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Communication: Shifted forces in molecular dynamics

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Simulations involving the Lennard-Jones potential usually employ a cutoff at $r = 2.5\sigma$. This communication investigates the possibility of reducing the cutoff. Two different cutoff implementations are compared, the standard shifted potential cutoff and the less commonly used shifted forces cutoff. The first has correct forces below the cutoff, whereas the shifted forces cutoff modifies Newton’s equations at all distances. The latter is nevertheless superior; we find that for most purposes realistic simulations may be obtained using a shifted forces cutoff at $r = 1.5\sigma$, even though the pair force is here 30 times larger than at $r = 2.5\sigma$. © 2011 American Institute of Physics. [doi:10.1063/1.3558787]

Molecular dynamics (MD) simulations solve Newton’s equations of motion by discretizing the time coordinate. The time-consuming part of a simulation is the force calculation. For a system of N particles this is an $O(N^2)$ process whenever all particles interact. In practice the interactions are negligible at long distances, however, and for this reason one always introduces a cutoff at some interparticle distance $r = r_c$ beyond which interactions are ignored.¹

The standard Lennard-Jones (LJ) pair potential is given by

$$u_{\text{LJ}}(r) = 4\epsilon \left[\left(\frac{\sigma}{r} \right)^{12} - \left(\frac{\sigma}{r} \right)^6 \right]. \quad (1)$$

Usually, a cutoff at $r_c = 2.5\sigma$ is employed; at this point the potential is merely 1.6% of its value at the minimum ($-\epsilon$). Although a cutoff makes the force calculation an $O(N)$ process, this calculation remains the most demanding in terms of computer time.

The present communication investigates the possibility of reducing the LJ cutoff below 2.5σ without compromising accuracy to any significant extent. Before presenting evidence that this is possible, it is important to recall that quantities depending explicitly on the free energy are generally quite sensitive to how large is the cutoff. Examples include the location of the critical point,² the surface tension,^{2,3} and the solid–liquid coexistence line.^{4,5} For such quantities even a cutoff at 2.5σ gives inaccurate results, and in some cases the cutoff must be larger than 6σ to get reliable results.³ Note, however, that if a simulation gives virtually correct particle distribution, the thermodynamics can be accurately calculated by first-order perturbation theory.⁶

We compared two cutoff implementations at varying cutoffs with the “true” LJ system, the latter being defined here by the cutoff $r_c = 4.5\sigma$. One cutoff is the standard “truncated and shifted potential” (SP for shifted potential), for which the radial force is given¹ by $[f_{\text{LJ}}(r) = -u'_{\text{LJ}}(r)$ is the LJ radial force]

$$f_{\text{SP}}(r) = \begin{cases} f_{\text{LJ}}(r) & \text{if } r < r_c \\ 0 & \text{if } r > r_c. \end{cases} \quad (2)$$

This is referred to as a SP cutoff because it corresponds to shifting the potential below the cutoff and putting it to zero above, which ensures continuity of the potential at r_c and avoids an infinite force here.

The “truncated and shifted forces” cutoff (SF for shifted forces)^{1,7} has the force go continuously to zero at r_c , which is obtained by subtracting a constant term:

$$f_{\text{SF}}(r) = \begin{cases} f_{\text{LJ}}(r) - f_{\text{LJ}}(r_c) & \text{if } r < r_c \\ 0 & \text{if } r > r_c. \end{cases} \quad (3)$$

This corresponds to the following modification of the potential: $u_{\text{SF}}(r) = u_{\text{LJ}}(r) - (r - r_c)u'_{\text{LJ}}(r_c) - u_{\text{LJ}}(r_c)$ for $r < r_c$, $u_{\text{SF}}(r) = 0$ for $r > r_c$. Use of a SF cutoff has recently become popular in connection with improved methods for simulating systems with Coulomb interactions.⁸

We simulated the standard single-component LJ liquid at the state point that in dimensionless units has density $\rho = 0.85$ and temperature $T = 1.0$.⁹ This is a typical moderate-pressure liquid state point.^{1,10} Other state points were also examined—including state points of the fcc crystal, at the liquid–gas interface, at the solid–liquid interface, and for a supercooled system—leading in all cases to the same overall conclusion. For this reason we report below results for just one state point of the LJ liquid and one of the Kob–Andersen binary LJ (KABLJ) liquid.¹¹ 2000 LJ particles were simulated using the standard central-difference constant temperature/energy (NVT/NVE) algorithms (Figs. 2, 3, and 4, 6, respectively); 1000 particles of the KABLJ liquid were simulated using the NVT algorithm (Fig. 5).

