Formation of multifractal population patterns from reproductive growth and local resettlement

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(Received 7 February 2005; revised manuscript received 25 July 2005; published 20 October 2005)

We consider the general character of the spatial distribution of a population that grows through reproduction and subsequent local resettlement of new population members. We present several simple one- and two-dimensional point placement models to illustrate possible generic behavior of these distributions. We show, numerically and analytically, that these models all lead to multifractal spatial distributions of population. Additionally, we make qualitative links between our models and the example of the Earth at Night image, showing the Earth's nighttime man-made light as seen from space. The Earth at Night data suffer from saturation of the sensing photodetectors at high brightness ("clipping"), and we account for how this influences the determined dimension spectrum of the light intensity distribution.

DOI: 10.1103/PhysRevE.72.046213 PACS number(s): 05.45.Df, 89.75.Kd, 87.23.Cc

I. INTRODUCTION

Growing populations exist in many different areas of interest. Although the most obvious examples are biological in nature (e.g., bacterial cultures, the human population), one can also consider the growth of technological development or the growth of urban infrastructure as more abstract examples. While the specific details of these systems can be quite different, the examples stated here all share the characteristic that their populations grow on some background space. A reasonable question to ask then is what types of spatial distributions result in these growing systems. Figure 1 is a version of the popular Earth at Night (EaN) image [1]. This is a composite image taken by orbiting satellites that shows the Earth's nighttime man-made lights as seen from space. The light intensities are brightest in areas that are known to be highly populated and developed, and thus the image can be crudely thought of as representing the spatial distribution of some combined measure of technological development and human population density. We see a very heterogeneous distribution, with areas of very high intensities as well as areas of almost no light

A useful characterization of the heterogeneity of a distribution is the distribution's dimension spectrum D_q . A grid of hypercubes, or boxes, of a fixed size with edge length ϵ is used to cover the distribution, where the dimension of the hypercubes corresponds to the dimension of the underlying space (e.g., squares for two dimensions and

line segments for one dimension). D_q is then calculated via [2,3]

$$D_q = \frac{1}{1 - q} \lim_{\epsilon \to 0} \frac{\ln I(q, \epsilon)}{\ln(1/\epsilon)},\tag{1}$$

where

$$I(q, \epsilon) = \sum_{i=1}^{N(\epsilon)} \mu_i^q.$$
 (2)

Here the sum is taken over $N(\epsilon)$ nonempty grid boxes and μ_i is the fraction of the particular quantity of interest (assumed to be non-negative—e.g., light intensity in Fig. 1) that is contained in box i.

The parameter q ($q \ge 0$ in this paper) can be varied continuously to vary the influence of high- and low- μ_i boxes, with larger values of q emphasizing the boxes with larger μ_i 's. In particular, when q=1 we have the *information dimension* D_1 which, applying L'Hospital's rule to Eqs. (1) and (2), is given by

$$D_1 = \lim_{\epsilon \to 0} \frac{\sum_{i=1}^{N(\epsilon)} \mu_i \ln \mu_i}{\ln \epsilon}.$$
 (3)

A distribution is *fractal* if it possesses a fractional (noninteger) dimension (i.e., if D_q is not an integer for some q). If D_q depends on q, then the distribution is said to be *multifractal*.

Generally, when measuring the dimension of a distribution of numerical and/or experimental data, there is a maximum resolution observable. In the EaN image, for example, the image's individual pixels are the smallest boxes that

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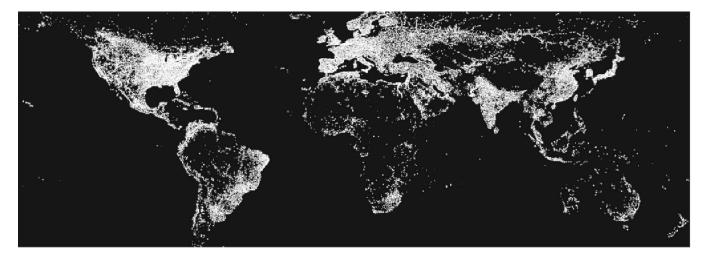


FIG. 1. A version of the Earth at Night image [1]. This is a composite image created with data from the Defense Meteorological Satellite Program of the Earth's nighttime man-made lights as seen from space. For the analysis, a gray scale $30\,000\times15\,000$ pixel image was used. The image pixels have intensity values ranging from 0 to 100.

can be used to cover the distribution in the calculation of D_a . In such cases the $\epsilon \rightarrow 0$ limit cannot be considered; instead one looks for a scaling range of ϵ where the quantity $(1-q)^{-1} \ln I(q,\epsilon)$ varies approximately linearly with $\ln(1/\epsilon)$, from which a dimension can be extracted as the slope of a straight line fitted to data in the ϵ scaling range. We illustrate this process by calculating the information dimension D_1 of North America (6000 × 3000 pixels) in the EaN image. Figure 2 shows a plot of $\sum_{i=1}^{N(\epsilon)} \mu_i \ln \mu_i$ vs $\ln \epsilon$ (with ϵ in units of pixels) for the image, where each μ_i is the fraction of the total intensity in box i. The good evidence of linearity in this plot over a scaling range in ϵ of order $e^4 \approx 50$ indicates that a fractal description makes sense. The solid line in Fig. 2 has a slope of 1.65 ± 0.02 , which we take as the value of D_1 (the error indicates the uncertainty involved in extracting the slope of the scaling region). Thus the light distribution over North America in the EaN image is fractal. (Other regions of the EaN image also exhibit fractality, with different values of D_1 .) In Sec. III we demonstrate that the EaN light distribution is also multifractal.

Yook *et al.* [4] have recently reported that the world's human population, as well as the population of internet routers and autonomous systems, is fractally distributed. Additionally, several studies [5–10] have demonstrated various fractal properties of urban settlements. We were thus led to consider the possibility of common, underlying mechanisms shared by growing populations that lead to fractal distributions.

One characteristic found in many growing populations is the existence of some generalized reproduction process, where existing members of the population generate more members. In addition to biological reproduction, in some nonbiological systems, concentrated populations can encourage the creation of more members. For example, areas that are technologically highly developed are likely to stimulate more development than areas that are not as technologically developed.

Generally, in such systems, new members of a population become situated near an existing member, but not at the exact same location. We model this process as a local "resettlement," whereby an existing member of the population produces one or more new members, who then resettle to a nearby location.

In this paper we demonstrate that these two ingredients, reproduction and local resettlement, can lead to multifractal spatial distributions for growing populations. We present several point placement models in one and two dimensions which implement these ideas and show, both numerically and analytically, that they lead to multifractal distributions.

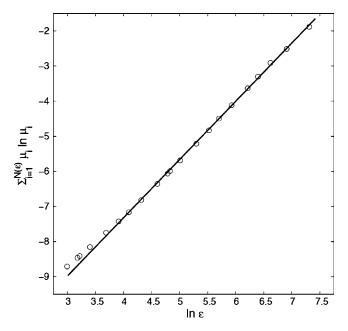


FIG. 2. Plot for calculating the information dimension D_1 [see Eq. (3)] for North America in the Earth at Night image. The quantity $\sum_{i=1}^{N(\epsilon)} \mu_i \ln \mu_i$ is plotted vs $\ln \epsilon$ (open circles), where ϵ is in units of pixels, resulting in a linear *scaling region*, the slope of which is a finite-scale approximation of D_1 . The solid line has a slope of 1.65 ± 0.02 , where the error value indicates the uncertainty involved in extracting the slope of the scaling region.

