Single particle jumps in a binary Lennard-Jones system below the glass transition

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We study a binary Lennard-Jones system below the glass transition with molecular dynamics simulations. To investigate the dynamics we focus on events (jumps) where a particle escapes the cage formed by its neighbors. Using single particle trajectories we define a jump by comparing for each particle its fluctuations with its changes in average position. We find two kinds of jumps: “reversible jumps,” where a particle jumps back and forth between two or more average positions, and “irreversible jumps,” where a particle does not return to any of its former average positions, i.e., successfully escapes its cage. For all investigated temperatures both kinds of particles jump and both reversible and irreversible jumps occur. With increasing temperature, relaxation is enhanced by an increasing number of jumps and growing jump lengths in position and potential energy. However, the waiting time between two successive jumps is independent of temperature. This temperature independence might be due to aging, which is present in our system. We therefore also present a comparison of simulation data with three different histories. The ratio of irreversible to reversible jumps is also increasing with increasing temperature, which we interpret as a consequence of the increased likelihood of changes in the cages, i.e., a blocking of the “entrance” back into the previous cage. In accordance with this interpretation, the fluctuations both in position and energy are increasing with increasing temperature. A comparison of the fluctuations of jumping particles and nonjumping particles indicates that jumping particles are more mobile even when not jumping. The jumps in energy normalized by their fluctuations are decreasing with increasing temperature, which is consistent with relaxation being increasingly driven by thermal fluctuations. In accordance with subdiffusive behavior are the distributions of waiting times and jump lengths in position.

I. INTRODUCTION

If a liquid is cooled and crystallization is avoided, one obtains a supercooled liquid. Upon further cooling the system falls out of equilibrium and results in a glass. During the transition from liquid to supercooled liquid to glass the thermodynamic properties change and even more drastic changes occur in the dynamics. The viscosity increases by many orders of magnitude upon cooling and the mean square displacement (MSD) as a function of time develops a plateau at intermediate times. The lower the temperature the longer the waiting time within the plateau until a second rise in the MSD occurs. One common explanation for the plateau is that while at high temperatures one has (at late enough times) normal diffusion, at lower temperatures particles are caged in, i.e., trapped by their neighbors, and spend longer time within this cage with decreasing temperatures. The second increase in the MSD indicates that after long enough waiting time the particles manage to escape their cage. This escape out of the cage (jump) is the focus of this paper.

To set the work of this paper in context, we review briefly previous and to our work related studies on the dynamics of supercooled liquids and glasses.

Central quantities in both experiments and computer simulations are the viscosity, diffusion constant, and MSD. The MSD, which is an average over single particle

\[ r^2(t) = \frac{1}{N} \sum_i \left| \mathbf{r}_i(t) - \mathbf{r}_i(0) \right|^2, \]

and variations thereof show jumps when viewed with fine enough time resolution, and when studied at low enough temperatures and for single configurations. Detailed studies of these jumps indicate collective motion.

One learns about the interplay between regular diffusion and hopping motion via four-point correlation functions and the van Hove correlation function

\[ G(r,t) = G_s(r,t) + G_d(r,t) \]

(Refs. 15–24) where \( G_s(r,t) \) is the self part and \( G_d(r,t) \) is the distinct part. The hopping shows up in an additional peak in \( G_s(r,t) \) and an increase in the amplitude at \( r \rightarrow 0 \) of \( G_d(r,t) \).

Recently attention has been drawn to the non-Gaussian tail of \( G(r,t) \) (Refs. 25–29) and the non-Gaussian parameter \( \alpha_2 \) (Refs. 5, 13, 25 and 29–35) which is the second coefficient of the cumulant series of \( F(q,t) \), the Fourier transform of \( G(r,t) \). They are signatures of nonexponential behavior, which might be either due to homogeneous complex dynamics and/or due to spatially heterogeneous dynamics.

The dynamics of glasses out of equilibrium, i.e., systems which have been quenched from high to low temperature,
displays additional complexity, since the system might "age." This means that the dynamics depends on the waiting time between the quench and the measurement which results in the violation of the fluctuation dissipation theorem.

Another fruitful approach to gain insight into the dynamics of supercooled liquids and glasses has been to investigate the energy landscape of the inherent structure that means of the instantaneous configurations which have been quenched to their local potential energy minimum. At the same temperature when the system starts slowing down drastically, the potential energy of the inherent structure undergoes a qualitative change. The long time behavior of this potential energy shows jumps between metabasins, where the latter are groups of strongly correlated local minima. The mean average waiting time between these jumps turns out to be dominated by the long times. Together with the diffusion constant allow an estimate of the cage size.

Direct studies of the cage have been done via three-time correlations, velocity-velocity correlations, the cage correlation function, and a passage time before a particle escapes. The results of these studies tell us about cage properties such as the cage size and waiting time within a cage (or jump rate).

Hand in hand with the experiments and simulations are the theoretical models for supercooled liquids and glasses. One of the very successful theories is the mode coupling theory of the glass transition (MCT), which describes the dynamics of supercooled liquids via nonlinear dynamics of coupled density modes and makes predictions for quantities such as $D$, $F(q, t)$, and the susceptibility $\chi^{\alpha}(\omega)$. The extended version of MCT includes hopping processes via the coupling to current fluctuations. The comparison of experiment and this theory for $\chi^{\alpha}(\omega)$ (Ref. 56) and $\alpha_2$ (Ref. 57) shows very good agreement.

The MCT is a theory for the glass transition for temperatures above the glass transition. Below the glass transition there exists no equivalent of a microscopic theory such as MCT. Some phenomenological theories which are more general for solids build in hopping either indirectly or in the free volume model or directly in hopping models.

