Evidence of a higher-order singularity in dense short-ranged attractive colloids

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We study a model in which particles interact through a hard-core repulsion complemented by a short-ranged attractive potential, of the kind found in colloidal suspensions. Combining theoretical and numerical work we locate the line of higher-order glass transition singularities and its end-point — named $A_4$ — on the fluid-glass line. Close to the $A_4$ point, we detect logarithmic decay of density correlations and sub linear power-law increase of the mean square displacement, for time intervals up to four order of magnitudes. We establish the presence of the $A_4$ singularity by studying how the range of the potential affects the time-window where anomalous dynamics is observed.

Recently, a great deal of interest has grown around dynamical phenomena arising in dense colloidal suspensions when particles interact through a hard-core repulsion followed by a rather short-ranged attractive potential [1–3]. Some of these unusual phenomena, initially predicted theoretically [4–6], have been observed experimentally [7–10] and in numerical simulations [11–14,23]. The novelty, as compared to the well-studied hard-sphere-colloids case, comes from the possibility of generating structurally arrested states with an additional physical mechanism, driven by the range of the attractive interaction. Indeed, in the case of hard-sphere, structural arrest is controlled by the well-known excluded volume cage effect, where particle motion is hindered by the presence of neighboring ones. This mechanism generates a localization length of the order of 10 % of the particle diameter $\sigma$. In the additional mechanism, the motion of the particles is restricted by the adhesiveness, i.e. the formation of non-permanent bonds due to the attractive part of the inter-particle potential; particles are confined by bonds and the corresponding localization length is fixed by the attraction range $\Delta$. When $\Delta \ll \sigma$, the interplay of the two localization mechanisms creates liquid states for packing fractions higher than those possible for a pure hard-sphere system. This means that the attraction stabilizes the liquid and leads to a reentry phenomenon, where this new liquid state can arrest into a glass by cooling as well as heating. The experimental and numerical confirmation of this prediction [9–14] suggests that the mode-coupling theory (MCT) [15], which was the basis of the cited theoretical work [4–6], can contribute to explain the slow dynamics in short-ranged attractive colloids.

These systems are characterized by three control parameters — the packing fraction $\phi$, the ratio $\theta$ of the thermal energy $k_B T$ to the typical well depth $u_0$ and the range $\Delta$ of the attractive potential. Within MCT, the phase diagram of this three-dimensional control parameter space (shown in Fig.1) is organized around a critical point $(\phi^*, \theta^*, \Delta^*)$, referred to as a type $A_4$ higher-order glass transition singularity in MCT classification. $A_4$ is the end-point of a line of higher-order singularities (of type $A_4$). From a physical point of view, $A_4$ is characterized for being the only singularity point accessible from the liquid phase. For $\Delta > \Delta^*$ no singular points are predicted by the theory, while for $\Delta < \Delta^*$ the $A_3$ singularity points are buried in the glass phase and their presence can be only observed indirectly [7,9,11,14]. Near the $A_4$ singularity, MCT predicts a structural relaxation dynamics, which is utterly different from that known for conventional glass-forming liquids. It is ruled by logarithmic variations of correlators as function of time and subdiffusive increase of the mean squared displacement (MSD). Theory makes precise predictions for the time-interval where this unusual dynamics is expected, as well as for its variation with changes of the $\phi$, $\theta$ and $\Delta$ parameters. Indications of logarithmic dynamics have been reported [7,9,11,14], but neither a systematic study of the dependence on the control parameters has been published nor the $A_4$ point has been located numerically. It is in fact difficult to discriminate between straight lines caused by some inflection points in the usual logarithmic representation of correlators and the logarithmic decay laws generated by the higher-order singularity. In this Letter, we establish the existence of the location of the $A_4$ point for a well defined model, amenable of simultaneous numerical and theoretical treatment. Reported data — with control parameters explicitly chosen close to the $A_4$ point — exhibit the mentioned laws over time intervals up to four orders of magnitude. More importantly, we show that the decay patterns vary with changes of the control parameters and wave-vectors as expected [16,19], properly testing the theoretical predictions.

