_{\text{stat}}(\boldsymbol{r}, \boldsymbol{v}_x, \boldsymbol{v}_y) \boldsymbol{v}$$
(8)

is the velocity field. The total granular temperature in two dimensions is defined as $T(\mathbf{r}) = [T_x(\mathbf{r}) + T_y(\mathbf{r})]/2$.

The stress tensor at position r and time t has a kinetic contribution and one that is due to interactions between the particles $\sigma(r,t) = \sigma^{kin}(r,t) + \sigma^{int}(r,t)$. The kinetic part is given by

$$\sigma_{ij}^{\text{kin}}(\mathbf{r}) = -m \int_{\mathbb{R}} dv_x \int_{\mathbb{R}} dv_y f_{\text{stat}}(\mathbf{r}, v_x, v_y)$$
$$\times [v_i - U_i(\mathbf{r})] [v_j - U_j(\mathbf{r})]. \tag{9}$$

If the particles interact through (finite) forces the contribution due to interactions is given by the correlations of the particles' relative positions and the forces between them. For hard-core interactions there are no forces, and one has to consider momentum transfer in a small time interval instead [59,60]: Suppose there is a collision at time t of particle k with another particle, then the change of momentum of particle k will contribute to the stress tensor an amount proportional to $l_i^k(t)\Delta p_i^k(t)$. Here $l_i^k(t)$ is the *i*th component of the vector of length a/2 pointing from the center of disk k to the contact point of the colliding disks, and $\Delta p_i^k(t)$ is the *j*th component of the change of momentum of particle k during this collision. To compute the collisional part of the stress tensor we need the total change of momentum in the time interval $[t-\Delta t,t]$ in cell V_r so we have to keep track of all collisions *n* occurring at times $t_n \in [t - \Delta t, t]$, for which at least one collision partner k_n (i.e., its center of mass) is located in cell V_r at time t_n [59],

$$\sigma_{ij}^{\text{int}}(\boldsymbol{r},t) = \frac{1}{\Delta t} \frac{1}{|V_{\boldsymbol{r}}|} \sum_{t_n} \sum_{k_n} l_i^{k_n}(t_n) \Delta p_j^{k_n}(t_n).$$
(10)

The particle number k_n of each such collision can take on one or two values, depending on whether one or both collision partners are located in cell V_r . In a second step $\sigma^{int}(r,t)$ is averaged over time, similar to the other hydrodynamic fields, in order to get the collisional part of the stress tensor in the stationary state,

$$\sigma_{ij}^{\text{int}}(\boldsymbol{r}) = \frac{1}{\tau} \int_{t_0}^{t_0 + \tau} dt \; \sigma_{ij}^{\text{int}}(\boldsymbol{r}, t). \tag{11}$$

The corresponding local pressure $p(\mathbf{r}) = -\text{tr}[\boldsymbol{\sigma}(\mathbf{r})]/2$ is defined as usual as the negative trace of the stress tensor divided by the space dimension.

Besides hydrodynamic fields we also measure velocity distributions in Sec. VI. They are readily obtained by integrating out the relevant variables in the stationary-state distribution function f_{stat} . A different method, which is better suited to determining high-velocity tails, for example, will be presented in Sec. VI.

Coarse-grained measurements in space and time of certain observables may depend on the coarse-graining resolution. For example, this was demonstrated for the stress tensor in shear-flow driven granular systems in [60,61]. Our measurements of observables in the stationary state, however, should not suffer from such effects for two reasons: First, we could not detect any significant nonzero local velocity field in the simulated systems and, second, because of the long-time average needed to obtain stationary-state quantities.

Moreover, in the simulations we have never found any significant dependence of the (long-time-averaged) hydrodynamic fields on y, the coordinate parallel to the driving walls. Yet, stripe states, which are homogeneous in y, but have enhanced density in the middle of the sample, are known [10,11] to exhibit instabilities with respect to density fluctuations in y. A marginal stability analysis [11] of granular hydrodynamics (for α close to 1) determines the conditions under which such phenomena occur. As far as a comparison can be made, our systems fall into the stable region. Hence, we choose the cells V_r as stripes along the y direction, for which we write V_x , and compute the hydrodynamic fields with spatial resolution in the x direction only. Thus, we also write $\rho(x)$ instead of $\rho(r)$ and change the notation accordingly for all other quantities. In the simulations the stripes V_r were all chosen to have equal width of $L_r/201$, the temporal resolution of our measurements was set to $\Delta t=1$, and the long-time average involves typically $10^7 - 10^9$ collision events.

Another instability of stripe states, which is related to oscillations of the center dense cluster in the x direction [62], will be addressed at the end of Sec. IV A.

III. DIMENSIONAL ANALYSIS AND BALANCE OF ENERGY

The model system contains three independent length scales: the diameter *a* of a disk and the box sizes, L_x and L_y .

In addition, there is one independent velocity scale, the driving velocity v_{drive} , and one independent mass scale, the mass of a disk *m*. Together with the initial positions and velocities of the particles this exhausts all dimensional quantities entering the time evolution of the system.

