An observation of a nascent fractal pattern in MD simulation for a fragmentation of an fcc lattice

Shinpei Chikazumi\textsuperscript{a,c,1} Akira Iwamoto\textsuperscript{b}

\textsuperscript{a}Institute of Physics, University of Tsukuba, Tsukuba, Ibaraki 305-0006, Japan
\textsuperscript{b}Japan Atomic Energy Research Institute, Tokai, Ibaraki, 319-1195 Japan
\textsuperscript{c}Department of Materials Science, Japan Atomic Energy Research Institute, Tokai, Ibaraki, 319-1195 Japan

Abstract

To seek for a possible origin of fractal pattern in nature, we perform a molecular dynamics simulation for a fragmentation of an infinite fcc lattice. The fragmentation is induced by the initial condition of the model that the lattice particles have the Hubble-type radial expansion velocities. As time proceeds, the average density decreases and density fluctuation develops. By using the box counting method, it is found that the frequency-size plot of the density follows instantaneously a universal power-law for each Hubble constant up to the size of a cross-over. This cross-over size corresponds to the maximum size of fluctuation and is found to obey a dynamical scaling law as a function of time. This instantaneous generation of a nascent fractal is purely of dynamical origin and it shows us a new formation mechanism of a fractal patterns different from the traditional criticality concept.

Key words: fractal, power law, molecular dynamics, fragmentation

PACS: 05.45.Df, 05.65.+b, 62.20.Mk

Since the pioneering work of Mandelbrot \cite{1} a large effort has been directed towards understanding the ubiquitous manifestation of fractal structure in nature. Since fractal structure is intimately related to a scaling property, a customary way of identifying the fractal is to check if the power-law relation of the frequency as functions of the size holds. Malcai et al. \cite{2} in this way have listed the fractal dimensions and the corresponding scaling ranges for a variety of physical systems. They pointed out that the finiteness of the fractal...
range is a typical feature for physical systems which is in sharp contrast with an ideal mathematical concept of fractal.

Let us therefore settle our target as “why and how a power-law relation in frequency-size relation appears frequently in nature no matter what its range”. It seems natural to adopt some fractal growth models to challenge this subject or to adopt some kinds of cellular-automata approaches. In these models, the power-law in frequency-size relation is established at an asymptotic time. Now the question is if we can detect a bud of fractal pattern in the short-time regime long before the asymptotic time. Anatomy of this formation process at short time is the subject of this paper.

In a typical fractal growth model of diffusion limited aggregation (DLA) \[3\], the short-time behavior is not studied directly. Its behavior is studied in the simulation study \[4\] and experimental study \[5\] of kinetic roughing model, in the model for fracture process \[6\] and in several cellular-automata approaches used to study the self-organized criticality \[7,8\]. These models in short-time regime are, however, concentrated on the phenomena of dynamical scaling relation \[9\] where the time-dependence of some growth parameter is investigated. The behavior of the growth parameter, however, is not an sufficient answer to the growth mechanism of fractal pattern asymptotically achieved.

In order to investigate the short-time behavior of the formation of fractal pattern, we propose a new simulation model for the fragmentation of an infinite fcc lattice. For this purpose, we adopt a box-counting method \[1\] to search for the fragmentation process instead of the fragment mass distribution. It is because the mass distribution is only defined for the low-density limit where a fragment is separated each other, which is not the case for the short-time regime where the density is not low enough to identify the fragments. In fact, we will see later that the box counting method is an appropriate method to investigate the bud of fractal, which starts as far as the system starts to expand from its saturation density.