The goal of this paper is to obtain via single particle trajectories direct information about jump processes such as jump sizes and waiting times between jumps. Specific examples of the dynamics of single particle jumps, in the form of a plot of one component of $r_i(t)$ (Refs. 16–18 and 62) or in the form of a two- or three-dimensional picture of the particle trajectory, are very helpful to get an idea of some qualitative features of jumps, such as the detailed geometry of jump processes. In this paper, however, we go beyond single examples by defining a systematic search algorithm for jump processes. For the case of single particle trajectories a similar approach has been used in the work, where the jump criterion is a minimum hopping distance, which is the same for all particles. In contrast, we use a relative criterion, where for each particle its size of fluctuations is compared to its jump size. We choose this relative criterion to be able to identify jumps of particles of different sizes and neighborhood. That means the criterion is adjusted to the cage size of each individual particle. As criterion for the occurrence of a jump we use the positions of particles instead of their energy. While both approaches are fruitful, we believe that many theoretical models and interpretations of simulations and experiments are based on an intuitive picture of the particle motion in real space. Similarly our goal is to mimic a careful observer of each particles’ motion in our system. We therefore use single particle trajectories, instead of a quantity which is an average over all particles (as in the work Refs. 3–10). For the case of single particle trajectories the distinction between vibrational and hopping motion turns out to be clearer by taking time averages than by using trajectories of the inherent structure.

In the following we define our model, and give details about the simulation (Sec. II). Our precise definition of a jump is given in Sec. III. We find two types of jumps (irreversible and reversible). The latter are distinguished for the rest of the paper (in distinction from Refs. 17, 19, 35 and 64). In Sec. IV we count as a function of temperature the number of jumping particles, and in Sec. V the number of visited different cages. In Sec. VI we investigate the times during a jump and between successive jumps. The jump size both in position and in potential energy are presented in Secs. VII and VIII. Section IX addresses the question of history dependence of the jump results. In Sec. X we conclude with a summary of our results, a comparison with the results of previous work and with our resulting picture of jump processes. We finish with open questions suggesting future work.

II. MODEL AND SIMULATION

We use a binary Lennard-Jones mixture of 800 $A$ and 200 $B$ particles with the same mass. The interaction potential for particles $i$ and $j$ at positions $r_i$ and $r_j$ and of type $\alpha, \beta \in \{A, B\}$ is

$$V_{\alpha\beta}(r) = 4\epsilon_{\alpha\beta}\left[\frac{\sigma_{\alpha\beta}}{r}\right]^{12} - \left[\frac{\sigma_{\alpha\beta}}{r}\right]^{6},$$

where $r = |r_i - r_j|$ and $\epsilon_{AA} = 1.0, \epsilon_{AB} = 1.5, \epsilon_{BB} = 0.5, \sigma_{AA} = 1.0, \sigma_{AB} = 0.8, \sigma_{BB} = 0.88$. We truncate and shift the potential at $r = 2.5 \sigma_{\alpha\beta}$. From previous investigations it is known that this system is not prone to crystallization and demixing. In the following we will use reduced units where the unit of length is $\sigma_{AA}$, the unit of energy is $\epsilon_{AA}$, and the unit of time is $\sqrt{m\sigma_{AA}^2/(48\epsilon_{AA})}$.

We carry out molecular dynamics (MD) simulations using the velocity Verlet algorithm with a time step of 0.02. The volume is kept constant at $V = 9.4^3 = 831$ and we use periodic boundary conditions. Previous simulations showed that for present day computer simulations the system falls out of equilibrium in the vicinity of $T = 0.435$. We analyze here simulations at $T = 0.15, 0.2, 0.25, 0.3, 0.35, 0.38, 0.4, 0.41, 0.42$, and 0.43 as they have been described in Ref. 67. We start with ten independent well equilibrated configurations at $T = 0.446$. Each of these configurations undergoes the following sequence of simulation runs. After an instantaneous quench to $T = 0.15$ we first run a (NVT) simu-
lution for $10^5$ MD steps. The temperature is kept constant by replacing the velocities of all particles by new velocities drawn from the corresponding Boltzmann distribution every 50 time steps. We then run the simulation for $10^6$ MD steps without the temperature bath (NVT) and subsequently raise the temperature instantaneously to the next higher temperature $T=0.2$. (NVT) and (NVE) runs of $10^5$ MD steps each follow and the temperature is again raised, and so forth. The final configurations of these (NVT) runs are the initial configurations for the (NVE) production runs (for $5 \times 10^6$ MD steps) presented in this paper. During each production run the positions of all particles (configurations) are stored every 2000 MD steps which are then used for analysis.

For the here investigated temperatures the relaxation times $\tau$ are much larger than the waiting time before the production runs (at $T=0.446$, $\tau \approx 8 \times 10^5$). We therefore study relaxation processes out of equilibrium and find aging effects, which is consistent with previous detailed studies of aging of the same binary Lennard-Jones system.\textsuperscript{40,41}

III. DEFINITION OF JUMP AND JUMPTYPE

We focus in this paper on the process of a particle escaping its cage, using single particle trajectories $\mathbf{r}_i(t)$ given by the periodically stored configurations.

To distinguish vibrations around an average position from a change in the average position, we average 20 consecutive positions $\mathbf{r}_i(t)$ to obtain $\mathbf{r}_i(m)$ as sketched in Fig. 1. We identify jumps by comparing changes in these averaged positions $|\Delta \mathbf{r}_i|$ with the fluctuations in position $\sigma_{i,\text{est}}$ for each particle $i$ where

$$|\Delta \mathbf{r}_i| = |\mathbf{r}_i(m) - \mathbf{r}_i(m-4 \times 20)|$$

(see Fig. 1) and $\sigma_{i,\text{est}}$ is defined in Ref. 70. We use in specific for $|\Delta \mathbf{r}_i|$ not consecutive but instead average positions which are four averages (each of 20 configurations) apart. This choice of time interval in $|\Delta \mathbf{r}_i|$ is to allow identification of not only sudden but also more gradual jumps. We define that a jump of particle $i$ occurs whenever

$$|\Delta \mathbf{r}_i|^2 > 20 \sigma_{i,\text{est}}^2. \tag{4}$$

Notice that we use a relative criterion, namely, for each particle $i$ a comparison of $|\Delta \mathbf{r}_i|$ with its $\sigma_{i,\text{est}}$. Our motivation for this relative criterion is that we would like to identify jumps of both $A$ and the smaller $B$ particles. Also even for particles of the same type their jump size might differ due to different cage sizes. $\sigma_{i,\text{est}}^2$ is an estimate for the cage size of each individual particle and is therefore used as criterion for the occurrence of a jump.