We would like to describe the system using dimensionless variables that do not depend on the initial conditions since we expect stationary states to be independent of them. Therefore, we will measure all lengths in units of particle diameter a, all times in units of a/v_{drive} , and all energies in units of $mv_{\rm drive}^2$. Note that there are no other time and energy scales. Thus, we introduce dimensionless variables: box sizes \tilde{L}_x $=L_x/a$ and $\tilde{L}_y=L_y/a$, granular temperatures $\tilde{T}_x=T_x/(mv_{\text{drive}}^2)$ and $\tilde{T}_v = T_v / (m v_{\text{drive}}^2)$, and stress tensor $\tilde{\sigma} = \sigma a^2 / (m v_{\text{drive}}^2)$. In the stationary state all dimensionless variables like T and $\tilde{\sigma}$ are independent of the driving velocity and only depend on position *r* and the remaining four independent dimensionless system parameters, which characterize the system completely: the number of disks N, the two edges of length L_x and L_{y} in units of *a*, and the coefficient of normal restitution α , which is a dimensionless material constant.

For simplicity of notation we refrain from indicating dimensionless quantities by a tilde from now on. This should not cause confusion, because quantities having a physical dimension will no longer be used in the rest of the paper (except in the Appendix).

We are interested in the macroscopic limit, which is taken such that $N \rightarrow \infty$ and $L_y \rightarrow \infty$ with fixed line density $\lambda = N/L_y$ of particles. Thus, in the macroscopic limit the number of system parameters is further reduced to L_x , λ , and α . It is important to keep L_x finite, otherwise balance of energy would not work: energy input occurs only at boundaries, while energy dissipation is a bulk phenomenon.

It is instructive to estimate the average or global granular temperature $\overline{T} = \int_{-L_x/2}^{L_x/2} dx \ \rho(x)T(x)$ by balancing the energy input at the walls and the energy loss due to particle collisions in the bulk [33,63]. As shown in the Appendix, we find that the granular temperature in the stationary state is independent of the initial data and given by

$$\bar{T} = \left(\frac{2}{\pi}\right)^3 \psi^{-2} \left(1 + \sqrt{1 + (\pi/2)^2 \psi}\right)^2, \tag{12}$$

where $\psi = \sqrt{2}\chi\lambda(1-\alpha^2)$ and χ stands for the pair correlation at contact of the corresponding elastic system, which we estimate by the Henderson approximation [Eq. (A6)]. In Fig. 2 we plot the global granular temperature \overline{T} as a function of ψ^{-1} and compare it with simulations. For small values of ψ^{-1} the simulation deviates significantly from this simple theory's prediction. For large values of ψ^{-1} (dilute and/or quasielastic systems) the agreement is reasonable. In addition, we show the result,

$$\bar{T} = \frac{1}{2\pi\psi^2} \left(1 + \sqrt{1 + \psi/2} \right)^2, \tag{13}$$

of a more refined calculation, which uses the pseudo-Liouville-operator approach to kinetic theory that will be



FIG. 2. (Color online) Global granular temperature \overline{T} as a function of the parameter ψ^{-1} defined in the line below Eq. (12). Comparison between simulations and the simple energy-balance argument, Eq. (12), as well as with the refined version, Eq. (13). The inset shows the same graph but on a nonlogarithmic scale.

presented elsewhere. Equation (13) yields better agreement with the data for intermediate values of ψ^{-1} .

It is worth noting that the same type of argument which led to Eq. (12) also predicts that there is no generic stationary state for systems with the multiplicative driving mechanism [Eq. (A1)], described by $v_{drive}=0$ and a coefficient of restitution bigger than 1 for particle–wall collisions. Instead, such systems either cool down or heat up according to Haff's law [64]; see Eq. (A9) in the Appendix. This was also confirmed by simulations (not presented).

IV. HYDRODYNAMIC FIELDS IN SIMULATIONS

Here in Sec. IV we discuss the results of our simulations for the hydrodynamic fields, computed from Eqs. (6)–(11). Due to the absence of a local velocity field U, these equations simplify accordingly. We have performed simulations for a wide range of system parameters, α , ϕ_0 , L_x , L_y , and present examples thereof below. We then go on in Sec. V to discuss the question of a local equation of state that relates p(x), $\rho(x)$, and T(x).

A. Density and temperature profiles

We first present data for the density and temperature to demonstrate that the system is strongly inhomogeneous even for collisions with α =0.9 and moderate densities. The parameters chosen are L_x =20, L_y =25, and N=256 so that the global area fraction is $\phi_0 = \pi a^2 N/(4L_x L_y)$ =0.4. In Fig. 3 we show the local area fraction $\phi(x)$, the x and y components of the granular temperature, $T_x(x)$ and $T_y(x)$, as well as the isotropic temperature $T(x)=[T_x(x)+T_y(x)]/2$. We note that all quantities are symmetric in x as expected. Except for a small area of approximately one disk diameter next to the walls (indicated by the vertical lines), the area fraction is a monotonic function for either half of the system with the maximal value in the middle of the system. The temperatures T_x , T_y ,



FIG. 3. (Color online) Spatial profiles of the area fraction ϕ and the granular temperatures T_x , T_y , and T from simulations (α =0.9, L_x =20, λ =10.24, N=256, ϕ_0 =0.4, corresponding to a mean free path of $\ell/L_x \approx 0.02$).

and T are monotonic as well with the lowest temperatures in the middle of the sample.