The dynamics of the fragmentation is caused in our model by the Hubble-like radial expansion that is imposed as the initial condition of the molecular dynamics simulation \[10,11,12,13\]. The Hubble expansion is the simplest and most symmetric driving force for the fragmentation since any point in the system is equivalent to any other point, a perfectly local equivalence for everywhere when we deal with an infinite system. Infinite system is chosen to eliminate the effect of surface and thus we concentrate on a fragmentation initiated by the fracture occurring in the volume. We perform a molecular dynamics simulation using the Lennard-Jones (L-J) potential with a cut-off length \(r_c = 2.5\) \[13\]. As the initial condition, the particles are located at the lattice points of the fcc lattice in the ground state, however, 10% of randomly selected lattice points are removed so that the initial density \(\rho\) becomes 0.95 in
a reduced unit (all quantities in what follow are measured in the reduced unit). This imperfection is introduced to prepare enough number of different initial states of the fragmentation. Our main results are essentially independent of the vacancy fraction, in the range 5% to 15%. The initial velocity of a particle at \( r \) is given by \( h r \) where \( h \) is a dimensionless parameter. Thus the system mimics a uniformly expanding Hubble universe. When solving the equation of motion for the infinite system represented by one primitive cell and 26 replica cells, we impose a generalized periodic boundary condition \[10,11,12,13\], under which a replica cell with the origin at \( R_i \) moves with the velocity \( hR_i \). This means that the time evolution is completely determined by the one parameter \( h \).

A complete nascent fractal structures emerge only for suitably selected value of \( h \). When \( h \) is too small, the effect of the attractive force dominates the expansion and one single droplet whose mass equals to the total number of particles in a unit cell is formed. Since this structure depends on the unit cell size, we judge it unphysical. When \( h \) is too large, the expansion dominates the response of the system and the solutions to the equation of motion result in a simple enlarged copy of the initial configuration. Since the initial configuration with near saturation density is almost uniform, the resultant structure has trivial dimension 3, even after the expansion. It is only with \( h \) in an intermediate range, namely \( 0.1 \leq h \leq 0.4 \) that our numerical simulations result in an emergence of intriguing fractal structures.

In the box-counting method, a unit cubic cell of the size \( R_{cell} \) of expanding matter is divided into smaller cubic boxes of the size \( \delta = R_{cell}/n \) where \( n \) is a positive integer. We count the number \( N(d) \) of small boxes that contain at least one particle. The fractal dimension \( D_f \) (box counting dimension) is introduced with the relation \( N(\delta) = R_{cell}^3/\delta^{D_f} \), which is rewritten \( \log[N(\delta)] = 3 \log[R_{cell}] - D_f \log[\delta] \). When \( \log[N(\delta)] \) is proportional to \( \log[\delta] \), \( D_f \) becomes constant and the relation between \( N(\delta) \) and \( \delta \) satisfies a power-law. In the following discussion, we will use the \( \log - \log \) plot of \( N(\delta) \) and \( \delta \) to obtain \( D_f \).

Figure 1 shows the time evolution of snapshots of the fractal structures and the results of the box counting method applied to these structures. The left column corresponds to \( h = 0.1 \) and the right column to \( h = 0.2 \). For each column, the time evolution proceeds from the top to the bottom. As time proceeds the density decreases. The three graphs presented are for the average densities \( \rho =0.20, 0.10, \) and \( 0.05 \). For our studies, we prepare 50 initial configurations each of which is composed of 1,235 particles (fcc lattice with 10% vacancies). \( N(\delta) \) is calculated as an average of these 50 configurations. Note that for \( \delta < 1 \), \( N(\delta) \) is always equal to the total number of particles because two particles cannot come closer than \( \delta \sim 1 \) due to the repulsive term of the L-J potential. From lines fitted by the method of least squares and from the location of the intersection point between the two lines we conclude that the
physical size of the fractal structure grows with decreasing density and that $D_f$ does not change significantly during the time evolution for a given $h$, that is, $1.95 \leq D_f \leq 1.98$ for $h = 0.1$ and $1.66 \leq D_f \leq 1.74$ for $h = 0.2$. The value of $D_f$ for $h = 0.1$ is larger than that for $h = 0.2$ and the range of scale where $D_f$ remains constant for $h = 0.1$ is wider than for $h = 0.2$. We checked if these results are independent of the number of particles in the unit cell and found that a cell composed of 9,878 particles leads to almost the same results. In the snapshot picture, we show an example of a configuration with 9,878 particles.

A similar behavior is observed in the case of more rapid expansion. In Fig.2, we show snapshots of the fractal evolution and the corresponding plot of the results of the box counting method for $h = 0.3$ and $h = 0.4$ at $\rho = 0.05$. From Figs.1 and 2, we conclude that a larger $h$ gives a smaller $D_f$ and narrower range of scale. Note that the intersection point of the 3-dimensional and $D_f$ lines corresponds to the maximum size of the density fluctuation in the system. Beyond this crossover point, the distribution of matter is uniform and its dimensionality is three.