When we apply the above jump definition, we find two types of jumps which we call “irreversible” and “reversible” jumps (see Fig. 2). A particle which undergoes an irreversible jump succeeds in escaping its cage (for the time window of the simulation) whereas a particle undergoing a reversible jump returns back to one of its previous average positions. Similar results have been found in previous simulations of other systems.\textsuperscript{3,4,6,8,17,21,62} However, the present work differs from these that we analyze all following quantities for the reversible and irreversible jumps separately. As sketched in Fig. 2, we call a jump reversible whenever case 1 or case 2 occurs. Case 1 corresponds to the situation of a particle undergoing multiple jumps and returning to one or more of any previous average positions (for details see Sec. V). If two average positions are equal then all jumps which happen between the previous position and the recurring position are called reversible jumps. In the example of Fig. 2 this means in case 1 that all shown jumps are reversible jumps. To increase our resolution in time we use not only the information of the averaged positions $\mathbf{r}_i(m)$ but also the complete information of $\mathbf{r}_i(t)$. In case 2 (see Fig. 2) the spikes in $\mathbf{r}_i(t)$ indicate returns to the average position before the jump.\textsuperscript{71} If a jump satisfies neither case 1 nor case 2, then it is called irreversible jump.

In the following we distinguish between the jump events, as they have been defined so far, and the corresponding jumping particles which may undergo multiple jumps of different types. Any reversible jump designates the corre-
FIG. 3. Number of jumping particles normalized by the number of corresponding particle type as a function of temperature $T$. Irreversible and reversible jumpers are distinguished.

responding particle as a reversible jumper for the entire time window.

IV. NUMBER OF JUMPING PARTICLES

We apply now the above definitions to identify all jumps occurring in our simulations. The number of identified jumps depends on the time interval of the production runs and, since the system is out of equilibrium, also on the waiting time before the production runs. The straightforward counting of jump events would give reversible jumpers (which sometimes jump many times) a larger weight. We therefore present the number of jumping particles. Figure 3 shows the number of jumping particles normalized by the number of corresponding particle type as a function of temperature $T$. The straightforward counting of jump events would give reversible jumpers (which sometimes jump many times) a larger weight. We therefore present the number of jumping particles. Figure 3 shows the number of jumping particles normalized by the number $N_{\alpha}$ of particles in the system of type $\alpha \in \{A, B\}$. With increasing temperature $T$ the number of jumping particles increases consistent with an increasing number of relaxation processes. A similar temperature dependence has been found indirectly via the participation ratio in the work. More surprisingly, we find that not only the small $B$ particles are jumping but also a considerable fraction of $A$ particles. However, the fraction of jumping $B$ particles is larger than the fraction of jumping $A$ particles due to the smaller size and therefore higher mobility of the $B$ particles. Furthermore, both irreversible and reversible jumps are not only occurring in a certain temperature range but at all temperatures.

Figure 4 illustrates the fraction of these jumping particles that are irreversible jumpers. At low to intermediate temperatures this ratio is, within the large error bars, roughly constant. At intermediate to larger temperatures irreversible jumpers become more likely than reversible jumpers with increasing temperature. This increase is even more pronounced for all temperatures if one considers the number of irreversible jump events (instead of jumping particles) divided by the number of all jump events (see inset of Fig. 4). Gaukel et al. come to a similar conclusion via a model for their simulation data. Their irreversible jumps become more likely with increasing temperature. We interpret this increase in the fraction of irreversible jumps as sketched in Fig. 5. Both irreversible and reversible jumps start out the same, with the jumping particle leaving the cage (formed by the neighboring particles) possibly through an opening in the cage. In the case of the irreversible jump, the entrance of the cage gets meanwhile blocked by a particle, loosely speaking the door gets closed, and the jumping particle can no longer return and has successfully escaped its cage. With increasing temperature all particles become more mobile, which increases the likelihood of the blockage of the entrance back into the cage which in turn leads to an increase in the fraction of irreversible jumps as shown in Fig. 4. We do not intend to make here a statement about the exact geometric process, for example, that a door made up of a single particle gets open and closed, but more generally the process of rearrangement of the cage. Interestingly enough, this blockage happens more often in the case of the larger $A$ than the smaller $B$ particles which leads to larger ratios for $A$ than $B$ particles in Fig. 4.

V. NUMBER OF AVERAGE POSITIONS

With each jump a particle either returns to one of its former average positions within a cage or to a new overall average position. We call the average positions before and after a jump $\langle \bar{r}_i \rangle_s$ and $\langle \bar{r}_i \rangle_f$ (for details about the time averages $\langle \cdot \rangle_t$ see Fig. 11 and Sec. VII). We use as criterion for two average positions to be the same that the distance between them $\Delta \langle \bar{r}_i \rangle = |\langle \bar{r}_i \rangle_f - \langle \bar{r}_i \rangle_s|$ and the average fluctuations before the jump $\langle \sigma_i^2 \rangle_s$ (for the definition of $\sigma_i^2$ see Ref. 70) satisfy $(\Delta \langle r_i \rangle)^2 \leq 5 \langle \sigma_i^2 \rangle_s$. Figure 6(a) shows an average of

![Diagram showing reversible and irreversible jumps](image)
the number of distinct average positions which are visited by a particle. According to the higher mobility of the smaller $B$ particles, they visit more average positions than the $A$ particles. For low temperatures most jumping particles visit only two average positions during our simulation. For intermediate to larger temperatures, however, not only increasingly more particles jump (see Sec. IV) but these particles also jump more often. The distribution of the number of different visited average positions at the highest temperatures as shown in Fig. 6(b), broadens with increasing $T$ such that some particles visit up to seven different average positions during the simulation run.