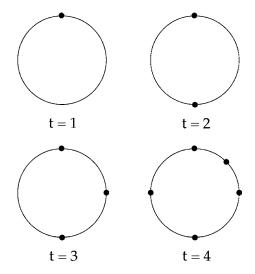


FIG. 3. Three representative steps of models 1 and 3.

One might suspect that fractal population densities arise due to strong inhomogeneities of the underlying space (e.g., inhomogeneities in the distribution of land fertility, natural resources, etc.). Our models initially (Sec. II) employ spaces that have *no inhomogeneity*, thus demonstrating that fractal population distributions can occur without inhomogeneity of the background space on which the population grows. We also consider generalizations to our models that include geographical inhomogeneities (Sec. IV) and demonstrate that, for the particular forms of the inhomogeneities that we employ, the multifractality of the distribution is unaffected.

In Sec. II we present our point placement models along with numerical results. (We emphasize that the set of models we introduce in Sec. II is not exhaustive, and many other similar models could be conceived.) Section III considers situations where a multifractal distribution is sensed by an instrument that saturates at a maximum measurable value (clipping). We show that such clipping occurs in the EaN data, affecting the determined D_q . In Sec. IV, we investigate the effect of adding geographical inhomogeneity to the underlying space of our models. In Sec. V we present analytical results for some of our models. We summarize our findings and conclude in Sec. VI. Appendix A provides details and background on the analytical calculation of the spectra of fractal dimensions for some of our models. Appendix B contains the proof of a theorem used in Appendix A (this theorem should be of very general utility, not restricted to our specific models, in analysis of the spectra of fractal dimensions).

II. MODELS AND RESULTS

A. One-dimensional random interval (model 1)

We begin with a very simple model that places points on the circumference of a circle. (For convenience we consider a circumference of unit length.) Our initial state is a single point on the circle's circumference. As the population (number of points on the circumference of the circle) grows, we

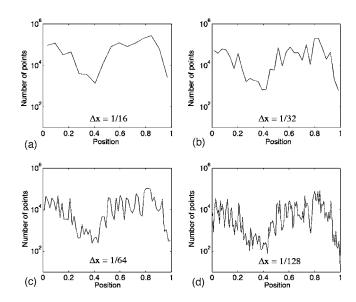


FIG. 4. Semilogarithmic plots of histograms of point locations on the unit length circumference of a circle for a point distribution ($\sim 10^6$ points) generated by model 1 with p=0.5, at time t=35, for various bin sizes Δx . (a) $\Delta x=1/16$, (b) $\Delta x=1/32$, (c) $\Delta x=1/64$, and (d) $\Delta x=1/128$.

consider the intervals of the circle's circumference between two adjacent points. At each subsequent discrete time step (each "generation") labeled t, a new point is "born" in the middle of each preexisting interval with the same given fixed birth probability p, with each new point bisecting its corresponding target interval. Several representative steps of this construction are illustrated in Fig. 3.

In this way, every interval has an equal probability to generate or attract new points at the same rate p. This also implies that regions with a high density of points (i.e., with a greater number of intervals) are more likely to attract new members than areas that are sparsely populated, creating a "rich get richer" phenomenon. We will see that all the models presented in this paper have this general characteristic. The simple rule of placing each new point in the center of an interval allows us to obtain analytical results, as will be shown in Sec. V A.

For p=1 each existing interval is bisected by a point at each time step, resulting in a uniformly distributed set of points on the circle circumference for all t. Thus, for more interesting point distributions, we turn our attention to models with p < 1. As an example, we numerically examine the p=0.5 case. Figure 4 shows semilogarithmic plots of point location histograms, along the circle circumference, with different bin sizes Δx for one realization of our point placement scheme with p=0.5, at time t=35. We observe a roughening of the plot as smaller histogram bins are used, thus giving a sense of the heterogeneity of the distribution. To quantify this heterogeneity, we calculate the dimension spectrum D_a of the distribution. We cover the distribution of points with intervals of a fixed length ϵ and assign to each interval i a measure μ_i equal to the fraction of the total points contained in i. Then, referring to Eq. (1), we plot the quantity $(1-q)^{-1} \ln I(q, \epsilon)$ vs $\ln(1/\epsilon)$ for a range of ϵ , so that the slope of the graph in the scaling range gives us D_a .

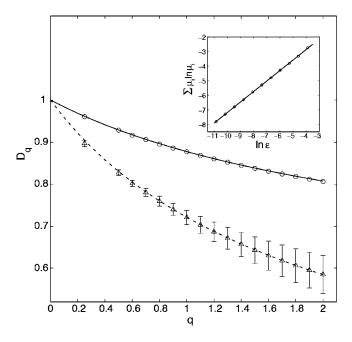


FIG. 5. Plots of D_q for model 1 with p=0.5 and $p\to 0$. The open circles (p=0.5) and triangles $(p\to 0)$ are the numerical values calculated from the slope of a line fit to $(1-q)^{-1} \ln I(q,\epsilon)$ vs $\ln(1/\epsilon)$ [see Eq. (1)], averaged over 20 realizations of point distributions (except for q=1, in which case the line is fit to $\sum_{i=1}^{N(\epsilon)} \mu_i \ln \mu_i$ vs $\ln \epsilon$ [see Eq. (3)]), with the error bars indicating the sample standard deviation for each value (in the p=0.5 case the error bars are all smaller than the open circles and we omit them). The solid line is the analytical result of Eq. (13) with p=0.5, while the dashed line corresponds to Eq. (15). As an example the inset shows the determination of D_1 for one point distribution. For p=0.5, the distributions are calculated at time t=35 and have between 1.2×10^5 and 2.8×10^6 points. For $p\to 0$ the numerical results were obtained for distributions of 10^6 points.

For $q \ge 0.5$ the plots exhibit a linear regime that allows for a relatively unambiguous determination of the dimension. However, for smaller q there tend to be shorter scaling regions, which makes the extraction of a reliable D_q more error prone. This is due to the fact that the distribution is made up of a finite collection of points. As the grid covering the distribution is made finer, grid intervals in the more sparsely populated areas begin to contain either 0 or 1 points. Furthermore, if ϵ is made small enough, eventually all the points in the distribution are covered by different boxes, and the dimension for the distribution is simply the dimension of a collection of individual points—i.e., zero. Thus, in order to get a meaningful dimension out of a finite distribution of points, a certain degree of coarseness of the covering grid is necessary. For small values of q, D_q relies heavily on boxes containing small amounts of the measure. Therefore, as ϵ is decreased, a scale is quickly reached where individual points in sparsely populated areas are covered by separate boxes, thereby destroying the overall scaling behavior. For larger q, the dense regions have a large enough influence on D_a that the effects of the finiteness of the distribution are not observed until very small scales are reached, allowing for a relatively larger scaling range.

Figure 5 is a plot of the numerically determined average

 D_q (open circles) for 20 realizations of point distributions generated by model 1 with $p\!=\!0.5$ at time $t\!=\!35$. (The number of points in each realization ranges from $1.2\!\times\!10^5$ to $2.8\!\times\!10^6$ points.) The solid line represents the analytical result for D_q derived in Sec. V A [Eq. (13) with $p\!=\!0.5$]. Both the numerical and analytical results demonstrate that this model generates multifractal distributions.

Notice that our analytical value for the box-counting dimension D_0 is 1. Unlike D_q for q positive, D_0 depends only on the limiting set of points and not their distribution. Our result that $D_0=1$ indicates that, although the point distribution is multifractal, the set filled by the points in model 1 as time t goes to infinity is not fractal. In fact, the entire circumference is filled with probability 1, which can be easily seen as follows. Since each interval has a probability (1-p) of not being bisected at each time step, the probability for an interval not to have been bisected by time t is simply $(1-p)^t$, which approaches zero as t approaches infinity. Therefore no empty intervals remain as $t \rightarrow \infty$. One can similarly argue for the other models in this paper (except model 7) that the limiting sets are not fractal; the heterogeneous point distributions they yield are reflected by the fractional values of D_q for q > 0.