Figure 1 shows the basics of the LJ system. In the upper figure the black curve gives the LJ pair potential $u_{\text{LJ}}(r)$ and the black dashed curve the radial distribution function $g(r)$, which has its maximum close to u ’s minimum. In the lower figure the black (lower) curve shows the LJ pair force $f_{\text{LJ}}(r)$. The red (upper) curve gives $f_{\text{SF}}(r)$ when a cutoff at 1.5σ is introduced; note that the shifted force differs significantly from the true force.

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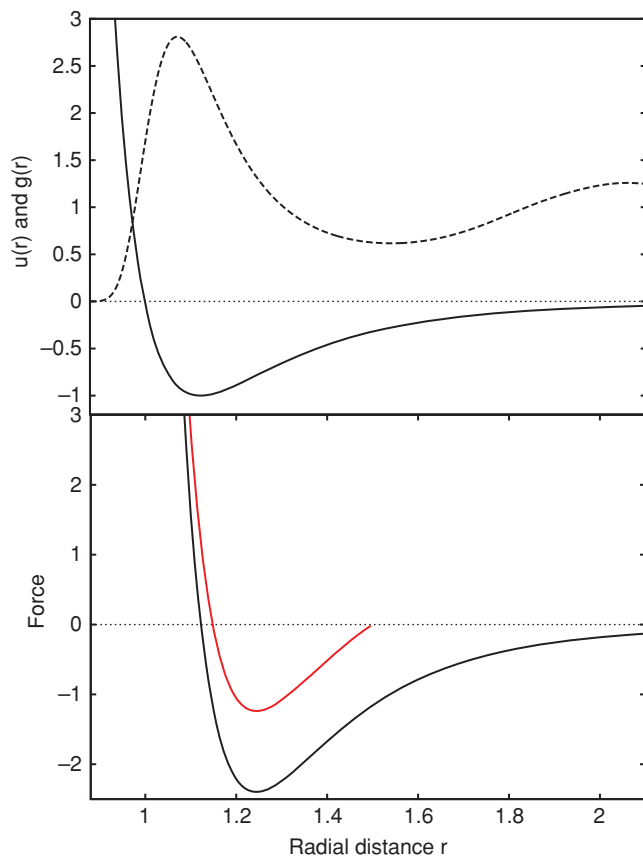


FIG. 1. (a) The Lennard-Jones potential (black full curve) and the radial distribution function $g(r)$ (black dashed curve) for a system at $\rho = 0.85$ and $T=1.00$ in dimensionless units. (b) The radial force $f_{LJ}(r) = -u'_{LJ}(r)$ (black). At $r = 1.5\sigma$ the force is 30 times larger than at $r = 2.5\sigma$. Also shown is the shifted force for a cutoff at 1.5σ (red, upper curve).

Figure 2 shows the true pair-distribution function (black) and the simulated $g(r)$ for three $r_c = 1.5\sigma$ cutoffs: SF (red, barely visible), SP (green, slightly lower at $r = 1.5\sigma$), and a smoothed SP cutoff ensuring the force and its first derivative go continuously to zero at the cutoff¹² [dashed (green) curve]. The curves deviate little, except near the cutoff where the smallest errors are found for a SF cutoff (inset).

In order to systematically compare the SP and SF cutoffs we studied the LJ liquid for a range of cutoffs. Figure 3 quantifies the difference between the computed $g(r)$ and the true radial distribution function, $g_0(r)$, by evaluating $\int_0^{4.5\sigma} |g(r) - g_0(r)| dr$. SF is the red (lower) curve, SP is the green (upper) curve. SF works better than SP for all values of r_c above the Weeks–Chandler–Andersen (WCA) cutoff at the potential energy minimum⁶ where SF = SP ($r_c = 2^{1/6}\sigma = 1.12\sigma$). Smoothing a SP cutoff has only a marginal effect compared to not smoothing it (results not shown). Applying first-order perturbation theory with the $g(r)$ obtained in a simulation with SF cutoff at $r_c = 1.5\sigma$ leads to a pressure that deviates only 1% from the correct value.