VI. TIMES

In this section we investigate the time scale of jumps. As sketched in Fig. 7 we determine the time duration of a jump $\Delta t_d = t_f - t_i$, the time before the first jump of a particle $\Delta t_{\text{head}}$, the time after the last jump of a particle $\Delta t_{\text{tail}}$, and the time between two successive jumps I and II to be $\Delta t_b$.

$$\Delta t_b = t_{i}^{II} - t_{i}^{I}$$

where $t_i$ and $t_f$ indicate the starting and ending time of a jump. Notice that $\Delta t_b$ is only defined if a jump particle jumps twice or more and that we analyze $\Delta t_{\text{head}}$ and $\Delta t_{\text{tail}}$ separately as presented shortly. For the distinction of irreversible and reversible jumps, we assign the jump type of $\Delta t_b$ according to the jump ending $\Delta t_{\text{tail}}$, for example, in Fig. 7 the jump type of $\Delta t_b$ is determined by jump II. This means that $\Delta t_b$ is a measure of the waiting time before a jump occurs. Figure 8 illustrates that $\Delta t_d \ll \Delta t_b$ which tells us in hindsight why we could identify jumps as rare events with the above described procedure. Since the time resolution is of the order of 1000 ($\sim 20 \times 2000 \times 0.02$), i.e., of the same order as $\Delta t_d$, we do not draw further conclusions about the temperature dependence of $\Delta t_d$. However, $\Delta t_b$ is well above our time resolution and small enough to be detected during our simulation run of length $1 \times 10^7$. In accordance with the
picture of reversible jumpers which try but do not succeed to escape their cage, these trials happen on a shorter time scale than irreversible jumps. However, to our surprise, $\Delta t_b$ seems to be independent not only of the particle type but also of temperature. The question arises if this temperature independence of $\Delta t_b$ is a consequence of the finite time window of our simulation. To take this finite time interval $t_{\text{tot}}$ into account, we make a crude approximation to correct the time intervals $\Delta t_b$, $\Delta t_{\text{head}}$, and $\Delta t_{\text{tail}}$ by assuming that jumps happen equally likely at any time in the window $[0, t_{\text{tot}}]$ (which is not accurate due to aging). The probability $P_{\text{corr}}(\Delta t)$ of finding the complete interval $\Delta t$ reduces to the probability of finding $\Delta t$ during $[0, t_{\text{tot}}]$ in the simulation

$$P_{\text{sim}}(\Delta t) = P_{\text{corr}}(\Delta t)\left(\frac{t_{\text{tot}}-\Delta t}{t_{\text{tot}}}\right)^c,$$

where $c$ is a normalization constant. We may approximate $c$ with $\langle t_{\text{tot}}/(t_{\text{tot}}-\Delta t) \rangle_{\text{sim}}^{-1}$ where $\langle \cdot \rangle_{\text{sim}}$ indicates an average over jump events of the simulation. We thus obtain

$$P_{\text{corr}}(\Delta t) = P_{\text{sim}}(\Delta t)\left(\frac{1}{\langle t_{\text{tot}}/(t_{\text{tot}}-\Delta t) \rangle_{\text{sim}}}\right)^{-1}.$$  

Similarly the average times $\langle \Delta t \rangle_{\text{sim}}$ may be approximately corrected by

$$\langle \Delta t \rangle_{\text{corr}} = \left(\frac{\Delta t}{\langle t_{\text{tot}}/(t_{\text{tot}}-\Delta t) \rangle_{\text{sim}}}\right)^{-1}. \tag{7}$$

Figures 9(a) and 9(b) show the resulting corrected times $\langle \Delta t \rangle_{\text{corr}}$ for irreversible and reversible jumps, respectively. For simplification of notation we drop $\langle \cdot \rangle_{\text{corr}}$ in the following. The tail and head times, $\Delta t_{\text{tail}}$ and $\Delta t_{\text{head}}$, reveal that we find aging, since $\Delta t_{\text{tail}} > \Delta t_{\text{head}}$, i.e., jumps are more likely to occur at the beginning of the simulation than later. This particular aging effect decreases with increasing temperature. $\Delta t_{\text{tail}}$ and $\Delta t_{\text{head}}$ are both larger than $\Delta t_b$ because $\Delta t_b$ includes only times of particles which jump multiple times whereas $\Delta t_{\text{head}}$ and $\Delta t_{\text{tail}}$ include also times of particles which jump only once. Interestingly, $\Delta t_b$ is (although approximately corrected) still independent of temperature. Together with our results of Secs. III–V we therefore find that with increasing temperature, relaxation processes are accelerated via more jumping particles and more multiple jumps, but the times between multiple jumps do not become shorter. This is contrary to previous results in which waiting times decrease with increasing temperature. In our system, however, the temperature independence seems to be true not only for the averages but even for the distribution $P_{\text{corr}}(\Delta t_b)$ (see Fig. 10). We interpret this temperature independence as being due to aging which is consistent with our data for $\Delta t_{\text{tail}}$ and $\Delta t_{\text{head}}$ and with the results of Doliwa and Heuer who find that the distribution of waiting times initially is temperature independent and becomes temperature dependent at later times. As shown in Fig. 10 for irreversible jumps of $A$ and $B$ particles, $P(\Delta t_b)$ approximately follows a power law $P(\Delta t_b) \propto \Delta t_b^{-\nu}$ with $\nu = 0.84$. Even though we expect the system to show normal diffusion on a time scale much larger than our simulation run, the intermediate dynamics is subdiffusive. Please notice however, that
a direct comparison of our result with the scheme presented in Fig. 1 of Ref. 76 is not possible because in our case the multiple jumps of one particle at different times as well as the jumps of different particles are not independent, which changes the dynamics.\textsuperscript{77}