The reason for the increased density next to the wall is the effective attractive potential of the wall due to entropic effects: Once a disk gets closer to the wall than one disk diameter, it can only receive hits from within the box but no hits from the direction of the wall. Thus the particle is pushed closer to the wall [65,66]. This effect is partially compensated by the driving walls, which add momentum to any particle hitting the walls.

To support this explanation we have also investigated systems of half the size $-0.5 \le x/L_x \le 0$, half the number of particles, with an elastic wall at x=0. The resulting hydrodynamic fields (not shown) are almost identical to the ones in Fig. 3 except for a small region of about one diameter close to x=0 where the aforementioned effect is particularly visible.

In systems exhibiting stripe states, we have sometimes observed an oscillatory instability of a central dense cluster in the *x* direction [62], particularly if α is low and L_x is large. However, these oscillations occur on far shorter time scales compared to the time interval τ , over which we average to obtain stationary-state quantities. Hence, these oscillations are completely averaged out in the data presented.

B. Stress tensor and pressure

In Figs. 4–6 we show the components of the stress tensor $\sigma_{xx}(x)$, $\sigma_{yy}(x)$, and $\sigma_{xy}(x)$ for two different sets of parameters. Figure 4 is typical of systems with fairly elastic collisions (α =0.9) and moderate densities (ϕ_0 =0.4), whereas Figs. 5 and 6 show rather dilute systems (ϕ_0 =0.15) at two different inelasticities, a moderate one (α =0.9 in Fig. 4) and a strong one (α =0.5). The off-diagonal component σ_{xy} of the stress tensor is always vanishingly small. The *xx* component σ_{xx} is constant within the sample except for at a boundary layer close to the driving walls, whereas Fig. 4 displays a dip



FIG. 4. (Color online) Spatial profiles of the stress tensor components σ_{xx} , σ_{yy} , σ_{xy} , and the product ρT for the system in Fig. 3. For all *x* away from the boundary layers, ρT is hard to distinguish from ρT_x and ρT_y (not shown).

of σ_{yy} in the center of the sample, which is more pronounced for the more dilute system in Fig. 5 and hardly visible in the strongly inelastic, dilute system of Fig. 6. Furthermore, σ_{yy} increases considerably over a broad range in Fig. 6 when moving from a driving wall towards the center of the system. Currently, the origin of both *x* dependences of σ_{yy} is not clear to us. While the stationary-state condition $\nabla \cdot \boldsymbol{\sigma} = 0$ and homogeneity in *y* require σ_{xx} and σ_{xy} to be constant in *x* on general grounds, this is not the case for σ_{yy} . We have carefully checked that the *x* dependence of σ_{yy} is not caused by any shear instability associated with a nonzero velocity field U(x) in the system. Thus, there are normal stresses present in the simulated systems, and they depend on *x*.

If the equation of state of the ideal gas were to hold, then the local pressure p(x) would be related to the local temperature and density according to $p(x)=\rho(x)T(x)$ $=(4/\pi)\phi(x)T(x)$. For our system the kinetic part of the stress tensor is diagonal and each component is simply related to



FIG. 5. (Color online) Same as Fig. 4, but for a less dense system (α =0.9, L_x =50, λ =9.6, N=240, ϕ_0 =0.15, $\ell/L_x \approx$ 0.03).



FIG. 6. (Color online) Same as Fig. 5, but for $\alpha = 0.5$.

the corresponding temperature $\sigma_{xx}^{kin}(x) = \rho(x)T_x(x)$ and $\sigma_{yy}^{kin}(x) = \rho(x)T_y(x)$. Hence the difference between the measured stress tensor and the ideal gas behavior is due to the collisional part of the stress tensor. It can be seen clearly from Figs. 4–6 that collisions contribute significantly to the stress tensor measured. Here they have been measured directly in the simulations and not just estimated by approximate theories as has been done elsewhere [30,31].

We have also estimated the global mean free path ℓ for our simulated systems according to Eq. (9) in Ref. [2], which expresses ℓ solely as a function of the global volume fraction ϕ_0 . For the denser system with $\phi_0=0.4$, shown in Figs. 3 and 4, this gives $\ell \approx 0.5$ (in units of the diameter of the disks, that is, $\ell/L_x \approx 0.025$). For the thinner systems with $\phi_0=0.15$, shown in Figs. 5 and 6, one gets $\ell \approx 1.65$, that is, $\ell/L_x \approx 0.033$. Thus, in all cases the mean free path ℓ is much smaller than the scales governing spatial variations of the hydrodynamic fields, which are of order L_x .

C. Density scaling

For low global area fractions, $\phi_0 \leq 0.01$, we observe scaling of the relative local area fraction $\phi(x)/\phi_0$ when plotted versus x/L_x . This is shown in Fig. 7, where we compare density profiles of systems with the same degree of inelasticity, $\alpha = 0.9$, and the same line density, $\lambda = 10.24$, but different values of the box width $1280 \leq L_x \leq 20\ 000$, corresponding to a global area fraction ϕ_0 between 6×10^{-3} and 4×10^{-4} . The relative local area fraction ϕ/ϕ_0 is seen to be a function of x/L_x only and does not depend separately on L_x and ϕ_0 . If the global area fraction were increased beyond $\phi_0 \approx 0.01$ (not shown), then data collapse would cease to hold. In addition, the height of the peak would be reduced and instead of the bell-shaped master curve in Fig. 7 one would get overall concave profiles, similar to the one in Fig. 3.