Figure 4 studies energy drift in long NVE simulations for $r_c = 1.5\sigma$. The SF cutoff (red, horizontal) exhibits no energy drift, whereas SP (green, diverging) does. Figure 4 also gives results when the force of a SP cutoff is smoothed¹² (green dashed, horizontal curve). This leads to much better energy conservation,¹ but the energy fluctuations are some-

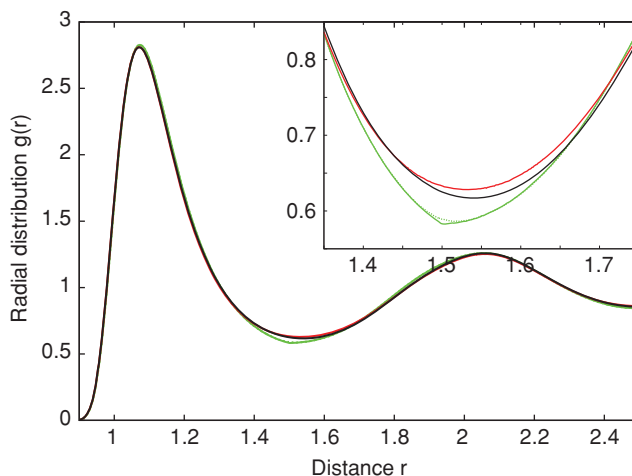


FIG. 2. Radial distribution function $g(r)$ for the “true” LJ system (black) and two cutoffs at $r_c = 1.5\sigma$. The red (barely visible) curve gives results for a SF cutoff, the green curve (slightly lower at $r = 1.5$) for a SP cutoff. The dashed (green) curve gives results for a SP cutoff with smoothing of the force and its derivative at the cutoff (Ref. 12); this, however, does not improve the SP results.

what larger than for a SF cutoff. The simulations indicate the existence of a hidden invariance in the central-difference algorithm for a continuous force function, deriving from a “shadow Hamiltonian.”¹³

Not only static quantities, but also the dynamics are affected little by replacing a 2.5σ SP cutoff with a 1.5σ SF cutoff. This is demonstrated in Fig. 5, which shows simulations of the incoherent intermediate scattering function of the supercooled KABLJ liquid.¹¹ For reference a WCA cutoff simulation is included (blue dashed curve, fastest relaxation), which was recently shown to be inaccurate despite the fact that the WCA radial distribution function is reasonably good for this system.¹⁴ A SP cutoff at $r_c = 1.5\sigma_{AA}$ gives too slow dynamics (purple dotted curve). Within the numerical uncertainties incoherent scattering functions are identical for

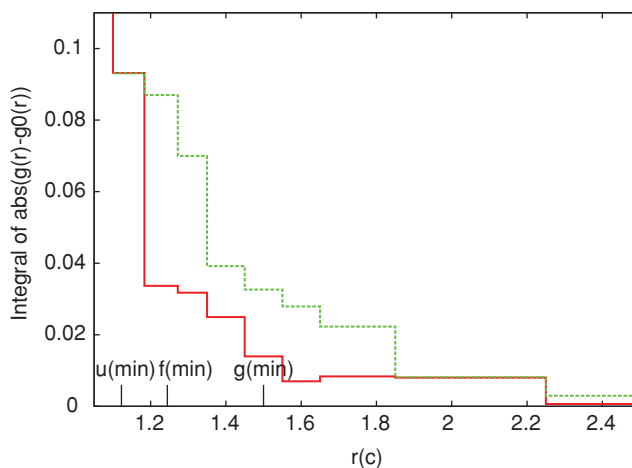


FIG. 3. Integrated numerical difference $\int_0^{4.5\sigma} |g(r) - g_0(r)| dr$ of the true radial distribution function, $g_0(r)$, and $g(r)$ for various cutoff distances r_c . The red (lower) curve gives results for the SF cutoff, the green (upper) for the SP cutoff. Smoothing a SP cutoff (Ref. 12) does not improve its accuracy (data not shown).