We simulate a 50% – 50% binary mixture of $N = 700$ hard spheres of mass $m$, with diameters $\sigma_{AA}$ and $\sigma_{BB}$ and ratio $\sigma_{AA}/\sigma_{BB} = 1.2$. The hard core between particles of different type $\sigma_{AB} = 0.5(\sigma_{AA} + \sigma_{BB})$. The hard core potential is complemented by an attractive square well potential of depth $u_0$, independent on the particle...
formed to reduce noise. Equilibration has been carefully parameter within the accuracy of our calculations. Av-
ing, in previous simulations [14]. As discussed in the follow-

\[ \Delta_{ij} \] differing in the width of the square well \( \Delta_{ij} \), \( S_1 \) and \( S_2 \) have \( \Delta_{ij} = 0.031\sigma_{ij} \), and \( \Delta_{ij} = 0.043\sigma_{ij} \) respectively \( (i = A, B) \). The system \( S_1 \) has been extensively studied in previous simulations [14]. As discussed in the follow-

\( \phi \) as a function of \( S \) through the vanishing of the diffusion (open symbols) and cal-

\( S \) is chosen to coincide with the critical amplitude \( \phi_c \) of the \( S_1 \) line. The MCT curves have been transformed according to the mapping discussed in the text. The \( T \) and \( \phi \) of the simulated point are indicated by the black dot. The inset shows the power-law fits for the isothermal diffusion coefficient \( D \) (data from Ref. [14]) for \( S_1 \) as a function of \( \phi - \phi_c \) (see text).

To estimate the predicted location of the \( A_4 \) point for the binary mixture square well model (and consequently select the parameters to be used in the simulations) we proceed as follows. First, we calculate the glass line \( \phi_c(T) \) for the mixture \( S_1 \) extrapolating diffusivity \( D \) data [14] according to \( D \sim (\phi - \phi_c(T))^{\gamma(T)} \), for eight different \( T \). Second, we calculate the MCT ideal glass transition line. The partial structure factors, the only input needed, are calculated with the Percus-Yevick approximation, numerically solving the Ornstein-Zernike equation [18]. Third, we map the theoretical glass line on the simulation line, following the procedure first used by Sperl [19]. Indeed MCT cannot reproduce the numerical values for \( \phi_c \). As shown in Fig.1, the linear transformation \( \phi \rightarrow 1.897 \phi - 0.3922 \) and \( T \rightarrow 0.5882 T - 0.225 \), al-

allows to superimpose the MCT result with the simulation data, in the studied region of \( T \) and \( \phi \). By calculating the MCT ideal glass transition line for several values of \( \Delta \), the theoretical location of the \( A_4 \) point for the binary mixture square-well model is obtained. Assuming that, to a first approximation, the same linear transformation holds also for the critical value of the well width, we find that the \( A_4 \) location maps to \( \Delta^*_{ij} = 0.043\sigma_{ij} \), \( \phi^* = 0.611 \) and \( \theta^* = 0.416 \). This allows us to select a state point with parameters close to the critical ones and perform a simulation (in thermodynamic equilibrium) close to \( A_4 \).

We choose [20] \( \theta = 0.4 \) and \( \phi = 0.6075 \) and compare, for the same \( T \) and \( \phi \), the dynamics for \( \Delta^*_{ij} \) (system \( S_2 \)) and for the close-by value \( \Delta_{ij} = 0.031\sigma_{ij} \) (\( S_1 \)).

Fig.2 shows the MSD for the larger species \( \langle \delta r^2_{ij} \rangle \) for both systems. We notice the presence of a subdiffusive regime at intermediate times, i.e. a variation according to \( \ln(\langle \delta r^2_{ij} \rangle) = f + a \ln(t/\tau) \) for about three decades for \( S_1 \), where \( a_1 = 0.44 \). The logarithmic regime extends by more than a decade if the attraction range is increased to that for system \( S_2 \). Simultaneously the exponent \( a \) decreases to \( a_2 = 0.28 \). Fig. 3 shows the density-fluctuation auto-correlation functions \( \Phi_q(t) \) for a representative set of wave vectors \( q \). The decay curves in Fig. 3 do not show the two steps scenario with a plateau characteristic of conventional glass forming liquids. Indeed it is impossible to fit these curves for large \( q \) with the standard stretched exponential function. Instead, there is a region of clear logarithmic decay at \( q^* \sigma_{BB} = 23.5 \) for \( S_1 \) and at \( q^* \sigma_{BB} = 16.8 \) for \( S_2 \). The time-intervals over which logarithmic decay is observed are of similar size as those for the MSD. For \( q < q^* \), the \( \Phi_q \) vs \( \ln t \) curves are concave, and for \( q > q^* \) convex.
To show that the above described features are consistent with the ones predicted by MCT, we cite the general asymptotic decay law for the correlation function of a generic variable $X$, near the higher-order singularity [16]

$$\Phi_X(t) = f_X - h_X \left[ B^{(1)} \ln(t/\tau) + B^{(2)} X \ln^2(t/\tau) \right]. \tag{1}$$