We can also consider model 1 in the limit where $p \rightarrow 0$. We think of this limit as the situation where at most one interval is chosen to be bisected from all available intervals at each time step. This allows us to reformulate the point placement scheme as follows. We begin with the same initial condition as the original model, which is a single point on the circle circumference. Then at each subsequent discrete time step we place a new point according to the following prescription: (a) a target point, or parent, is chosen from all of the preexisting points with equal probability; (b) the new point, or child, is placed in the middle of one of the two empty intervals adjacent to the target point with equal probability. Steps (a) and (b) are repeated until the desired system size is reached.

To show that this point placement scheme coincides with the $p \rightarrow 0$ limit of model 1, we plot the numerically determined average D_q for 20 realizations of 10^6 point distributions generated by it in Fig. 5 (open triangles) and observe that the data agree with the $p \rightarrow 0$ limit of the analytical result for D_q (dashed line) calculated in Sec. V A [Eq. (15)].

Thus we see that model 1 can create point distributions with D_q ranging from the trivial D_q =1 curve, corresponding to p=1, to the D_q curve associated with $p \rightarrow 0$.

B. Two-dimensional square (model 2)

Model 2 is a two-dimensional analogue of model 1, where points are placed on the unit square $(0 \le x \le 1, 0 \le y \le 1)$, instead of the circumference of a circle. Beginning with a single point at (0.5,0.5), we consider reproduction by a process mimicking cell division. At each discrete time step, each preexisting point, or parent, has a probability p of dividing into four points, or children, where these four points disperse to occupy (resettle) the centers of the four equal size squares obtained by partitioning the original square associated with the parent point; see Fig. 6.

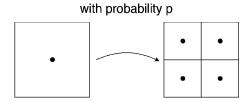


FIG. 6. Illustration of the fourfold "cell division" step in model 2.

Note that this construction can be specified without reference to points, but only to the squares they occupy: At each discrete time step, each square is subdivided into four equal new squares with probability p. After the desired iterations of this area partitioning procedure, the corresponding population distribution can be produced by simply placing a point in the center of each square.

Since p=1 results in a homogeneous planar distributions of points $(D_q=2)$, we again consider p<1 for more interesting point distributions. Choosing p=0.5, we plot in Fig. 7 a distribution of points generated by this model at time t=15. The points form a heterogeneous distribution with dense clusters in some regions and relatively sparsely populated areas in other regions. Figure 8 is a plot of the numerically determined average D_q (open circles) for 20 realizations of point distributions generated by model 2, with t=15 and p=0.5. The solid line represents the analytical result for D_q derived in Sec. V B.

As we did for model 1, we can explore the $p \rightarrow 0$ limit by considering the construction of the point distribution one point at a time. The appropriate process is as follows: At each time step a "parent" point is chosen out of all preexisting points with equal probability and produces four "children" points according to the fourfold cell division process described above. The open triangles in Fig. 8 are the average D_q of 20 independent 10^6 point distributions generated by this construction, while the dashed line is the $p \rightarrow 0$ limit of

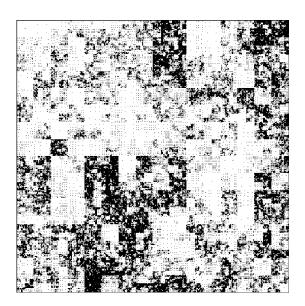


FIG. 7. A plot of a point distribution on the unit square generated by model 2 with t=15 and p=0.5.

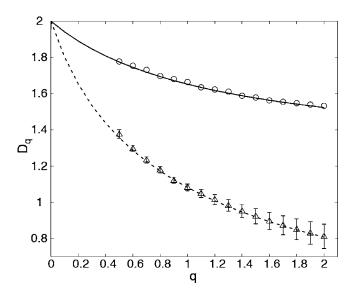


FIG. 8. Plot of D_q for model 2 with $p\!=\!0.5$ and $p\!\to\!0$. The open circles $(p\!=\!0.5)$ and triangles $(p\!\to\!0)$ are the average numerical values of D_q (see caption in Fig. 5) obtained from 20 realizations of point distributions generated by this model and the error bars are the sample standard deviations for each value (in the $p\!=\!0.5$ case the error bars are all comparable in size to the open circles and we omit them). The solid line is the analytical result of Eq. (21), and the dashed line is Eq. (23). For $p\!=\!0.5$, the distributions are calculated at time $t\!=\!15$ and have between $1.2\!\times\!10^5$ and $2.9\!\times\!10^6$ points. For $p\!\to\!0$ the numerical results were obtained for distributions of 10^6 points.

the theoretical result (Sec. V B). Thus, model 2 creates point distributions with D_q ranging between D_q =2 for p=1 and the D_q associated with $p \rightarrow 0$, and hence, other than the p=1 case, results in multifractal point distributions.

C. Two-dimensional triangle (model 3)

Model 3 is a slight variation of model 2. Here triangles, as opposed to squares, are the basis of the construction. We begin with a point in the middle of a single equilateral triangle with unit edge length. At each time step each preexisting triangle (parent) is divided into four identical equilateral triangles (children) with probability p, with the edge lengths of the child triangles equal to half the edge length of the original parent triangle (see Fig. 9) and with points placed in the centers of the child triangles. We show in Sec. V B that model 3 has the identical analytical expression for D_q as model 2 and, thus, also results in multifractal distributions.

D. Two-dimensional random square (model 4)

Model 4 is a two-dimensional point placement scheme similar to the $p \rightarrow 0$ limit construction of model 2, where

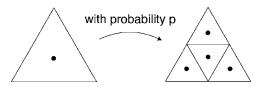


FIG. 9. Illustration of the "reproductive" step in model 3.

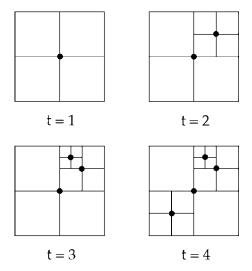


FIG. 10. Three representative steps of models 4 and 6.

points are placed on the unit square, $(0 \le x \le 1, 0 \le y \le 1)$. We begin with a single point at (0.5,0.5). Imagining vertical and horizontal lines drawn through this point, we see that the point is at the vertex of four neighboring squares. At the next time step, we choose one of the squares at random with equal probability and place a point in its middle. We then divide this chosen square into four smaller squares by horizontal and vertical lines through the newly added point. Then we continue this process: At each subsequent discrete time step we pick a target point (parent) from all preexisting points with equal probability and add a new point (child) to the center of one of the target point's four adjacent squares, also chosen with equal probability. Several representative steps of this construction are illustrated in Fig. 10.

Notice that some squares have two potential parent points at their corners and some squares have only one. The former squares are twice as likely to be subdivided as the latter squares, unlike the $p \rightarrow 0$ limit of model 2, in which all squares are equally likely to be subdivided.

Figure 11 shows a distribution generated by this model's point placement scheme, where we have successively magnified dense regions of the distribution to illustrate the large differences in densities that result at various length scales. Figure 12 is a plot of the average D_q of 20 independent distributions generated by this model, where the solid line is the analytical result (37) derived in Sec. V C. Both the nu-

merical and analytical results confirm that model 4 generates multifractal distributions.