The hydrodynamic approach to quasielastic, driven granular gases in Sec. II D in Ref. [2] predicts for low-density systems that the mean free path in units of L_x depends only on the line density according to $\ell/L_x=1/(\sqrt{8\lambda})$. When computed for the situation in Fig. 7, one gets the small numerical



FIG. 7. (Color online) Master curve for rescaled local area fraction in different low-density systems with the same line density (α =0.9, λ =10.24, $\ell/L_x \approx 0.03$). The dotted line corresponds to the theoretical prediction of Ref. [2].

value $\ell/L_x \approx 0.03$, indicating that ℓ is not a relevant length scale for the master curve of the rescaled spatial density profiles. The theory of Ref. [2] also makes a prediction for the relative local area fraction $\phi(x)/\phi_0$ in terms of the solution of a first-order differential equation. This differential equation includes one free parameter, which we have fitted in order to match the solution (dotted line in Fig. 7) at x=0 with the master curve from our simulations. The agreement is reasonable, showing that inelasticities of $\alpha=0.9$ are at the borderline of the scope of this otherwise powerful approach to quasielastic, driven granular gases.

V. LOCAL EQUATION OF STATE

Having measured hydrodynamic fields in the steady state of the system, one is led to search for a relation among them. Such a constitutive equation is needed, for example, in all hydrodynamic approaches to driven granular gases in order to obtain a closed set of equations [2,11,30]. In Sec. IV it was shown that a driven granular gas is intrinsically inhomogeneous. Therefore it is only natural to investigate how the *local* values of the granular temperature T(x), pressure p(x), and density $\rho(x)$ [respective area fraction $\phi(x) = \rho(x)\pi/4$ in our units] are related to each other. To do so we observe in Fig. 3 that T(x), p(x), and $\phi(x)$ are all symmetric in x. Moreover, and this is crucial, $\phi(x)$ is monotone in x for either sign of x (except for a boundary layer of approximately one diameter in width close to a driving wall, which we ignore). Therefore one can invert the function $\phi(x)$ for positive x. Upon inserting $x = \phi^{-1} [\phi(x)]$ into the local pressure and temperature, we arrive at the constitutive equation,

$$\frac{p(x)}{\rho(x)T(x)} = G(\phi(x)), \tag{14}$$

with some function *G*. Figure 8 shows parametric plots on a semilogarithmic scale of the function *G* with the values of $p(x)/[\rho(x)T(x)]$ plotted against those of $\phi(x)$ for all *x* (except



FIG. 8. (Color online) Semilogarithmic parametric plots of function $G(\phi)$ from the local equation of state, Eq. (14), for different systems (α =0.9, L_y =25, 0.01 $\leq \ell/L_x \leq$ 0.05). Also shown are the theoretical predictions from our Eqs. (15) and (16), and Eq. (4) in Ref. [67].

for those in the boundary layer mentioned above). In order to utilize a broad range of ϕ values, Fig. 8 contains data from 11 different systems, all of which have the same coefficient of restitution α =0.9 and the same height L_y =25, but different widths L_x and different global area fractions ϕ_0 . For not too large values of ϕ these data merge quite nicely, indicating that there is only a weak dependence of *G* on the global system parameters L_x and ϕ_0 in the corresponding parameter range. In this case constitutive equation (14) can be interpreted as the local equation of state of the system. The horizontal line in Fig. 8 marks ideal-gas behavior, from which *G* deviates due to the collisional contribution to the pressure. These deviations increase significantly with increases of ϕ .

The dotted line in Fig. 8 corresponds to the function,

$$G(\phi) = 1 + (1 + \alpha)\phi\chi, \tag{15}$$

where χ stands for the pair correlation function at contact of the associated elastic hard-sphere gas in thermal equilibrium [73]. Since χ is not known exactly, we estimate it by the Henderson approximation [Eq. (A6)] for numerical purposes. In the context of granular gases, Eq. (15) occurred already in [39]. In Ref. [68] the pair correlation function was studied in a homogeneously driven inelastic system with periodic boundary conditions. It was found to be nearly independent of the coefficient of restitution and approximated well by the Henderson approximation [Eq. (A6)]. In Ref. [31] this form of G was used in the local equation of state (14) to theoretically predict p(x) from simulations of T(x) and $\phi(x)$ in a driven granular gas. Figure 8 reveals that this works generally quite well for up to rather high local area fractions $\phi(x) \leq 0.5$. In even denser systems agreement still holds for the well fluidized parts. Deviations from Eq. (15) start to occur when entering the transition zones to the frozen-out stripe of particles in the center of these high-density systems. The dash-dotted line in Fig. 8 corresponds to the interpolation formula [2],



FIG. 9. (Color online) Same as Fig. 8, but for more inelastic systems (α =0.5, L_y =25, 0.01 $\leq \ell/L_x \leq$ 0.1).