VII. JUMP LENGTHS

In the preceding sections we have investigated how many particles jump and how often jumps occur. Next we study how far these particles jump. To be able to test if our qualitative results are dependent on the definition of jump length, we use three quantities: $\Delta R_{i\text{f}}$, $\Delta R_{\text{max}}$, and $\Delta R_{\text{avg}}$ as sketched in Fig. 11. For a jump of particle $i$ the jump distance of the jump starting at $t_i$ and ending at $t_f$ (for the definition of $t_i$ and $t_f$ see Ref. 74) is

$$\Delta R_{i\text{f}} = |\vec{r}(t_f) - \vec{r}(t_i)|$$

and the maximal distance being detected during the jump is

$$\Delta R_{\text{max}} = \max_t |\vec{r}(t) - \vec{r}(t - 3200)|$$

(3200=4×20×2000×0.02 as in Ref. 70 and Sec. III), where $t$ is varied over times $t$ which satisfy $t_{i,\text{detect}} < t < t_{i,\text{detect}}.\textsuperscript{74}$ The third length $\Delta R_{\text{avg}}$ is less dependent on fluctuations than $\Delta R_{i\text{f}}$ and $\Delta R_{\text{max}}$. It is the distance of the overall average positions before and after the jump

$$\Delta R_{\text{avg}} = |\langle \vec{r} \rangle_t - \langle \vec{r} \rangle_t|.$$  

For the averages $\langle \cdot \rangle_t$ and $\langle \cdot \rangle_i$ the positions $\vec{r}(t)$ are averaged over times before and after the jump excluding a broadened time window around the jump $[t_i-800,t_i+800]$. Figure 12(a) shows that the different jump lengths have the same qualitative behavior, namely, as one might expect, an increase with increasing temperature $T$. We therefore show in the following mainly results for only $\Delta R_{\text{avg}}$ but find very similar behavior in the case of $\Delta R_{i\text{f}}$ and $\Delta R_{\text{max}}$. An increase in jump size with increasing temperature has been seen in previous simulations.\textsuperscript{3,19,35,72,73} However, in the case of Refs. 3, 72, and 73 the jump size is for a jump in the MSD, i.e., an average over all particles, which includes two effects: an increasing number of jumping particles and an increase in

the jump size of single particle jumps. Our approach has the advantage of keeping these two effects separate. Furthermore, in our case the single particle jumps are larger than in Refs. 3, 35, 72, and 73.

Figure 12(b) is the same as Fig. 12(a) for $\Delta R_{\text{avg}}$ but broken down by both jump and particle type. The smaller $B$ particles are jumping further than the $A$ particles. Furthermore reversible jumps are shorter than the irreversible jumps, which is due to two features in the distribution $P(\Delta R_{\text{avg}})$. As illustrated in Fig. 13 both irreversible and reversible jumps have a peak at about 0.8 and 1.0 for $A$ and $B$ particles, respectively. These peaks are relatively similar for reversible and irreversible jumps, taking aside that for irreversible jumps this peak position is slightly shifted to the right and slightly broadened (partly due to multiple jumps which are not resolved in time). This similarity and the peak positions around unity are consistent with our picture that

\begin{figure}[h]
\centering
\includegraphics[width=0.8\textwidth]{figure11}
\caption{Fig. 11. Sketch for the definitions of $\Delta R_{i\text{f}}$, $\Delta R_{\text{max}}$, and $\Delta R_{\text{avg}}$.}
\end{figure}

\begin{figure}[h]
\centering
\includegraphics[width=0.8\textwidth]{figure12}
\caption{(a) The jump sizes $\Delta R_{i\text{f}}$, $\Delta R_{\text{max}}$, and $\Delta R_{\text{avg}}$ (sketched in Fig. 11 and defined in the text) as function of temperature $T$. The averages are over both irreversible and reversible jumps of both $A$ and $B$ particles. (b) The jump size $\Delta R_{\text{avg}}$ as function of temperature $T$ separately for irreversible and reversible jumps of $A$ and $B$ particles.}
\end{figure}
irreversible and reversible jumps start out the same, namely, with a particle jumping out of its cage.

The increase of $\Delta R_{\text{avg}}$ with increasing temperature is a consequence of both a shift of the peak position and a broadening of the distribution of jump sizes. This is illustrated in the inset of Fig. 13 for irreversible jumps of $A$ particles and we find similar distributions for irreversible jumps of $B$ particles. At the largest investigated temperature $T=0.43$ some of the particles move as far as four-particle spacings. Some of these large jumps might correspond to multiple jumps of a single particle during single events but are also oscillating within their cages with a larger than average amplitude.

For irreversible jumps of size unity and larger we find that the distribution for all temperatures follow roughly a power law $P(\Delta R)\sim \Delta R^{-\mu}$ (see Fig. 14) with $\mu=6.3$ for $A$ and $\mu=4.8$ for $B$ particles. This is in accordance with sub-diffusive behavior for intermediate times as presented in Sec. VI for the distribution of waiting times.$^{26}$

As commented on earlier (see Secs. I and III), our definition for a jump is not based on a specified size but instead a multiple of the fluctuations of the particle. Let us therefore next look at the fluctuations. Figure 15 shows a comparison of the fluctuations of jumping particles $\sigma_{\text{jump}}$ and of nonjumping particles $\sigma_{\text{nojump}}$. Also included in Fig. 15 is an average over all particles $\sigma_{\text{est}}$ of the estimates for fluctuations $\sigma_{\text{est}}$ as they have been used for the jump identification (see Sec. III and Ref. 70). $\sigma_{\text{est}}$ is very similar to $\sigma_{\text{nojump}}$ and represents the fluctuations of an average particle. For the average of $\sigma_{\text{jump}}$ we exclude fluctuations during the jump: for jumping particle $i$ we take

$$\sigma_{\text{jump}}^i = \left( \frac{\langle \sigma_i^2 \rangle + \langle \sigma_i^2 \rangle}{2} \right)^{1/2},$$

where $\sigma_i$ and $\langle \cdot \rangle_i$ are defined as in Ref. 70 and earlier in this section, respectively. Although jump times are excluded, the fluctuations of jumping particles are clearly larger than the fluctuations of nonjumping particles. This means that jumping particles are not only moving farther than an average particle during single events but are also oscillating within their cages with a larger than average amplitude.