A summary of fractal structure for $h = 0.1$ and $h = 0.2$ is shown in Fig.3 where the fitted lines used to determine $D_f$ in Fig.1 are depicted together with additional lines for $\rho = 0.40$. For each figure, the four lines corresponding to $\rho = 0.40, 0.20, 0.10, 0.05$ are almost overlapping for small values of $\delta$. The intersection points of these lines with the four non-overlapping 3-dimensional lines in Fig.3a and Fig3b are shifted toward larger $\delta$ values with decreasing density. The overlap of the four lines for small value of $\delta$ is the most important result of this study. It means that there is a nontrivial physical process that tunes the box counting so regularly. This feature is in sharp contrast with the results for $h \to \infty$ as shown in fig.3c. In this case, the effect of expansion dominates the system response and the density snapshot is a trivial enlarged copy of the initial configuration that simply results in a parallel shift of the 3-dimensional lines in the figure. It is remarkable that the fractal structures born already at densities near $\rho = 0.4$ are maintained at lower densities. The approximate constancy of $D_f$ and the overlap of $D_f$ lines as functions of time clearly shows that there is a special correlations governing the whole system. Although we have not yet fully understood the mechanism behind this phenomenon, it is clearly caused by the critical balance between the externally imposed expanding motion and the internal potential.

When we follow the time evolution beyond $\rho = 0.05$, the least square fit of two lines to the data becomes worse. This degradation occurs when the internal expansion of connected aggregations of particles almost stops. After this time, only the space between the fragments expands. This is another reason that we used the word critical balance since we need an on-going competition between the expanding external field and internal system forces in order to get persistent fractal behavior. As a result, the range of scale for which we
get a good least-square fit for the fractal dimension $D_f$ is limited to about one-order of magnitude.

Finally, we will see the meaning of the cross-over point. As was discussed before, the box-size of this crossover point is a measure of the maximum size of the density fluctuation at a given time. The time-dependence of this size is shown in Fig. 4, where the cross-over box-size $\delta_{cross}$ for $h = 0.1$ and for $h = 0.2$ are plotted together with the cell size $R_{cell}$ as a function of the quantity $1 + ht$. The calculated points are shown by various symbols corresponding to four system densities, $\rho = 0.4, 0.2, 0.1$ and 0.05 from the left to right, and lines are fitting to these four points. The linear dependence of $R_{cell}$ on the abscissa is the original assumption of boundary motion. On the other hand, the fitting to cross-over size yields two curves

$$\delta_{cross} = 1.222(1 + ht)^{2.6883} \quad \text{for} \quad h = 0.1, \quad (1)$$
$$\delta_{cross} = 1.252(1 + ht)^{2.0436} \quad \text{for} \quad h = 0.2. \quad (2)$$

These power-law dependence of the quantity on time is known as the dynamical scaling phenomena [9,4,6,5]. The dynamical exponent depends sensitively on the expansion speed and is larger for slower expansion. Their nature seems to be quite different from the corresponding quantities examined so far [9,4,6,5]. We like to stress here that when we extrapolate to $t=0$, the value of $\delta_{cross}$ becomes about 1.2, which nearly coincides with the smallest size of the system fluctuation. We can say, in other word, that the maximum size of the fluctuation starts to increase in coincidence with the start of expansion. The bud of a fractal starts to grow from the saturation density already.

In conclusion, we have shown that a set of many-body coupled Newton equations together with an externally imposed initial condition of uniform expansion for an fcc lattice directly leads to the birth and evolution of fractal structures for specific ranges of the initial expansion rate. The fractal dimension obtained with the box counting method remains constant up to a cross-over scale, which obeys the dynamical scaling law as a function of time. The most important finding is that the nascent fractal is generated from the very beginning of the expansion, which might give a hint to the original motivation of this study, “why the fractal-like structure prevails in nature?”. 
We thank S. Chiba for encouragement and support and P. Möller for critical readings of the original manuscript. S.C. is grateful to supports from a Research fellowship of the Japan Society for the Promotion of Science for Young Scientists.