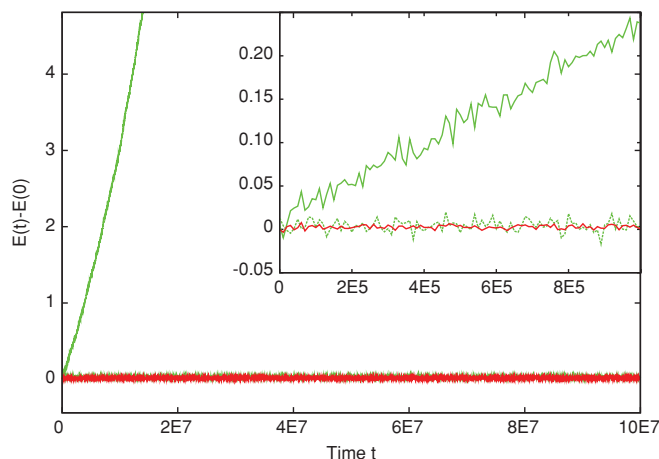


FIG. 4. Energy drift as a function of time for long simulations (10^8 time steps of length 0.005) with a cutoff at 1.5σ . The red (horizontal) curve gives results for the SF cutoff, the green (diverging) for the SP cutoff, and the green dashed (horizontal) curve for a smoothed SP cutoff. Smoothing a SP cutoff stabilizes the algorithm, but the fluctuations are still somewhat larger than for a SF cutoff. The inset shows the initial part of the simulation.

the “true” system, a SP cutoff at $r_c = 2.5\sigma_{AA}$, and a SF cutoff at $r_c = 1.5\sigma_{AA}$. Similar results were found for the single-component LJ liquid’s dynamics. We conclude that a SF cutoff at $r_c = 1.5\sigma$ generally works well for both statics and dynamics of LJ systems.

Why does a cutoff, for which the forces are modified at all distances (SF), work better than when the forces are correct below the cutoff (SP)? A SF cutoff modifies the pair force by adding a constant force for all distances below r_c ; at the same time SF ensures that the pair force goes continuously to zero at $r = r_c$. Apparently, ensuring continuity of the force—and thereby that $u''(r)$ does not spike artificially at the cutoff—is more important than maintaining the correct pair force below the cutoff. How large is the

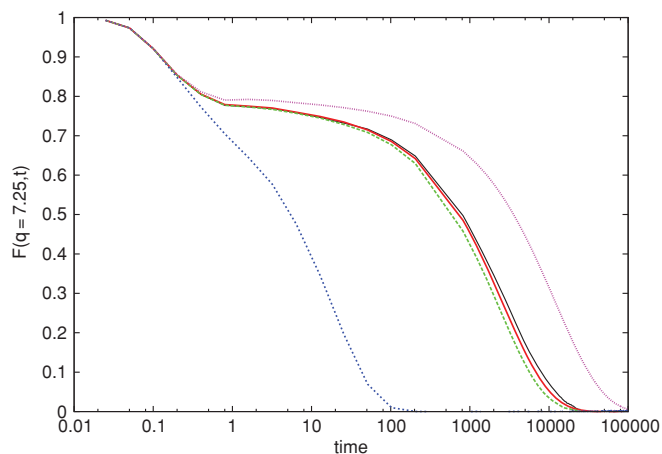
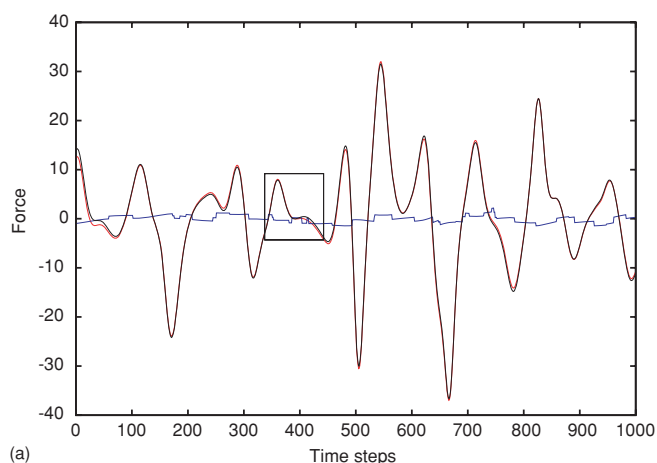
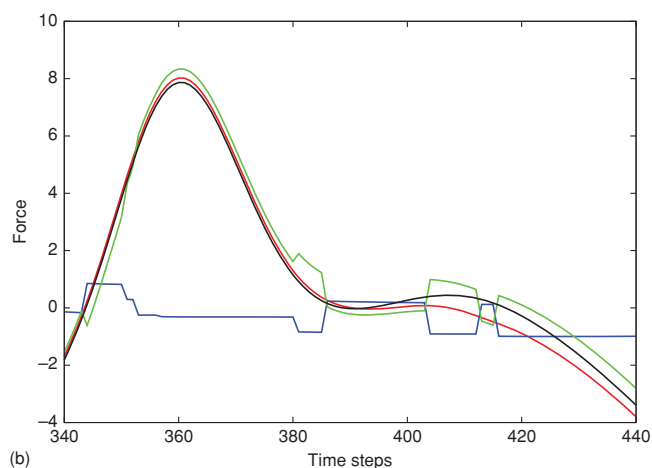