Figure 16(a) shows an average of jump size relative to the particle’s fluctuation $\Delta R_{\text{avg}}/\sigma_{\text{jump}}$ for jumping particles $i$. $\Delta R_{\text{avg}}/\sigma_{\text{jump}}$ seems independent of temperature for reversible jumps and for irreversible jumps slightly increasing with increasing temperature.

The distribution of $\Delta R/\sigma$ gives us the opportunity to estimate the influence of the cutoff in our definition of a jump. Similar to our approach for the search of a jump, we use $\sigma_{\text{est}}$, and obtain $P(\Delta R_{\text{avg}}/\sigma_{\text{est}})$ as shown in Fig. 16(b) for irreversible jumps of $A$ particles. We find a peak at $(\Delta R_{\text{avg}}/\sigma_{\text{est}}) = 5.5$ (for $B$ particles = 6.5), i.e., larger than our cutoff at $(\Delta R/\sigma) = \sqrt{20} = 4.5$. This gives us the hope that we have included the major contribution of jumps consistent
The single particle potential energy of particle $i$,$\Delta E_i$ is defined in Eq. (2). To obtain in addition the energies of the inherent structures we minimized the configurations $\{\mathbf{r}_i(t)\}$ via a steepest decent procedure. With the thus obtained configuration $\{\mathbf{r}_i^0(t)\}$ we determine

$$E^0(t) = \frac{1}{N} \sum_{i=1}^{N} \sum_{j \neq i} V_{ab}[r_{ij}^0(t)].$$

(14)

and

$$E_i^0(t) = \sum_{j \neq i} V_{ab}[r_{ij}^0(t)].$$

(15)

Similar to the jump size in position (see Fig. 17 and Sec. VII) we then determine

$$\Delta E_{av} = \langle E(t) \rangle_i - \langle E(t) \rangle_i,$$

(16)

and

$$\Delta E_{if} = E(t_f) - E(t_i).$$

(17)

and similarly for the energies of Eqs. (13)–(15).

We find in most of our analysis that $\Delta E_{av}$ and $\Delta E_{if}$ (and all other equivalents for $E^0, E_i$, and $E_i^0$) are showing the same qualitative behavior, and therefore present in the following results for $\Delta E_{av}$ using Eq. (16) and Eqs (12)–(15).

Figure 18(a) shows $\Delta E_{av}$ as a function of temperature averaged separately over irreversible and reversible jumps of $A$ and $B$ particles. Notice that the energy jumps are small compared to energy values $\langle E \rangle_{i,\alpha}^\infty - 7$. The irreversible jumps lower the energies more than reversible jumps and more so for $A$ than $B$ particles, because a jump of a larger $A$ particle results in a bigger change in environment than for a smaller $B$ particle. The lowering of the total potential energy happens mostly at intermediate temperatures and less at the lowest and highest investigated temperatures. This might be an aging effect: at low temperatures the system is basically frozen in, at intermediate temperatures the system has time to partially age, and at higher temperatures the system has already aged before the production run starts. This effect is lost when we look at the jumps in single particle energies $\Delta E_{i,av}(T)$ which is basically zero for all jumps. We therefore find that the total potential energies $\Delta E_{av}$, $\Delta E_{if}$, $\Delta E_0$, and $\Delta E_{if}$ show a more systematic dependence on temperature than their single particle equivalents, $\Delta E_{i,av}$, etc. For the absolute values $|\Delta E_{av}|$, etc., and all following.

FIG. 17. Sketch for the definitions of $\langle E \rangle_i$ and $\langle E \rangle_i.$
quantities, however, we find the opposite, i.e., the single particle equivalents $u_{DEi,avg}$, etc., show a more pronounced behavior. We therefore show from now on the results which use Eqs. (13) and (15) only.

We next look at the absolute value $|\Delta E_{i,avg}^0|$ as a function of temperature [see Fig. 18(b)]. As one might expect, the irreversible jumps show larger changes in energy than reversible jumps due to larger jumps in position [Fig. 12(b)] and therefore more change in the environment of the jumping particle. And similarly the larger $A$ particles in comparison with the smaller $B$ particles experience a larger change in energy because $A$ particles are surrounded by more neighbors, i.e., a larger environment. With increasing temperature the absolute value of the jump in energy [Fig. 18(b)] is increasing (with the one exception of irreversible $B$ particles) which is consistent with the increase in jump position [Fig. 12(b)].

We define the fluctuations in energy $\sigma_{E_i}$ at time $t=m \times 2000 \times 0.02$ similar to the fluctuations in position,$^{70}$ where $\bar{E}_i(m')$ is the time average $\frac{1}{20} \sum_{m'=m-9}^{m+10} E_i(m')$. The energy fluctuations are also increasing with increasing temperature and are larger for $A$ than $B$ particles [see Fig. 19(a)]. The fluctuations of irreversible and reversible jumps are, however, very similar, which is consistent with the picture that reversible jumps are “failed” irreversible jumps and in that sense start out the same way as irreversible jumps.

Figure 19(b) is a comparison of energy fluctuations of jumping (both irreversible and reversible and of both $A$ and $B$) particles $\sigma_{E_{i,jump}}$ and of nonjumping particles $\sigma_{E_{i,\text{nojump}}}$.
In \( \sigma_{E_{\text{jump}}} \) we exclude the times during the jump by using the time windows \( \langle \cdot \rangle_t \). We find that the energy fluctuations of jumping particles are larger than for nonjumping particles which is consistent with the corresponding fluctuations in position (see Fig. 15).