References

[1] B. B. Mandelbrot, The Fractal Geometry of Nature, Freeman, New York, 1983.

[2] O. Malcai, D. A. Lider, O. Biham, D. Avnir, Scaling range and cutoffs in empirical fractals, Phys. Rev. E 56 (1997) 2817.

[3] T. A. Witten, L. M. Sander, Diffusion-limited aggregation, a kinetic critical phenomenon, Phys. Rev. Lett. 47 (1981) 1400.

[4] J. J. Ramasco, J. M. Lópo, M. Rodríguez, Generic dynamic scaling in kinetic roughening, Phys. Rev. Lett. 84 (2000) 2199.

[5] N. M. Hasan, J. Mallett, S. dos Santos Filho, A. Pasa, W. Schwarzacher, Dynamic scaling of the surface roughness of cu deposited using a chemical bath, Phys. Rev. B 67 (2003) 081401.

[6] G.-P. Zheng, M. Li, Dynamic scaling for avalanches in disordered systems, Phys. Rev. E 63 (2001) 036122.

[7] H. J. Jensen, Self-Organized Criticality, Cambridge University Press, 1998.

[8] D. L. Turcotte, Self-organized criticality, Rep. Prog. Phys. 62 (1999) 1377.

[9] F. Family, T. Vicsek, Scaling of the active zone in the eden process on percolation networks and the ballistic deposition model, J. Phys. A 18 (1985) L75.

[10] B. L. Holian, D. E. Grady, Fragmentation by molecular dynamics: The microscopic “big bang”, Phys. Rev. Lett. 60 (1988) 1355.

[11] S. Toxvaerd, Fragmentation of fluids by molecular dynamics, Phys. Rev. E 58 (1998) 704.

[12] W. T. Ashurst, B. L. Holian, Droplet formation by rapid expansion of a liquid, Phys. Rev. E 59 (1999) 6742.

[13] S. Chikazumi, A. Iwamoto, First order phase transition of expanding matter and its fragmentation, Phys. Rev. C 65 (2002) 067601.

[14] P. J. Kraulis, a program to produce both detailed and schematic plots of protein structures, J. Appl. Crystallogr. 24 (1991) 946.
Fig. 1. Snapshots [14] of fractal structures and log-log plots of results of the box counting method for six sets of $h$ and $\rho$ values. The left column is for $h = 0.1$ and the right for $h = 0.2$. For each column, the time evolution proceeds from the top ($\rho = 0.2$) to the bottom ($\rho = 0.05$) where $\rho$ stands for the average density. In the log-log plot, $\delta$ is the size of the box used for counting and $N(\delta)$ is the number of boxes that contain at least one particle. Also plotted are vertical bars corresponding to the variance in the 50 sample cases. They are nearly the size of the dots. The lines in the log-log plot are obtained from a least square fit, the slopes of which are fractal dimensions. The images are drawn rescaled to the same size. The size before rescaling were $R_{cell} = 36.7$ ($\rho = 0.20$), 46.2 ($\rho = 0.10$) and 58.2 ($\rho = 0.05$).

Fig. 2. Snapshots [14] of fractal structures and log-log plots of results of the box counting method at average density $\rho = 0.05$ for $h = 0.3$ (top) and $h = 0.4$ (bottom). Various quantities plotted are the same as those of the bottom of Fig.1.
Fig. 3. Summary of fractal structure growth in expanding matter. The top two figures show the lines appearing in Fig. 1 together with additional lines for $\rho = 0.40$. The bottom figure is for $h \to \infty$ in which case the dimensionality of the space is 3 at any scale, a typical feature expected for the enlarged copies.

Fig. 4. The size of the crossover point $\delta_{\text{cross}}$ together with the cell size $R_{\text{cell}}$ as a function of the quantity $1 + ht$ where $t$ is the elapsed time. The calculated values are denoted by symbols corresponding to $\rho = 0.4, 0.2, 0.1, 0.05$ from the left to the right. Lines are fitting to these four symbols which expression are also given in the figure.