E. One-dimensional larger interval (model 5)

We return now to one dimension with a model similar to the construction of the $p \rightarrow 0$ limit of model 1, but with a different local resettlement rule. Referring to the $p \rightarrow 0$ point placement prescription of model 1, the initial state and step (a) are identical while step (b) is modified. Instead of a new point being randomly placed in one of the adjacent intervals of the chosen target point with equal probability, the new point is placed in the larger of the two intervals adjacent to the target point, and if the two intervals are equal in size, one is chosen at random.

In this way, the new points seek out more sparsely populated local intervals to settle within. This strategy is reasonable in a context such as human resettlement, where the settlers may try to combine the convenience of being close to others with the possible advantages associated with more space. We expect this resettlement scheme to yield distributions that are more homogeneous than those created by the $p \rightarrow 0$ limit of model 1 and, hence, yield greater D_a .

We numerically obtain D_q and find the dimension spectrum to be multifractal, having the same qualitative shape as the $p \rightarrow 0$ result for D_q for model 1. However, due to the more uniform nature of the distributions created by this model, the specific values of D_q are larger at each q (e.g., $D_1 = 0.84 \pm 0.01$ (averaged over 20 distributions of 10^6 points, where the error is the sample standard deviation) for this model vs our analytical result $D_1 \approx 0.72$ for the $p \rightarrow 0$ limit of model 1 [Eq. (14)]).

F. Two-dimensional sparse square (model 6)

Model 6 is a two-dimensional analog of model 5. The initial state and step (a) of its construction are the same as in model 4. Our aim is to extend step (b) of model 5, the choosing of the larger interval for settlement, to two dimensions, and for that purpose, we make the following observation. In the square-based two-dimensional point placement scheme of model 4, there are two different types of squares which we refer to as type-I and type-II squares. A type-I square only has one of its vertices occupied by a point, while a type-II square has two opposite vertices occupied by points. (See

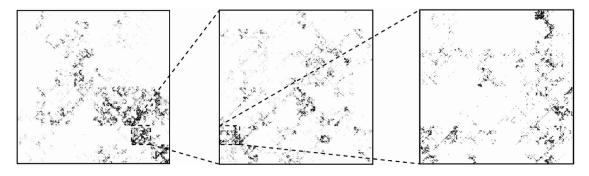


FIG. 11. A plot of a 4×10^5 point distribution on the unit square generated by model 4. Dense regions are magnified to illustrate the heterogeneity of the point densities in the distribution at different scales.

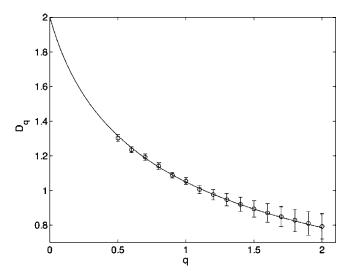


FIG. 12. Plot of D_q for model 4. The open circles are the average numerical values of D_q (see caption in Fig. 5) obtained from 20 realizations of 10^6 point distributions generated by this model and the error bars are the sample standard deviations for each value. The solid line is the analytical result of Eq. (37).

Fig. 13 for an illustration of type-I and type-II squares.) Thus a type-I square can be regarded as more sparsely populated than a type-II square of equal size. With this in mind, step (b) for model 6 is to place the new node in the middle of the largest square adjacent to the target point. In case of a tie, choose a type-I square over a type-II square, with any subsequent ties being resolved by an equal probability random choice from the remaining candidate squares.

We plot a distribution generated by this model in Fig. 14. We find this distribution to be multifractal, with a D_q that is similar in shape to, but everywhere larger than, the D_q of model 4 (except for q=0 where the two coincide with D_0 =2). In particular, the information dimension is D_1 =1.30±0.02 (averaged over 20 distributions of 10^6 points), as compared to our analytical result of D_1 \simeq 1.05 for model 4 [Eq. (38)].

G. Two-dimensional unstructured (model 7)

Looking at Fig. 7 (model 2), Fig. 11 (model 4), and Fig. 14 (model 6), we notice square-shaped artifacts and diagonal

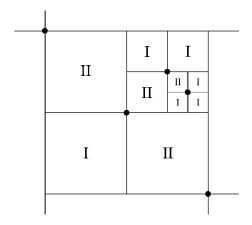


FIG. 13. Illustration of type-I (labeled I) and type-II (labeled II) squares, where black dots indicate populated vertices.

point formations that arise due to the structured, square-based nature of those point placement schemes. Model 7 is a two-dimensional point placement scheme that aims to avoid such artifacts, while maintaining the two ingredients of reproduction and local resettlement. We begin with a single point and assign to this point a *child distance* d_c =1/4. At each subsequent discrete time step t, each preexisting point (parent) is selected to "reproduce" with probability p, where the reproduction consists of a child point being placed at a random location on the circumference of a circle centered on the parent point and with radius equal to the parent point's d_c . A child distance equal to half the parent point's d_c is assigned to the child point.

Model 7 results in multifractal point distributions for *all* values of p. This is in contrast to the trivial dimension spectrum D_q =1 for the p=1 limit of model 1 and the trivial D_q =2 result for the p=1 limit of models 2 and 3. In addition, while all the models presented thus far possess an integer-valued box counting dimension D_0 equal to the dimensionality of their respective ambient spaces, $D_0 \neq 2$ for all p in model 7.

In a manner similar to the $p \rightarrow 0$ limits of models 1 and 2, we can define the $p \rightarrow 0$ limit of model 7 and create our point distribution one point at a time. The construction is as follows: At each time step a parent point is chosen from all preexisting points with equal probability. The parent point then produces a child according to the reproductive step described above in the general p case. A distribution generated by the $p \rightarrow 0$ limit of model 7 is shown in Fig. 15. As intended, there are no grid-type artifacts visible. The $p \rightarrow 0$ limit of model 7 produces point distributions with $D_1 = 1.36 \pm 0.05$ (averaged over 20 distributions of 10^6 points).

H. Discussion

Table I summarizes our results for D_1 for our models and for the Earth at Night. Table II summarizes gives qualitative

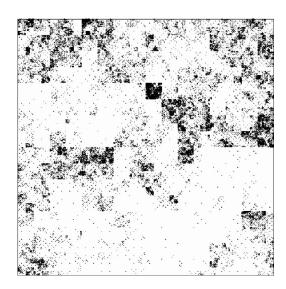


FIG. 14. A plot of a 10⁵-point distribution on the unit square generated by model 6.

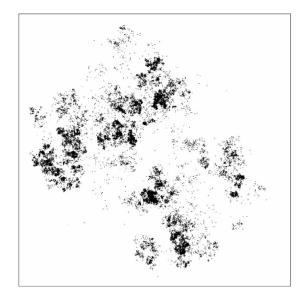


FIG. 15. A plot of a 10^5 -point distribution on the unit square generated by the $p \rightarrow 0$ construction of model 7.

descriptions of the reproduction and local resettlement mechanisms of our seven models.

Our result that the different strategies of reproduction and local resettlement used in these models (Table II) all lead to multifractal spatial patterns of population density strongly suggests that multifractality may be a generic feature in real situations in which processes involving reproduction and local resettlement take place.

We emphasize that the spirit of our approach is very minimalist. In many real cases multiple interacting complex processes undoubtedly influence the determination of population patterns. For example, several important factors for the EaN image are geography (mountains, rivers, deserts, etc.), politics, societal and cultural factors, economics, etc. Our models show that considerations of underlying spatial heterogeneities are not required for explaining the existence of fractally heterogeneous distributions: even very simple dynamics incorporating reproduction and local resettlement is sufficient.