FIG. 5. The AA particle incoherent intermediate scattering function for the KABLIJ liquid in the highly viscous regime ($T = 0.45$, $\rho = 1.2$, $q = 7.25$). In order from the left the figure shows results for: a WCA cutoff (blue dashed curve), a SP cutoff at $2.5\sigma_{AA}$ (green dashed curve), a SF cutoff at $1.5\sigma_{AA}$ (red full curve), the “true” system (black full curve), and a SP cutoff at $1.5\sigma_{AA}$ (purple dotted curve). Smoothing a SP cutoff (Ref. 12) does not improve its accuracy (data not shown).



(a)



(b)

FIG. 6. (a) The x component of the force on a typical particle during 1000 time steps. The black curve gives the true force, the red curve (barely visible on top of the black curve) the force for a SF cutoff at 1.5σ . The blue curve (fluctuating around zero) gives the sum of the x coordinates of the constant “shift” terms of Eq. (3). (b) Details after 340 steps. The green (upper) curve gives the SP force ($r_c = 1.5\sigma$). Only true and SF forces are smooth functions of time.

change induced by the added constant force of the SF cutoff? Figure 6 shows the x component of the force on a typical particle as a function of time ($r_c = 1.5\sigma$). The black curve gives the true force, the red, barely visible curve the SF force, and the blue, fluctuating curve the SF correction term. Although the true and SF individual pair forces differ significantly (Fig. 1), the difference between true and SF total forces is small and stochastic ($\sim 3\%$). This reflects an almost cancellation of the correction terms deriving from the fact that the nearest neighbors are more or less uniformly spread around the particle in question. It was recently discussed why adding a linear term ($\propto r$) to a pair potential hardly affects dynamics¹⁵ and statistical mechanics.¹⁶ For a given particle’s interactions with its neighbors the linear terms sum to almost a constant because, if the particle is moved, some nearest-neighbor distances increase and others decrease, and their sum is almost unaffected.

Figure 6(b) shows details of Fig. 6(a); we here added the SP force for the same cutoff (green, upper curve). Both SP and correction terms are discontinuous; they jump whenever

a particle pair distance passes the cutoff. Altogether, Fig. 6 shows that not only does the sum of the constant forces on a given particle from its neighbors cancel to a high degree, so do the interactions with particles beyond the cutoff. The result is that the particle distribution is affected little by the long-range attractive forces, a fact that lies behind the success of perturbation theory.^{6,17,18}

In summary, when a SF cutoff is used instead of the standard SP cutoff, errors are significantly reduced. Our simulations suggest that a SF cutoff at 1.5σ may be used whenever the standard SP cutoff at 2.5σ gives reliable results; this applies even though the pair force at $r = 1.5\sigma$ is 30 times larger than at $r = 2.5\sigma$. A cutoff at 1.5σ is large enough to ensure that all interactions within the first coordination shell are taken into account (Fig. 2). Use of a 1.5σ SF cutoff instead of a SP cutoff at 2.5σ leads potentially to a factor of $(2.5/1.5)^3 = 4.7$ shorter simulation time for LJ systems.