By normalizing \( |\Delta E_{i,\text{avg}}|/\sigma_{E_i} \) with the fluctuations \( \sigma_{E_i} \) we obtain Fig. 20(a). Similar to the case of normalized jumps in position [Fig. 16(a)] we find that \( |\Delta E_{i,\text{avg}}|/\sigma_{E_i} \) is for reversible jumps basically independent of temperature. For irreversible jumps, however, \( |\Delta E_{i,\text{avg}}|/\sigma_{E_i} \) decreases significantly with increasing temperature. This indicates that with increasing temperature the jumps are increasingly more driven by fluctuations. Contrary to the equivalent in position [Fig. 16(a)], the normalized energies are significantly smaller, in the range 0.5–1.4 rather than 3–7, and the distribution of \( |\Delta E_{i,\text{avg}}|/\sigma_{E_i} \) [see Fig. 20(b)] is monotonous decreasing with a less far reaching tail than the equivalent in position [Fig. 16(b)].

**IX. DEPENDENCE ON HISTORY**

We are presenting in this paper out of equilibrium results and have been finding aging effects. The question arises how the history of the simulation influences the particle jump statistics. In this section we address this question by comparing results of simulation runs with three different histories.

The data presented so far were initially equilibrated at \( T = 0.446 \) and were then quenched to the investigated temperatures \( (T = 0.15, 0.20, 0.25, 0.30, 0.35, 0.38, 0.40, 0.41, 0.42, 0.43) \) according to the procedure described in Sec. II. To investigate the effect of aging we performed a second set of simulations for the quenches from \( T = 0.446 \) to \( T = 0.15, 0.25, 0.35, 0.40, 0.43 \). The history of ten independent configurations is the same as before but the (NVE) production runs are instead of for \( 5 \times 10^5 \) MD steps for \( 3 \times 10^7 \) MD steps. We analyze the last \( 5 \times 10^6 \) MD steps, and refer to these results as \( “T_{\text{init}} = 0.446 \text{(aged)}” \) and to the previously presented data as \( “T_{\text{init}} = 0.446 \text{(not aged)}.” \) A third set of simulations is performed to study the influence of the starting temperature from which the system is quenched. To obtain ten independent configurations with starting temperature \( T = 0.5 \) we equilibrate at \( T = 3.0 \) and choose configurations at least \( 5 \times 10^3 \) time units apart. We then cool linearly in time \( t \) \( (T = T_0 - \gamma t) \) from \( T_0 = 3.0 \) to \( T = 0.5 \) with \( \gamma = 1.25 \times 10^{-5} \) (with \( \Delta t = 0.02 \) and therefore for \( 10^7 \) MD steps) at \( T = 0.5 \) we equilibrate for \( 10^7 \) MD steps with \( \Delta t = 0.02 \). We then quench the system to the investigated temperatures \( T = 0.15, 0.20, 0.25, 0.30, 0.35, 0.38, 0.40, 0.41, 0.42, 0.43 \) by running (NVT) simulations for \( 10^5 \) MD steps with \( \Delta t = 0.02 [T_{\text{init}} = 0.5 \text{(not aged)}] \) and consecutive (NVE) production runs for \( 5 \times 10^5 \) MD steps with \( \Delta t = 0.02 \). The history of this simulation set therefore differs from the nonaged \( T_{\text{init}} = 0.446 \) set both in the initial temperature and also in the direct quench to the investigated temperatures instead of the stepping procedure described in Sec. II.

As illustrated in Fig. 21 the number of jumping particles normalized by the number of particles for \( T_{\text{init}} = 0.446 \text{(aged)} \) and for \( T_{\text{init}} = 0.5 \text{(not aged)} \) have qualitatively the same tem-
perature dependence as before (see Fig. 3). As one might expect we find, however, quantitatively a dependence on the history (see Fig. 22). Directly after the quench we find more jumps than after a waiting time of \(5 \times 10^5\) time units. The difference becomes less for temperatures \(T \lesssim 0.41\). Quenches from the higher initial temperature \(T = 0.5\) result for all investigated temperatures in more jumps.

Independent of the history, an increasing fraction of these particles are irreversible jumpers with increasing temperature (see Fig. 23). This increase is less pronounced for the \(T_{\text{ini}} = 0.5\) simulations and steeper for the aged \(T_{\text{ini}} = 0.446\) data. This dependence on the history is again vanishing at high temperatures.

Let us next investigate the quantity which mostly raised the question of history dependence: \(\Delta t_b\), the time between jumps. According to Fig. 24 we find that \(\Delta t_b\) is for \(T \lesssim 0.3\) not only temperature independent but also independent of the initial temperature and the aging time. However, this dependence on the waiting time needs further investigation.\(^{80}\)

As shown in Fig. 25 we also find an independence of history for the jump length \(\Delta R_{\text{avg}}\) which might indicate that the jumps are not only characteristic for the relaxation process with one specific history but might be a more general feature.

X. CONCLUSIONS

We study the dynamics of a binary Lennard-Jones system below the glass transition. Our focus is on jump processes, which we identify via single particle trajectories. Two kinds of jumps are found: “reversible jumps,” where a particle jumps back and forth between one or more average positions, and “irreversible jumps,” where a particle does not return to any of its former average positions, i.e. success-

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**FIG. 22.** A comparison of the number of jumping particles normalized by the number of particles for three simulation sets with different histories. We include here both irreversible and reversible jumps of both A and B particles.

**FIG. 23.** Same as Fig. 4 but for both A and B particles and for simulation sets with three different histories.

**FIG. 24.** Same as Fig. 9(a) and (b) (in the inset) but for only \(\Delta t_b\) and for simulations with three different histories.