III. EFFECT OF CLIPPED DATA ON MULTIFRACTALITY

We now revisit the North American part of the Earth at Night image (Fig. 1) and analyze its fractal properties. Figure 16 is a plot of D_q calculated from the measured light distribution, where the error bars indicate the uncertainty involved in extracting the slopes of the various scaling regions. Although we observe a multifractal D_q curve, studies of human population distributions [8,10] suggest that the a steeper D_q should result. We now show how this discrepancy can be resolved.

Figure 17 is a log-log plot of the histogram of pixel intensities for North America in the EaN image. A spike is observed for a range of intensity values near the maximum value. This suggests that some regions with high light intensities caused some of the photoelectric cells of the satellites' sensors to saturate, thereby clipping the intensity values at a maximum allowable value. Additionally, it is known that, when a photoelectric cell is subjected to high intensities, it can trigger surrounding cells to register more light, resulting in the so-called "blooming" effect, which could account for the broadness of the observed high-intensity spike. We note that the saturated pixels make up about 8% of the total number of nonzero-intensity pixels. We expect that blooming only changes intensities by a marginal amount, whereas clipping can make a drastic change. Thus the effect of blooming on D_a should be negligible and we ignore it.

To illustrate the effects that this type of clipping has on multifractal distributions, we apply a scheme that mimics the clipping in the EaN image to a point distribution generated by one of our models. For this purpose we choose a (3×10^6) -point distribution generated by the $p \rightarrow 0$ limit of model 2. (Similar results are obtained using our other models.) First we roughly match the smallest scale of our distribution to that of the EaN image. Since the part of the EaN image we analyze is made up of 3000×6000 pixels, we cover our distribution with a fixed size grid of simulated pixels in which a simulated pixel corresponds to a grid square with edge length of 2^{-12} , which is about $1/\sqrt{3000 \times 6000}$. Thus, while each pixel of the EaN image represents a light intensity value, each simulated pixel contains a number of individual points. The clipping procedure

TABLE I. A summary of numerical and theoretical results for the point placement models and the Earth at Night image (Fig. 1).

Model	Dimensionality of space	D_1	D_q (theory)
1	1	Eq. (14)	Eq.(13)
$1 (p \rightarrow 0)$	1	$1/(2 \ln 2) \simeq 0.72$	Eq. (15)
2,3	2	Eq. (22)	Eq. (21)
$2,3 \ (p \to 0)$	2	$3/(4 \ln 2) \approx 1.08$	Eq. (23)
4	2	$8/(11 \ln 2) \approx 1.05$	Eq. (37)
5	1	0.84 ± 0.01	_
6	2	1.30 ± 0.02	_
$7 (p \rightarrow 0)$	2	1.36 ± 0.05	_
Earth at Night (North America)	2	1.65±0.02	_

TABLE II. Summary of reproduction and local resettlement strategies employed in models 1–7.

Model	Reproduction	Resettlement	
1	parent has one child	child settles in the middle of chosen interval	
2	parent reproduces by four-fold cell division	children share equally the original square area occupied by their parent	
3	parent has three children	parent remains in place; children settle equal shares of parent's triangular area	
	or		
	parent reproduces by four-fold cell division	children share equally the original triangular area occupied by their parent	
4	parent has one child	child settles randomly chosen adjacent square	
5	parent has one child	child settles largest adjacent interval	
6	parent has one child	child settles largest and sparsest adjacent square	
7	parent has one child	child settles in a random direction	

involves picking a maximum number of points allowed in a simulated pixel and changing the number of points in any simulated pixel containing more than this amount to the maximum amount. In order for the clipped point distribution to exhibit an appreciable range for the numbers of points in the individual simulated pixels we choose the maximum clipping value such that the total number of clipped simulated pixels make up only 4% of the total nonempty simulated pixels (this is in contrast to the 8% observed in the EaN image). For this particular distribution, any simulated pixel containing more than 40 points has its count reset to 40.

Figures 18(a) and 18(b) are histograms showing the distribution of the number of points contained in individual simulated pixels before and after the clipping is applied, respectively. We see that the clipped histogram is qualitatively similar to the histogram for the EaN image (Fig. 17). Next,

we calculate D_q for the point distribution before and after clipping (open circles and open triangles in Fig. 19, respectively). We see that the clipping procedure transforms the unclipped D_q to a flatter curve, making it more similar to the D_q for North America in the EaN image (Fig. 16). Thus, we are able to conclude that the EaN image characteristics are consistent with a multifractal light intensity distribution with a steep D_q sensed by an instrument that saturates at a maximum measurable intensity value.

IV. INHOMOGENEITY

Our models can be generalized to include geographical inhomogeneity. For example, one can imagine that the underlying space is supplemented by a fertility field and that reproduction is more likely in regions of higher fertility. To investigate the effects of such inhomogeneities on the D_q of

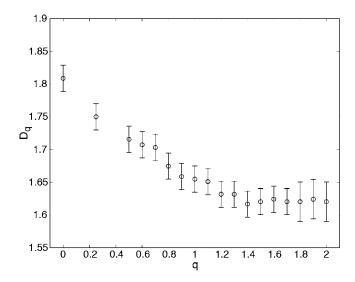


FIG. 16. Plot of D_q for North America in the Earth at Night image. The error bars reflect the uncertainty involved in determining the slopes of the scaling regions in the plots of the quantity $(1-q)^{-1} \ln I(q, \epsilon)$ vs $\ln(1/\epsilon)$ [see Eq. (1)].

a point distribution, we modify the $p \rightarrow 0$ construction of model 2 by use of a fertility field F(x,y) which we use to construct a space-dependent, parent node selection probability. If the location of point i is (x_i,y_i) , then the selection probability for this point is taken to be $P_i=F(x_i,y_i)/\sum_j F(x_j,y_j)$ (this is in contrast to the original prescription of choosing the parent point from all preexisting points with equal probability). We investigate two forms for the fertility field. First we use a smooth field

$$F_s(x, y) = 1 - 0.3\{\cos(2\pi x) + \sin(2\pi y)\}\tag{4}$$

on the unit square $(0 \le x \le 1, 0 \le y \le 1)$. We generate point distributions with this modified version of the $p \to 0$ limit of model 2 and find that, while the regions of larger F_s in the center of the square are much more dense with points than the regions of smaller F_s near the corners, there is, neverthe-

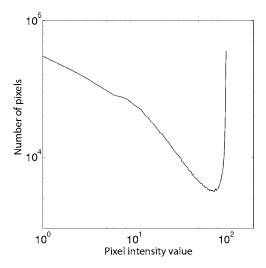


FIG. 17. Histogram of individual pixel intensities for North America in the Earth at Night image.

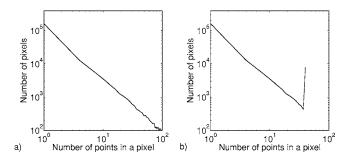


FIG. 18. Histograms of the number of points in individual simulated pixels (boxes in an ϵ =1/4096 grid) used to cover a (3×10^6) -point distribution generated by the $p\to0$ limit of model 2 (a) for the original distribution and (b) after the original distribution was clipped by allowing a maximum of 40 points in any simulated pixel.

less, no discernible difference in D_q compared to the homogeneous case. That is, both the values of D_q and the quality and extent of the scaling ranges remain unchanged.