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¹M. P. Allen and D. J. Tildesley, *Computer Simulation of Liquids* (Oxford Science Publications, Oxford, 1987); D. Frenkel and B. Smit, *Understanding Molecular Simulation* (Academic, New York, 2002).

²B. Smit, *J. Chem. Phys.* **96**, 8639 (1992).

³P. Grosfils and J. F. Lutsko, *J. Chem. Phys.* **130**, 054703 (2009).

⁴C. Valeriani, Z. J. Wang, and D. Frenkel, *Mol. Simul.* **33**, 1023 (2007).

⁵A. Ahmed and R. J. Sadus, *J. Chem. Phys.* **133**, 124515 (2010).

⁶J. D. Weeks, D. Chandler, and H. C. Andersen, *J. Chem. Phys.* **54**, 5237 (1971); J. H. R. Clarke, W. Smith, and L. V. Woodcock, *ibid.* **84**, 2290 (1986); J. D. Weeks, K. Vollmayr, and K. Katsov, *Physica A* **244**, 461 (1997); F. Cuadros, A. Mulero, and C. A. Faundez, *Mol. Phys.* **98**, 899 (2000).

⁷S. D. Stoddard and J. Ford, *Phys. Rev. A* **8**, 1504 (1973); J. J. Nicolas, K. E. Gubbins, W. B. Street, and D. J. Tildesley, *Mol. Phys.* **37**, 1429 (1979); J. G. Powles, W. A. B. Evans, and N. Quirke, *ibid.* **46**, 1347 (1982).

⁸D. Wolf, P. Keblinski, S. R. Phillpot, and J. Eggebrecht, *J. Chem. Phys.* **110**, 8254 (1999); D. Zahn, B. Schilling, and S. M. Kast, *J. Phys. Chem. B* **106**, 10725 (2002); C. J. Fennell and J. D. Gezelter, *J. Chem. Phys.* **124**, 234104 (2006).

⁹For MD details see S. Toxvaerd, *Mol. Phys.* **72**, 159 (1991). The unit length, energy, and time used are, respectively, σ , ϵ , and $\sigma\sqrt{m/\epsilon}$.

¹⁰L. Verlet, *Phys. Rev.* **159**, 98 (1967).

¹¹W. Kob and H. C. Andersen, *Phys. Rev. Lett.* **73**, 1376 (1994).

¹²The radial force $f_{\text{SP}}(r)$ was smoothed in the interval $r_c \pm \delta$ by replacing it with the function $f(r) = f(r_c - \delta)h(x)$ with $h(x) = 1 + cx - (3 + 2c)x^2 + (2 + c)x^3$, where $x = (r - r_c + \delta)/2\delta$ and $c = 2\delta u''(r_c - \delta)/u'(r_c - \delta)$. This ensures that f_{SP} and f'_{SP} go smoothly to zero in the cut interval. In the simulations $\delta = 0.025$ (other values lead to similar conclusions).

¹³S. Toxvaerd, *Phys. Rev. E* **50**, 2271 (1994).

¹⁴L. Berthier and G. Tarjus, *Phys. Rev. Lett.* **103**, 170601 (2009); U. R. Pedersen, T. B. Schröder, and J. C. Dyre, *ibid.* **105**, 157801 (2010); Y. S. Elmatad, D. Chandler, and J. P. Garrahan, *J. Phys. Chem. B* **114**, 17113 (2010).

¹⁵N. P. Bailey, U. R. Pedersen, N. Gnan, T. B. Schröder, and J. C. Dyre, *J. Chem. Phys.* **129**, 184508 (2008); T. B. Schröder, N. P. Bailey, U. R. Pedersen, N. Gnan, and J. C. Dyre, *ibid.* **131**, 234503 (2009).

¹⁶N. Gnan, T. B. Schröder, U. R. Pedersen, N. P. Bailey, and J. C. Dyre, *J. Chem. Phys.* **131**, 234504 (2009).

¹⁷J. A. Barker and D. Henderson, *J. Chem. Phys.* **47**, 4714 (1967).

¹⁸S. Toxvaerd, *J. Chem. Phys.* **55**, 3116 (1971).