$$G(\phi) = \frac{\phi_c + \phi}{\phi_c - \phi},\tag{16}$$

for $0 \le \phi < \phi_c$, which connects the behavior of dilute (van der Waals) and dense (ordered) elastic hard-sphere systems. Here $\phi_c = \pi/(2\sqrt{3}) \approx 0.91$ denotes the area fraction for ordered closed packings in two dimensions. Equation (16) was applied to quasielastic granular gases in Ref. [2], and even for our simulations with $\alpha = 0.9$ there is agreement with the local data in the low-density regions up to $\phi(x) \leq 0.4$. In addition, the most dense regions of regularly ordered, frozenout particles in the center of the high-density systems are described correctly, too. Yet another interpolation formula for G, which is rather accurate for an elastic hard-sphere gas even in the vicinity of the freezing transition, was put forth in Eq. (4) in Ref. [67] and is depicted by the dashed line in Fig. 8. The crossover between fluidized and frozen-out behavior in our inelastic, driven systems originates from the transition zone at the border of the solidified stripe of particles in the system center. It is very smooth in comparison to the coexistence region of the freezing transition for an equilibrated elastic gas. Moreover, the location of the crossover depends on the global system parameters. Since both interpolation formulas, Eq. (16) and that from Ref. [67], were tailored for elastic systems in thermal equilibrium, it is surprising to find such good agreement with the local data for driven inelastic hard-sphere gases with restitution $\alpha = 0.9$.

When lowering the coefficient of restitution α from 0.9 to 0.8 (not shown), there are only little changes to the plot in Fig. 8. In particular, the spread of the data pertaining to different systems increases, even for low densities. Still one may be inclined to interpret Eq. (14) as a local equation of state—at least approximately. This situation is intermediate to the previous one with α =0.9 and the following one for strongly inelastic systems with coefficient of restitution α =0.5. Figure 9 reveals major discrepancies in the local pressure which are due to the systems with global area fractions above $\phi_0 \approx 0.2$. The discrepancies occur even at positions in the sample where the local area fractions are below 0.2 and

where all the more dilute systems, i.e., those with $\phi_0 \leq 0.2$, agree reasonably with the proposal, Eq. (15). The concept of a local equation of state is therefore no longer sustainable for such strongly inelastic systems, and Eq. (14) merely plays the role of a local constitutive equation, which depends in addition on the global system parameters.

VI. ABSENCE OF SCALING FOR VELOCITY DISTRIBUTIONS

Finally, we examine the local velocity distributions of the driven granular gas in the stationary state. We distinguish between the local distribution,

$$f_x(x,v_x) = \frac{1}{\rho(x)} \int_{\mathbb{R}} dv_y f_{\text{stat}}(x,v_x,v_y), \qquad (17)$$

at position x of the velocity component v_x in the direction of the driving and the distribution,

$$f_{y}(x,v_{y}) = \frac{1}{\rho(x)} \int_{\mathbb{R}} dv_{x} f_{\text{stat}}(x,v_{x},v_{y}), \qquad (18)$$

of the velocity component v_y perpendicular to the direction of driving. By definition, these velocity distributions are normalized to unity.

In order to determine f_x and f_y from the simulation we use two different methods. The first one extracts them directly according to their definitions (17) and (18) from f_{stat} , which was determined in the manner described in Sec. II C. One disadvantage of this method is that it yields a smeared-out velocity distribution, which is spatially averaged over the width of strip V_x centered around x. This is of practically no importance in the middle of the simulation box, but strongly disturbing for resolving the subtleties which occur close to the driving walls that are presented in Sec. VI A. The second way of measuring the velocity distributions avoids this problem: To get f_i (*i*=x or y) we keep track of all those particles which pass the line parallel to the y axis at position x within a very long time interval of length τ and whose *i*th component of the velocity lies in a small interval of width Δv around v_i . Let us enumerate these particles by n_i and denote their velocity components by $v_x^{(n_i)}$ and $v_y^{(n_i)}$. Then one has (within the limits $\tau \rightarrow \infty$ and $\Delta v \rightarrow 0$)

$$\rho(x)f_i(x,v_i) = \frac{1}{\tau \Delta v L_y} \sum_{n_i} \frac{1}{|v_x^{(n_i)}|},$$
(19)

because for i=x, respectively i=y, the right-hand side of (19) is equal to

$$\int_{\mathbb{R}} dv_y \frac{|F_x(x, v_x, v_y)|}{|v_x|} \quad \text{respectively} \quad \int_{\mathbb{R}} dv_x \frac{|F_x(x, v_x, v_y)|}{|v_x|}.$$
(20)

Here, $F_x(x, v_x, v_y) = v_x f_{\text{stat}}(x, v_x, v_y)$ denotes the *x* component of the (differential) current density in the stationary state at position *x* of particles with velocity components v_x and v_y .

Compared to the first method of measuring velocity distributions, this one is also statistically more effective for de-



FIG. 10. (Color online) Rescaled velocity distributions f_x for different *x*. [System parameters: α =0.9, L_x =20, λ =10.24 (*N*=256, ϕ_0 =0.4).]

termining rare events, such as the high-velocity tails in Sec. VI B. However, as far as $f_x(x,v_x)$ is concerned for small v_x , the second method is inferior to the first, because Eq. (19) assigns a large weight to the relevant events and therefore amplifies statistical fluctuations too.