Next, to explore whether the above result can be attributed to the local smoothness of F_s , we consider a rough fertility field

$$F_r(x,y) = 1 + 0.1[f_w(x) + f_w(y)], \tag{5}$$

$$f_w(z) \equiv -\sum_{j=0}^{\infty} \alpha^{-j} \cos(2\pi\beta^j z), \tag{6}$$

with α =1.5 and β =3. For $\alpha < \beta$, the function $f_w(z)$ is a "Weierstrass function" (see Fig. 20); it is rough in the sense that, although it is continuous, it is nondifferentiable, and the graph of $f_w(z)$ vs z is a fractal curve [the fractal dimension of the curve is $2-(\ln \alpha)/(\ln \beta) \approx 1.63$ [11]]. With α =1.5, F_r

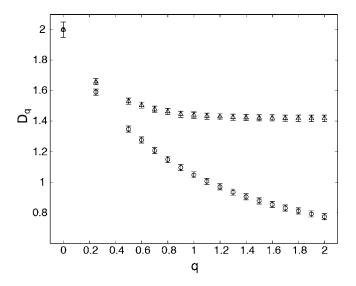


FIG. 19. Plot of D_q for a (3×10^6) -point distribution generated by the $p \to 0$ limit of model 2 before the clipping procedure discussed in Sec. III is applied (open circles) and after it is applied (open triangles). The error bars reflect the uncertainty involved in determining the slopes of the scaling regions in the plots of the quantity $(1-q)^{-1} \ln I(q,\epsilon)$ vs $\ln(1/\epsilon)$ [see Eq. (1)].

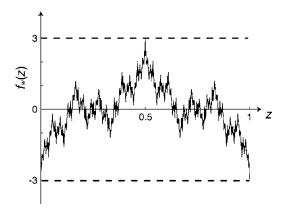


FIG. 20. A plot of the Weierstrass function $f_w(z)$ [Eq. (6)]. In our numerical implementation the infinite upper limit on the summation in Eq. (6) is replaced by j_{max} , where j_{max} is chosen such that the smallest value of ϵ used in determining D_q is at least 1 order of magnitude larger than the smallest scale of the roughness, $3^{-j_{max}}$.

has the same range of variation as F_s (i.e., 0.4–1.6). Applying this fertility field to the $p \rightarrow 0$ limit of model 2 results in the same values and scaling ranges for D_q as in the homogeneous case and the smooth fertility field case [Eq. (4)]. This suggests that, while inhomogeneities of the underlying space may dictate certain aspects of the distribution of a growing population—for example, where the points are more likely to settle—its fractality can be due mainly to reproduction and local resettlement processes.

V. THEORY

A. Model 1

Here we derive an analytical expression for the fractal dimension spectrum D_q of population distributions generated by the one-dimensional point placement scheme of model 1.

We define N(k,t) to be the average (over different realizations of model 1) number of intervals of length 2^{-k} at time t. The initial condition for this model (at t=1) is a single point on the unit length circumference of a circle, implying $N(k,1) = \delta_{k0}$, where δ_{kj} is the Kronecker delta. At each time t, each preexisting interval is bisected with probability p. Thus, with probability p an interval of length 2^{-k} is replaced by two intervals of lengths $2^{-(k+1)}$. The discrete time evolution equation for N is then

$$N(k,t+1) = 2pN(k-1,t) + (1-p)N(k,t), \tag{7}$$

where pN(k-1,t) is the average (over different realizations of model 1) number of intervals of length $2^{-(k-1)}$ bisected at time t+1 and (1-p)N(k,t) is the average number of 2^{-k} length intervals that remain unchanged at t+1. We now define $N(t) = \sum_k N(k,t)$ to be the average number of intervals (of all sizes) at time t. Taking the sum over k of Eq. (7) we obtain the simple recursion relation

$$N(t+1) = (1+p)N(t),$$
(8)

which, since N(1)=1, gives the result $N(t)=(1+p)^{t-1}$ for the evolution of the average total number of intervals.

In Appendix A, we introduce the partition function formalism for calculating D_q . We define the quantity $\tau = (q-1)D_q$ and derive the following expression for $q(\tau)$ for model 1 [Eq. (A10)]:

$$q(\tau) = 1 + \lim_{t \to \infty} \frac{\ln\langle 2^{k\tau} \rangle_t}{\ln N(t)}, \quad \langle 2^{k\tau} \rangle_t = \frac{\sum_k N(k, t) 2^{k\tau}}{N(t)}. \tag{9}$$

Multiplying both sides of Eq. (7) by $2^{k\tau}$ and then summing over k, we obtain the recursion relation

$$\langle 2^{k\tau} \rangle_{t+1} = \frac{1}{1+p} (2^{\tau}2p + 1 - p) \langle 2^{k\tau} \rangle_t, \tag{10}$$

which, since $\langle 2^{k\tau} \rangle_{t=1} = 1$, gives

$$\langle 2^{k\tau} \rangle_t = \left(\frac{1}{1+p} (2^{\tau} 2p + 1 - p) \right)^{t-1}.$$
 (11)

Substituting Eq. (11) into Eq. (9), we obtain

$$q = \frac{\ln(2^{\tau}2p + 1 - p)}{\ln(1 + p)}.$$
 (12)

Then, since $D_q = \tau(q)/(q-1)$, we invert Eq. (12) for $\tau(q)$ and obtain the following result for D_q :

$$D_q = \frac{\log_2\left(\frac{p-1+(1+p)^q}{2p}\right)}{a-1}.$$
 (13)

In particular, using L'Hospital's rule,

$$D_1 = \frac{(1+p)\log_2(1+p)}{2p}. (14)$$

For p=1 we get the expected result $D_q=1$, while for the $p \rightarrow 0$ limit, we obtain

$$D_q = \frac{\log_2(1+q) - 1}{q - 1} \tag{15}$$

and

$$D_1 = \frac{1}{2 \ln 2} \simeq 0.72. \tag{16}$$

B. Models 2 and 3

Here we derive the analytical D_q for models 2 and 3. We adopt the area partitioning viewpoint presented in Sec. II B, where we focus on the squares containing the points rather than the points themselves. We begin by defining N(k,t) to be the average number of squares with edge length 2^{-k} at time t. The initial state (at t=1) is a single square (with a point at its center) with unit edge length, giving us $N(k,1) = \delta_{k0}$. At each time t, each preexisting square is divided into four equal-sized squares with half the edge length of the original square with probability p (Fig. 6). Thus, with probability p a square of edge length 2^{-k} is replaced by four squares of edge lengths $2^{-(k+1)}$. The discrete time evolution equation for N(k,t) is then

$$N(k,t+1) = 4pN(k-1,t) + (1-p)N(k,t).$$
 (17)

We define $N(t) = \sum_k N(k,t)$ to be the average number of squares (of all sizes) at time t and take the sum over k of Eq. (17) to obtain

$$N(t+1) = (1+3p)N(t), (18)$$

which, since N(1)=1, gives the result $N(t)=(1+3p)^{t-1}$ for the evolution of the average total number of intervals. Interchanging the squares in the above derivation with equilateral triangles, one can see that the same evolution equation holds for the triangles in model 3.