**FIG. 25.** The jump size \(\Delta R_{\text{avg}}\) of both A and B particles as a function of temperature for simulations with three different histories.
fully escapes its cage of neighbors. Both irreversible and reversible jumps of A and B particles occur at all temperatures. With increasing temperature more particles jump, more average positions are visited, the jump size both in position and in the absolute value of the potential energies $|\Delta E|, |\Delta E^0|, |\Delta E_i|$, and $|\Delta E^0_i|$ (total and single particle, not minimized and minimized) increases, and the fluctuations in position $\sigma_R$ and potential energy $\sigma_E$ increase. The fluctuations are larger for jumping particles than for nonjumping particles even if fluctuations during the jump are not part of the average, which indicates that jumping particles are not only during jump times more mobile than nonjumping particles. The ratio $|\Delta E_i|/\sigma_{E_i}$ of irreversible jumps decreases with increasing temperature. This confirms a commonly used assumption that with increasing temperature the irreversible jumps become more driven by fluctuations.

With increasing temperature also irreversible jumps become proportionally more frequent than reversible jumps. We interpret this such that irreversible and reversible jumps are similar in that sense that a particle tries to escape its cage. In the case of reversible jumps the particle finds its way back into the cage whereas in the case of an irreversible jump the path back into the cage becomes blocked due to rearrangements of the cage. At larger temperature these rearrangements of the cage become more likely (since, for example, the fluctuations increase) and therefore irreversible jumps occur more often. Irreversible and reversible jumps show in most quantities qualitatively the same behavior, such as their temperature dependence of jump size in position and energy, and differ only in size.

The most surprising result of our work is that the times between successive jumps are independent of temperature. This is most likely due to aging, which could mean on the time scale of our simulation that with increasing temperature the irreversible jumps with increasing temperature, since this indicates that not only the jumping particle itself but also its cage are dependent on temperature. There might, however, be also collective jumps, where each particle jumps a too small amount to be detected by our search algorithm. It remains to be studied of how much importance such processes are in our system.

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2 Examples of the mean square displacement as a function of time (logarithmically plotted) are given in W. Kob and H. C. Andersen, Phys. Rev. E 51, 4626 (1995).


E. Leutheusser, Phys. Rev. A 54, 3049 (1996). Details can be found in Refs. 65 and 69.


The ten initial configurations differ drastically in their thermal history and are therefore completely independent. They were started at \( T = 5.0 \) and, after various cooling and reheating periods, were equilibrated at \( T = 0.464 \). Details can be found in Refs. 65 and 69.

T. Gleim and W. Kob, Phys. Rev. Lett. 81, 4404 (1998). As described in Sec. II, we start with the periodically (every 2000 MD steps of size step \( \tau = 0.02 \)) stored configurations \( \{ r_i(t) \} \) where \( t = m \times 2000 \times 0.02 \). For the single particle fluctuations we use the average \( \bar{r}_i(t) = \frac{1}{n} \sum_{j=1}^{n} r_j(t) \), their squares \( \bar{r}_i^2(t) = \frac{1}{n} \sum_{j=1}^{n} r_j^2(t) \), and \( \sigma_i^2(t) = \bar{r}_i^2(t) - \bar{r}_i(t)^2 \) at times \( m = 10, 30, \ldots \). We first make a preliminary estimate of the fluctuations \( \sigma_{\text{pre}} = \sqrt{n \bar{r}_i^2(t)} \) and then obtain the final estimate for the fluctuations \( \sigma_i(t) \) by redrawing the average over \( \sigma_i^2(t) \) but by averaging only over all \( \sigma_i^2(t) \) for which \( \sigma_i^2(t) < 5 \bar{r}_i^2(t) \) to roughly exclude jumps from the average.

We define \( r_i(t) \) to be a spike if \( r_i^2(t) - \bar{r}_i^2(t) > \kappa \sigma_i^2(t) \), \( \kappa \) is defined as in Sec. VII. Similar to case 1 we say that the spike \( r_i(t) \) returns to the previous average position if \( r_i^2(t) - \bar{r}_i^2(t) < \kappa \sigma_i^2(t) \).


To determine the starting time \( t_i \) and ending time \( t_f \) of a jump we use an approach similar to our definition for the occurrence of a jump [Eq. (4)]. We compare the differences in averaged positions \( \Delta r_i(t) \) and the fluctuations \( \sigma_{\text{pre}}^2 \) [see Ref. 70]. If a jump has been detected at time \( t_{\text{detect}} \) (because \( \Delta r_i^2(t_{\text{detect}}) > 2 \sigma_i^2(t_{\text{detect}}) \)) then the starting time \( t_i \) is the largest time \( t \) for which \( \Delta r_i^2(t) > 5 \sigma_i^2(t) \) within the boundaries of \( t_{\text{lowest}} < t_i < t_{\text{highest}}, \) where in the case of a first jump \( t_{\text{lowest}} \) corresponds to the beginning of the production run and in the case of the existence of a previous jump with \( t_i^p \) then \( t_{\text{lowest}} < t_i^p < t_i \). Similarly after a jump has started it is detected to be finished at the earliest time \( t_{\text{detect}} \) for which \( \Delta r_i^2(t_{\text{detect}}) \leq 5 \sigma_i^2(t_{\text{detect}}) \). The finishing time of the jump \( t_f \) is then the maximum of \( t_{\text{detect}} \) and \( t_{\text{lowest}} = 3200 \) where we subtract \( 3200 = (4 \times 20 \times 2000 \times 0.02) \) because \( \Delta T = \bar{T}_f(t_f) - \bar{T}(t_f - 3200) \) [see Eq. (3) and Ref. 70]. If the condition for \( t_{\text{detect}} \) never occurs, then \( t_i \) is equal to the length of the simulation run. To determine \( \Delta r_i \) in case 2 of Fig. 2, we use for the first spike the time between the previous \( t_i \) and the time of the spike and for successive spikes their time differences.


In Figs. 19(a) and 19(b) we use the time average windows \( \langle \cdot \rangle_1 \). In Fig. 19(a) we then average over jumps of a certain kind. In the case of a particle with multiple jumps this means that the times after the first jump and before the last jump are counted twice. In Fig. 19(b), however, we compare the behavior of particles (instead of jump events) and therefore count these time windows only once.