A. Effects of the discontinuity at a driving wall

Particle-number conservation at a driving wall requires the incoming particle flux at the wall to be equal to the outgoing flux. For the velocity distribution f_x this implies the boundary condition [30],

$$f_{x}(\mp L_{x}/2, v_{x}) = \Theta(\pm v_{x} - v_{\text{drive}}) \left(1 \mp \frac{v_{\text{drive}}}{v_{x}}\right)$$
$$\times f_{x}(\mp L_{x}/2, -v_{x} + v_{\text{drive}}), \quad (21)$$

which must hold for all $v_x > 0$. Here $\Theta(z)=1$ for $z \ge 0$, respectively $\Theta(z)=0$ for z < 0, denotes Heaviside's unit-step function and $v_{\text{drive}}=1$ in our units chosen. The boundary condition (21) relates the distribution f_x of velocities prior to a collision with the wall to that after a collision with the wall. In contrast, the velocity distribution f_y must obey the usual reflection symmetry in v_y everywhere in the system, that is,

$$f_y(x,v_y) = f_y(x,-v_y),$$
 (22)

for all $|x| \leq L_x$ and all $v_y > 0$.

We measured the velocity distribution $f_x(x, v_x)$ at 25 different positions in a moderately inelastic system with coefficient of restitution $\alpha = 0.9$. Figure 10 shows the rescaled velocity distribution $\tilde{f}_x[x, v_x/\sqrt{T_x(x)}] = f_x(x, v_x)\sqrt{T_x(x)}$, measured using the first method. At the driving walls \tilde{f}_x is seen to obey the boundary condition (21). When moving from a driving wall towards the center of the system, the gap in \tilde{f}_x gradually gets smeared out, and \tilde{f}_x becomes more and more symmetric. Clearly, even for this moderately inelastic system \tilde{f}_x does not scale for different x. The two extreme cases for $x = -L_x/2$ and x = 0 are shown together in Fig. 11. The data for \tilde{f}_x at the wall were obtained with the second



FIG. 11. (Color online) Two rescaled velocity distributions f_x from Fig. 10: at the left wall and in the middle of the system.

method for measuring velocity distributions. This result is in agreement with direct-simulation-Monte Carlo results in Ref. [30] and molecular-dynamics simulations in Ref. [31].

B. High-velocity tails

Our main result for the velocity distributions is the nonscaling and multiformity of their tails. In order to observe these phenomena, extensive simulations for capturing rare events are required and the data have to be analyzed on a logarithmic scale. In contrast, on a *linear* scale the rescaled velocity distributions seem to collapse approximately, as was observed previously in e.g., Refs. [2,69]. As an example we show in Fig. 12 the rescaled distributions $\tilde{f}_y(x, v_y/\sqrt{T_y(x)})$ = $f_y(x, v_y)\sqrt{T_y(x)}$ in the middle of the sample (x=0) and at a



FIG. 12. (Color online) Combined plot of rescaled velocity distributions of the same system as in Figs. 10 and 11: \tilde{f}_x in the middle of the sample (green crosses), \tilde{f}_y at a driving wall (red squares), and \tilde{f}_y in the middle (blue triangles). Also shown is a centered Gaussian with unit variance (black solid line). The inset shows the center part of the main graph with the (1433/2964/1340) data points being smoothed [running average over (41/85/38) data points].

driving wall $(x=\pm L_r/2)$ for the same moderately inelastic system (α =0.9) as that in Figs. 10 and 11. Clearly, the curves are reflection symmetric in v_{y} . For comparison, we have also included the symmetric distribution \tilde{f}_x in the middle of the system. Deviations are visible for small velocities only, as the inset of Fig. 12 shows. The corresponding Gaussian (solid line) in Fig. 12 suggests that the velocity distributions are close to but not identical to a Maxwellian. The approximate data collapse, observed on this level of accuracy in Fig. 12, even continues to hold if the coefficient of restitution of the gas is varied to not too large an extent. In considerably more inelastic systems, such as for $\alpha = 0.5$, this is not true any longer. For example, the peak of \tilde{f}_x measured in the middle of the system would be considerably broader and flatter than the ones of \tilde{f}_{v} in the center and at the wall (not shown).

In contrast, Fig. 13(a) shows the same data in Fig. 12 on a *semilogarithmic* scale and also includes velocities of much higher absolute values. From Fig. 13(a) it is evident that scaling does not hold in the high-velocity tails of the distributions either. Similar observations were made before in, e.g., Refs. [31,45,46]. The type of decay in the high-velocity tails is different for \tilde{f}_x and \tilde{f}_y , and also depends on the position in the sample, the coefficient of restitution α , and the global area fraction ϕ_0 . This is illustrated by examples of different systems in Fig. 13. In the insets we show $|\ln \tilde{f}_i|^{-1}$ on a double-logarithmic scale in order to determine the decay exponent β defined by

$$\ln \tilde{f}_{y}(0, v_{y} / \sqrt{T_{y}(0)}) \overset{|v_{y}| \to \infty}{\sim} - |v_{y} / \sqrt{T_{y}(0)}|^{\beta}$$
(23)