Here $\tau$ abbreviates a time-scale which diverges if the state approaches the singularity. The formula is obtained by asymptotic solution of the MCT equations, using the parameter differences $\phi^* - \phi, \theta^* - \theta, \Delta^* - \Delta$ as small quantities, say, of order $\epsilon$. The coefficient $B^{(1)}$ is of order $\sqrt{\epsilon}$ while $B^{(2)}$ is of order $\epsilon$. The amplitude $h_X$ is independent of $\epsilon$. The first term $f_X$ is the sum of the non-ergodicity parameter of variable $X$ at the singularity and a correction of order $\epsilon$. Terms of order $\epsilon^{3/2}$ are neglected [16]. We can interpret $\ln(h^2_A)$ also according to Eq. (1). Hence, the straight line for $S_1$ with slope $a_1$ shown in Fig. 2 is consistent with the assumption that $S_1$ is close to the singularity. Changing from $S_1$ to $S_2$, the slope $a = h_{BB} B^{(1)}$ has to decrease and the range of validity of the leading order description has to expand. This is demonstrated impressively by the data.

![FIG. 3. A-type particle density correlators for $S_1$ (top panel) and $S_2$ (bottom panel), showing six different wave vectors (from top to bottom $q_{BB} = 6.7, 11.7, 16.8, 23.5, 33.5$ and 50.3). For correlators indicated by an arrow, a logarithmic behavior is observed within the selected time window. The dashed lines show the fits according to Eq.1. The vertical dashed lines indicate the fitting interval.](image)

![FIG. 4. The fitting parameters $f_q, H^{(1)}_q$ and $H^{(2)}_q$ of the asymptotic logarithmic law, for $S_1$ and $S_2$. The central panel also shows that a multiplication by 2.43 of $H^{(1)}_q$ for the $S_2$ system gives $H^{(1)}_q$ for $S_1$, confirming the factorization of $H^{(1)}_q$. At the wave vector where $H^{(2)}_q = 0$ ($q_1^2 \sigma_{BB} = 23.5$ for $S_1$ and $q_2^2 \sigma_{BB} = 16.8$ for $S_2$) correlation functions display a pure logarithmic decay.](image)

The fit parameters for the coefficient $H^{(1)}_q \equiv h_q B^{(1)}$ are reported in the middle panel of Fig. 4. The decrease of $H^{(1)}_q$ upon changing from $S_1$ to $S_2$ is in agreement with the prediction that $B^{(1)}$ tends to zero with $\epsilon$ approaching zero. As shown in Eq.1, $H^{(1)}_q$ factorizes into a control-parameter independent factor $h_q$, depending on $q$, and a control-parameter dependent factor $B^{(1)}$, independent of $q$. This implies that the $q$-dependence of $H^{(1)}_q$ should be the same for $S_1$ and $S_2$. As shown in Fig.4, this property is verified by the data. The same property does not apply to $H^{(2)}_q \equiv h_q B^{(2)}$ because of the $q$-dependence of $B^{(2)}$. Moreover, the fitting results confirm that $H^{(2)}_q$ is smaller than $H^{(1)}_q$, as expected being the first of order $\epsilon$ and the second of order $\sqrt{\epsilon}$. The wave-vector value where $B^{(2)}_q = 0$ allows to identify the characteristic length scale associated to the pure logarithmic decay (see Fig.3). Such length scales are much shorter than the typical first neighbor shell.

To explicitly test the mentioned factorization, Fig. 5 shows a set of correlators rescaled to $\tilde{\Phi}_q(t) = (\Phi_q(t) - f_q)/H^{(1)}_q$. MCT predicts that in leading order $\epsilon$ all $\tilde{\Phi}_q(t)$ vs $\ln(t)$ curves collapse on the logarithmic decay law.

To show that the above described features are consistent with the ones predicted by MCT, we cite the general asymptotic decay law for the correlation function of a generic variable $X$, near the higher-order singularity [16]
\[ -\ln(t/\tau). \] This is indeed shown for a large time interval that expands upon approaching the singularity, as it is demonstrated by comparing results for \( S_1 \) and \( S_2 \) [22].

Results reported in this Letter have an intrinsic value associated to the observation, in a particularly simple system — a square well potential —, of a particularly complex dynamics over more than four decades in time. They show that MCT, a theory essentially developed to address the problem of the excluded volume glass transition, is able — without any modification — to handle the logarithmic dynamics and to provide an interpretation scheme in term of \( A_1 \) point. The numerical results and the comparison with the theoretical predictions do constitute, in fact, a stringent verification that the logarithmic dynamics in the density autocorrelation functions and the subdiffusive behavior in the MSD can be fully rationalized by MCT. Indeed, data in Figs.3 and 4 are in agreement with MCT predictions (Fig. 12 and 13 of Ref. [19]). The simplicity of the model and the complexity of the dynamics suggest that the short-range attractive colloids have the potentiality to become a benchmark for the development of extended theories of the glass transition.

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