In Appendix A we show that Eq. (9), which allowed us to calculate $q(\tau)$ in terms of $\langle 2^{k\tau} \rangle_t$ for model 1, holds for models 2 and 3 as well. Following the procedure in Sec. V A, we multiply both sides of Eq. (17) by $2^{k\tau}$ and sum over k, solving the resulting recursion relation to obtain

$$\langle 2^{k\tau} \rangle_t = \left(\frac{1}{1+3p} (2^{\tau} 4p + 1 - p) \right)^{t-1}.$$
 (19)

Substituting Eq. (19) into Eq. (9), we obtain

$$q = \frac{\ln(2^{\tau}4p + 1 - p)}{\ln(1 + 3p)}.$$
 (20)

Thus, for models 2 and 3 we obtain the multifractal dimension spectrum

$$D_{q} = \frac{\log_{2}\left(\frac{p-1+(1+3p)^{q}}{4p}\right)}{q-1}$$
 (21)

and the information dimension

$$D_1 = \frac{(1+3p)\log_2(1+3p)}{4p}.$$
 (22)

For p=1 we get $D_q=2$, while for the $p \rightarrow 0$ limit, we obtain

$$D_q = \frac{\log_2(1+3q) - 2}{q-1} \tag{23}$$

and

$$D_1 = \frac{3}{4\log 2} \simeq 1.08. \tag{24}$$

C. Model 4

The calculation of D_q for model 4 is based on the classification of squares into types I and II, discussed in Sec. II F. We begin by defining $N_1(k,t)$ and $N_2(k,t)$ to be the average (over different realizations) number of type-I and type-II squares, respectively, with edge length 2^{-k} at time t. The initial state (t=1) is a single point at (0.5,0.5). The point divides the unit square into four equal squares of edge length 1/2, with each square having one point on one of its vertices. Thus $N_1(k,1)=4\delta_{k1}$, while $N_2(k,1)=0$ for all k. At each time t, the new point is equally likely to appear in

each type-I square, but is twice as likely to appear in each type-II square, because there are two parent points that can produce offspring in each type-II square. In other words, a type-I square with edge length 2^{-k} is chosen with probability $N_1(k,t)/\Sigma_j[N_1(j,t)+2N_2(j,t)]$, while a type-II square of the same size is chosen with probability $2N_2(k,t)/\Sigma_j[N_1(j,t)+2N_2(j,t)]$. When a type-I square is chosen, the new point is placed in its center, destroying the original square and creating three type-I squares and one type-II square, all with half the edge length of the original square. On the other hand, when the new point is in a type-II squares, it is replaced by two type-I squares and two type-II squares, all having half the edge length of the original square. Notice that either way, the quantity $\Sigma_j[N_1(j,t)+2N_2(j,t)]$ increases by 4 between time t and t+1, so that $\Sigma_j[N_1(j,t)+2N_2(j,t)]=4t$.

The discrete time evolution equations for both types of squares are then

$$N_{1}(k,t+1) = N_{1}(k,t)\left(1 - \frac{1}{4t}\right) + N_{1}(k-1,t)\frac{3}{4t} + 2N_{2}(k-1,t)\frac{2}{4t},$$
(25)

$$N_2(k,t+1) = N_2(k,t) \left(1 - \frac{2}{4t}\right) + N_1(k-1,t) \frac{1}{4t} + 2N_2(k-1,t) \frac{2}{4t}.$$
 (26)

In Appendix A we show that for model 4 [Eq. (A14)],

$$q(\tau) = \lim_{t \to \infty} \frac{\ln[T_1(t) + 2^q T_2(t)]}{\ln t},$$
(27)

where $T_i(t) = \sum_k 2^{k\tau} N_i(k,t)$. Multiplying Eqs. (25) and (26) by $2^{k\tau}$ and summing over k, we obtain

$$T_1(t+1) = T_1(t) \left(1 + \frac{3 \times 2^{\tau} - 1}{4t}\right) + T_2(t) \frac{4 \times 2^{\tau}}{4t},$$
 (28)

$$T_2(t+1) = T_2(t) \left(1 + \frac{4 \times 2^{\tau} - 2}{4t} \right) + T_1(t) \frac{2^{\tau}}{4t}.$$
 (29)

At this point, we make the continuous time approximation $T_i(t+1)-T_i(t)\approx dT_i(t)/dt$, valid for large t, which allows us to represent Eqs. (28) and (29) as the matrix differential equation

$$\frac{d\mathbf{T}(s)}{ds} = \mathbf{M} \cdot \mathbf{T}(s), \tag{30}$$

where

$$\mathbf{T}(s) = \begin{bmatrix} T_1(s) \\ T_2(s) \end{bmatrix},\tag{31}$$

$$\mathbf{M} = \begin{bmatrix} 3 \times 2^{\tau} - 1 & 4 \times 2^{\tau} \\ 2^{\tau} & 4 \times 2^{\tau} - 2 \end{bmatrix}, \tag{32}$$

and $s = (1/4) \ln t$.

Equation (30) is solved in terms of the eigenvalues λ_a, λ_b and the eigenvectors $\mathbf{T}_a, \mathbf{T}_b$ of the matrix \mathbf{M} , giving

$$\mathbf{T}_i \sim e^{\lambda_i s} = t^{\lambda_i/4},\tag{33}$$

for i=a,b, with λ_a,λ_b distinct and real and $\lambda_a>\lambda_b$. We can express the quantity $T_1(t)+2^qT_2(t)$ in Eq. (27) in terms of a linear combination of the components of these eigenvectors, whose long-time behavior is dominated by the behavior of T_a , the eigenvector associated with the larger eigenvalue λ_a . Thus Eq. (27) gives us the simple result

$$q = \lambda_a/4. \tag{34}$$

The eigenvalues of M are the roots of its characteristic polynomial,

$$\lambda^2 + \lambda(3 - 7 \times 2^{\tau}) + 2 - 10 \times 2^{\tau} + 8 \times 2^{2\tau} = 0.$$
 (35)

This equation is also quadratic in 2^{τ} , allowing us to solve, for $\tau(\lambda)$,

$$\tau(\lambda) = \log_2 \left[\frac{1}{16} (10 + 7\lambda \pm \sqrt{17\lambda^2 + 44\lambda + 36}) \right]. \quad (36)$$

We substitute the relationship from Eq. (34) into Eq. (36) and, since $\tau(q) = (q-1)D_q$, we resolve the ambiguity of the \pm sign by requiring that $\tau(q=1)=0$. This results in

$$D_q = \frac{1}{q-1} \log_2 \left[\frac{1}{8} (5 + 14q - \sqrt{68q^2 + 44q + 9}) \right]$$
 (37)

for the dimension spectrum of model 4. In particular, applying L'Hospital's rule to Eq. (37), the information dimension is

$$D_1 = \frac{8}{11 \ln 2} \simeq 1.05. \tag{38}$$

VI. CONCLUSION

The main conclusion of this paper is that reproduction and local resettlement processes may lead to multifractal spatial distributions for growing populations. We introduced a number of point placement models in one and two dimensions and showed that the models resulted in multifractal distributions. Furthermore, we have demonstrated a qualitative similarity between the example of the Earth at Night image and clipped versions of distributions generated by our models. We thus suggest that the mechanism by which our models create multifractal distributions may be operative in the growth of real systems.

ACKNOWLEDGMENTS

We thank Dan Lathrop and Rowena Ball for useful conversations. This work was supported by the ONR (Physics) and by NSF (Contract Nos. PHYS 0098632 and DMS 0104087).

APPENDIX A: CALCULATION OF MODEL DIMENSION SPECTRA

1. Partition function formalism

The partition function formalism [12,13] is an alternative to the fixed-sized grid method of Eq. (1) for calculating a dimension spectrum of a measure. We demonstrate here how it allows us to relate the various interval and square-size distribution functions found in models 1, 2, 3, and 4 to the dimensions of the point distributions generated by them.