for \tilde{f}_y in the middle of the sample. The exponent is defined accordingly for \tilde{f}_{y} at the wall and \tilde{f}_{x} in the middle of the sample. For the moderately inelastic systems with $\alpha = 0.9$ in the first row of Fig. 13, the asymptotics have been clearly reached. We note that β is different for the different distributions, and also depends on the global area fraction ϕ_0 . Upon lowering α (top to bottom in Fig. 13) and/or decreasing the global area fraction ϕ_0 (left to right), the tails of f_v get more and more populated, that is, β decreases. In very simple terms, this may be understood from the fact that (i) the largest typical velocities are always of the order of $v_{\text{drive}} = 1$ (also see Fig. 14) and (ii) that $v_{\text{drive}} / \sqrt{T_i(x)}$ increases up to 20 with decreases of α and ϕ_0 . Hence, a Maxwellian velocity distribution would not be able to supply enough probability to particles with velocities of the order of v_{drive} , instead higher-populated tails are needed. This argument suggests different behavior in different velocity regions so that the distributions cannot be fitted to the functional form (23) over the entire range of velocities [45]. Indeed, such a behavior can be seen in Figs. 13(d)-13(f). The final asymptotics could not always be deduced from the simulations, even though our data include velocities which are up to 40 times bigger than the appropriate granular velocities $\sqrt{T_i}$. This applies to \tilde{f}_{v} in the middle of the sample in Fig. 13(e), where we suspect that the final asymptotics have not been reached. An asymptotic analysis of f_x in the middle of the

sample is even more problematic due to particles that reach the system center from a driving wall without undergoing a collision. These particles give rise to the side peaks of \tilde{f}_x in Figs. 13(e) and 13(f), which has prevented us from determining β in these cases.

Figure 13 contains smoothed data like those in the inset of Fig. 12, and the scales of the horizontal axes are determined by the square root of the appropriate granular temperatures. For comparison, Fig. 14 shows the unsmoothed data corresponding to Fig. 13(f), plotted directly versus v_i (in units of v_{drive}). The side peaks of \tilde{f}_x in the middle of the sample are due to the above-mentioned particles which fly from a driving wall to the system center without undergoing a collision. For the vast majority of particles $|v_y|$ is less than 80% of the driving velocity. It is in the region of this highest velocity observed for \tilde{f}_y in the middle of the sample, that \tilde{f}_x abruptly changes its slope.

The exponent β has also been determined experimentally in a strongly driven gas so that gravity effects are small [16]. A value of $\beta \approx 1.55$ was measured for a gas with coefficient of restitution $\alpha \approx 0.93$. It was found to be remarkably independent of the global area fraction ϕ_0 , which was varied from 0.05 to 0.3. We have also simulated the system of Ref. [16] with zero gravity (not shown) and reproduced those authors' value for β . Even though the experimental data cover a wide range of velocities, some of the interesting phenomena discussed in Fig. 13 cannot be observed in this range. For the same reason the significance of the particular value $\beta \approx 1.55$ should not be overestimated.

Simulations of a driven granular gas in a circularly shaped box also yield stretched Gaussian tails [45]. In addition, evidence is given that β depends only on the coefficient of restitution and the average ratio of the number of particle– wall collisions and particle–particle collisions. For homogeneously driven systems the tail of the velocity distribution was theoretically predicted [23] to be governed by the decay exponent β =3/2. Simulations of homogeneously driven systems [70,71] observed the exponent β =3/2 only for unrealistically low values of the coefficient of restitution.

VII. SUMMARY AND OUTLOOK

In this paper we explored the steady-state properties of a granular gas, driven by vibrating walls. We have measured the full stress tensor, including the collisional contribution. This allowed us to obtain a constitutive equation directly from the simulations. The constitutive equation relates local density, pressure, and temperature. For small inelasticities it can be regarded as the local equation of state of the gas because it is to a large extent independent of the global system parameters. For strongly inelastic systems this interpretation cannot be sustained, instead the constitutive relation depends on the global volume fraction and sample geometry. We have also measured local velocity distributions, whose high-velocity tails were found to depend on the position in the sample as well as on the coefficient of restitution and the global volume fraction. Moreover, the tails are different for the two velocity components parallel and perpendicular to



FIG. 13. (Color online) Rescaled velocity distributions \tilde{f}_x in the middle of the sample (green crosses), \tilde{f}_y at a driving wall (red squares), and \tilde{f}_y in the middle (blue triangles) like in Fig. 12, but here smoothed data are shown on a semilogrithmic scale and velocities of much higher absolute values are included. The solid line is a centered Gaussian with unit variance. (a)–(f) Different systems with decreasing global area fraction ϕ_0 (left to right) and decreasing coefficient of restitution α (top to bottom). The remaining system data are $L_x=20$, λ =10.24(N=256) for the left column and $L_x=50$, $\lambda=9.6(N=240)$ for the right column. The insets show $|\ln \tilde{f}_i|^{-1}$ on a double-logarithmic scale to determine the decay exponent β from Eq. (23). The dashed line is the best linear fit to \tilde{f}_x in the middle of the sample the solid line to \tilde{f}_y in the middle, and the dotted line to \tilde{f}_y at the wall. (a) System corresponding to that in Fig. 12.



FIG. 14. (Color online) Velocity distributions for the system in Fig. 13(f), but with unsmoothed data and without rescaling of the horizontal axis.

the driving walls. To conclude, the stationary state of a driven granular gas is in general *nonuniversal*—in contrast to the corresponding elastic system. This is not unexpected because the driven granular gas is a nonequilibrium state. Furthermore, driving the system by energy input through the walls is effective only if the distance between the driving walls is finite so that the sample geometry enters naturally.

We plan to extend our studies in various directions. It is straightforward to also consider rotational motion of the disks by generalizing the collision rules to include tangential restitution as well as Coulomb friction. Again, the question arises as to what extent the additional parameters affect the local equation of state. In this context it would be interesting to study the effects of gravity, which is also important for comparison with experiments. Finally, one might also investigate other driving mechanisms, like vibrations with finite frequency and nonzero amplitude. Work along these lines is in progress.

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APPENDIX: SIMPLE ENERGY-BALANCE ARGUMENT

Here we derive the energy balance [Eq. (12)] by using arguments like those in Refs. [33,63]. We do this in a slightly more general setting than needed for Eq. (12) and allow in addition for inelastic collisions with the wall, characterized by a coefficient of restitution α_w . Appropriate generalization of the collision rule (4) includes both the driving velocity and the coefficient of restitution with the wall,

with

$$\Delta \boldsymbol{v}_{\text{pw}} = \left[-(1 + \alpha_w) \boldsymbol{v}_x \pm \boldsymbol{v}_{\text{drive}} \right] \boldsymbol{e}_x. \tag{A1}$$

The special case of $v_{\text{drive}}=0$ and $\alpha_w > 1$ provides an alternative driving mechanism, which, however, does not give rise to a stationary state, as will be shown below. In order to be able to treat this limit and for ease in reading of the formulas, we refrain from using dimensionless physical units in this appendix.

 $\boldsymbol{v}' = \boldsymbol{v} + \Delta \boldsymbol{v}_{\text{pw}}$

The average energy gain ΔE_{pw} due to a single particlewall collision is estimated from Eq. (A1) by averaging the kinetic energy before and after the collison with a Maxwellian velocity distribution with (global) temperature *T*. This gives

$$\Delta E_{\rm pw} = \frac{m}{2} \left(v_{\rm drive}^2 + 4\alpha_{\rm w} \sqrt{\frac{T}{2\pi m}} v_{\rm drive} - (1 - \alpha_{\rm w}^2) \frac{T}{m} \right). \tag{A2}$$

The collision frequency of particles with the left (right) wall is estimated by

$$f_{\rm pw} = \frac{N}{L_x} \sqrt{\frac{T}{2\pi m}},\tag{A3}$$

where we have assumed the density to be spatially homogeneous throughout the system.

When two disks collide in the bulk, the average change in total energy is computed from Eqs. (2) and (3) in a manner similar to that for Eq. (A2) by averaging pre- and postcollisional kinetic energy, which yields

$$\Delta E_{\rm pp} = -\frac{1-\alpha^2}{2}T.$$
 (A4)

Finally, the number of particle–particle collisions per time is given approximately by Enskog's collision frequency [23],

$$f_{\rm pp} = \frac{N}{L_x} \lambda a \chi \sqrt{\frac{T\pi}{m}},\tag{A5}$$

where χ is the pair correlation function at contact of a corresponding elastic gas in thermal equilibrium. Since χ is not known exactly, we resort to the widely used Henderson approximation [72],

$$\chi \approx \frac{1 - 7\phi/16}{(1 - \phi)^2},$$
(A6)

for numerical purposes. It may be viewed as a heuristic rational approximation to the virial expansion of χ and is the two-dimensional equivalent to the Carnahan–Starling approximation [66] for a three-dimensional hard-sphere gas. Additional higher-order terms to the Henderson approximation, which are proportional to $\phi^3/(1-\phi)^4$, have turned out to be irrelevant for our purposes and will therefore not be taken into account.

Summing over energy loss in the bulk and energy gain at the right and left walls, we obtain for the total change in granular temperature,

$$\frac{dT}{dt} \approx \frac{2f_{\rm pw}E_{\rm pw} + f_{\rm pp}E_{\rm pp}}{N}$$
$$= \frac{m}{L_x}\sqrt{\frac{T}{2\pi m}} \left(v_{\rm drive}^2 + 4\alpha_{\rm w}\sqrt{\frac{T}{2\pi m}}v_{\rm drive} - \frac{\pi\psi_{\rm eff}T}{2m}\right).$$
(A7)

Here the dimensionless parameter,

$$\psi_{\rm eff} = \psi - \frac{2}{\pi} (\alpha_w^2 - 1),$$
 (A8)

is given in terms of $\psi = \sqrt{2\lambda a \chi (1 - \alpha^2)}$, which was already introduced in the text below Eq. (12). (We note that λa is the dimensionless line density employed there.) We briefly discuss two special cases:

(a) For $v_{\text{drive}}=0$ and $\alpha_w > 1$ no stationary state is reached in general. Both energy gain and loss increase like $T^{3/2}$, resulting in Haff's law,

$$\frac{dT}{dt} = -\frac{\psi_{\rm eff}}{2L_x} \sqrt{\frac{\pi}{2m}} T^{3/2},\tag{A9}$$

which has been discussed extensively in a different context of freely cooling granular gases. Here the temperature continues to decrease or increase depending on whether dissipation or driving wins.

(b) For $\alpha_w = 1$ and $v_{\text{drive}} > 0$ the granular temperature adjusts to the driving so that the stationary state with dT/dt=0 is characterized by the quadratic equation,

$$1 + 2\sqrt{\frac{2\bar{T}}{\pi}} - \frac{\pi\psi}{2}\bar{T} = 0.$$
 (A10)

Solution of Eq. (A10) for the dimensionless global temperature $\overline{T} = T/(mv_{\text{drive}}^2)$ is given in Eq. (12).

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