We cover the measure of interest with a disjoint covering $\{S_i\}$, $i=1,2,\ldots,N$, where each element S_i of the covering set has a diameter ϵ_i less than or equal to δ . (The diameter ϵ_i is the largest possible distance between two points in S_i .) The partition function is defined as

$$\Gamma_q(\tau, \{S_i\}, \delta) = \sum_{i=1}^N \mu_i^q / \epsilon_i^\tau, \tag{A1}$$

where μ_i is the measure of S_i . For a given δ the covering $\{S_i\}$ is now chosen such that Eq. (A1) is maximized (for q > 1) or minimized (for q < 1), which defines

$$\Gamma_{q}(\tau, \delta) = \begin{cases} \sup_{S_{i}} \Gamma_{q}(\tau, \{S_{i}\}, \delta), & \text{for } q > 1, \\ \inf_{S_{i}} \Gamma_{q}(\tau, \{S_{i}\}, \delta), & \text{for } q < 1. \end{cases}$$
(A2)

Then, letting $\delta \rightarrow 0$, we define

$$\Gamma_q(\tau) = \lim_{\delta \to 0} \Gamma_q(\tau, \delta).$$
 (A3)

The quantity $\Gamma_q(\tau)$ experiences a jump from 0 to $+\infty$, as τ is increased, at a critical value that we denote $\tau(q)$. The dimension \widetilde{D}_q of the measure is then defined as

$$\widetilde{D}_q = \tau(q)/(q-1). \tag{A4}$$

In practice, it is difficult to determine whether a particular covering $\{S_i\}$ is optimal in the sense of Eq. (A2). However, one can often compute the correct dimension by considering specific coverings, in the following sense. Consider a sequence of disjoint coverings $\{S_i^{(m)}\}$ where covering m of the sequence has a maximum diameter $\delta^{(m)}$ ($\epsilon_i^{(m)} < \delta^{(m)}$ for every component i of covering m) and $\delta^{(m)}$ converges to zero as m approaches infinity. If the limit

$$\hat{\Gamma}_q(\tau) = \lim_{m \to \infty} \Gamma_q(\tau, \{S_i^{(m)}\}, \delta^{(m)}) \tag{A5}$$

exists, then like $\Gamma_q(\tau)$ it experiences a jump from 0 to $+\infty$, as τ increases, at a critical value $\tau = \hat{\tau}(q)$. In terms of the sequence $\{S_i^{(m)}\}$ of coverings (not necessarily optimal), we can then compute a dimension spectrum

$$\hat{D}_q = \hat{\tau}(q)/(q-1). \tag{A6}$$

Since the sequence of coverings used to compute \hat{D}_q may be suboptimal, we have $\hat{\tau}(q) \ge \tau(q)$ for q > 1 and $\hat{\tau}(q) \le \tau(q)$ for q < 1. Thus $\hat{D}_q \ge \tilde{D}_q$ in each case. Assuming the limit in Eq. (1) that defines D_q to exist, then D_q can be computed by the partition function formalism above using

equal-size coverings $(\epsilon_i^{(m)} = \delta^{(m)})$ for every i and m). Thus $D_q \ge \widetilde{D}_q$ as well. While there do exist examples for which $\widetilde{D}_q \ne D_q$ for q < 1, the two dimensions are typically found to coincide in analytical examples with physical bases. Furthermore, for q > 1 the two dimensions always coincide.

Theorem. For every probability measure μ for which the limit in Eq. (1) exists, $\tilde{D}_q = D_q$ for all q > 1. (If one defines D_q as a liminf, then this equality holds for all μ .)

We prove this theorem in Appendix B.

The theorem demonstrates that, for q > 1 at least, the dimension spectrum computed according to the partition function formalism from a particular sequence of coverings is not that sensitive to the type of covering; equal-size coverings yield the same spectrum as a sequence of coverings that is optimal in the sense of Eq. (A2). While it may be possible to compute a different value of \hat{D}_q from a sufficiently suboptimal sequence of coverings, we conjecture that for the coverings we consider below,

$$\hat{D}_a = \tilde{D}_a = D_a. \tag{A7}$$

Furthermore, we conjecture that Eq. (A7) holds not only for q > 1, but for all $q \ge 0$.

For our purposes, the application of Eq. (A7) for $q \le 1$ is supported by the following reasoning. In all of the cases below, the formula we obtain for \hat{D}_q is an analytic function of q (for our region of interest $0 \le q \le 2$). If $\hat{D}_q = D_q$ for q > 1, it follows that if D_q is continuous at q = 1, then $\hat{D}_1 = D_1$. Furthermore, if D_q is analytic, then $\hat{D}_q = D_q$ for q < 1 as well.

2. Application to the models

The practical implication of Eq. (A7) for us is significant. It means that we can choose any sequence of coverings whose maximum diameter converges to zero, and this choice can be made in a manner that *facilitates analytic computations*. The result will be the same as that for an equal-cubesize covering [Eq. (1)] or an optimal covering [Eq. (A2)].

For model 1, we choose the intervals between the points at time t as a covering for the distribution. We regard each interval as covering the equivalent of one point since it contains two half points at each of its ends. Hence, on average,

 $\mu_i = [\Sigma_k N(k,t)]^{-1} = 1/N(t)$ for all i, and Eq. (A5) becomes²

$$\hat{\Gamma}_{q}(\tau) = \lim_{t \to \infty} [1/N(t)]^{q} \sum_{i} \epsilon_{i}^{-\tau}, \tag{A8}$$

where the $\delta \to 0$ limit is replaced by $t \to \infty$, since, as t increases, the size of the largest interval decreases to zero with probability 1. Since $\ln \hat{\Gamma}_q(\tau) = -\infty$ for $\tau < \hat{\tau}(q)$ and $\ln \hat{\Gamma}_q(\tau) = \infty$ for $\tau > \hat{\tau}(q)$, we equate $\ln \hat{\Gamma}_q(\tau)$ to zero and obtain q as a function of the transition value $\hat{\tau}$,

$$q(\hat{\tau}) = \lim_{t \to \infty} \frac{\ln \sum_{i} \epsilon_{i}^{-\hat{\tau}}}{\ln N(t)} = \lim_{t \to \infty} \frac{\ln \sum_{k} N(k, t) 2^{k\hat{\tau}}}{\ln N(t)}.$$
 (A9)

We then have [see Eq. (9)]

$$q(\hat{\tau}) = 1 + \lim_{t \to \infty} \frac{\ln\langle 2^{k\hat{\tau}} \rangle_t}{\ln N(t)},\tag{A10}$$

where

$$\langle 2^{k\tau} \rangle_t = \frac{\sum_k N(k,t) 2^{k\tau}}{N(t)}.$$
 (A11)

Each square in model 2 and triangle in model 3 contains, on average, $[\Sigma_k N(k,t)]^{-1} = 1/N(t)$ of the total measure. Thus, like we derived for model 1, we obtain

$$q(\hat{\tau}) = 1 + \lim_{t \to \infty} \frac{\ln \langle 2^{k\hat{\tau}} \rangle_t}{\ln N(t)}.$$
 (A12)

For model 4 we use the two types of squares (types I and II) to cover the distribution. A type-I square has one point on one of its vertices and hence covers the equivalent of a quarter of a point, while a type-II square, with two points on its vertices, covers the equivalent of half a point. Thus, since at time t we have t total points, the measure contained in a type-I square is 1/4t, while that contained in a type-II square is 1/2t. Thus Eq. (A5) becomes

$$\hat{\Gamma}_{q}(\tau) = \lim_{t \to \infty} \frac{1}{(4t)^{q}} \left(\sum_{i=1}^{N_{1}} \epsilon_{i}^{-\tau} + 2^{q} \sum_{i=1}^{N_{2}} \epsilon_{i}^{-\tau} \right), \tag{A13}$$

where the first summation is over all type-I squares and the second is over all type-II squares. Taking the logarithm of this expression and equating it to zero, we obtain

$$q(\hat{\tau}) = \lim_{t \to \infty} \frac{\ln[T_1(t) + 2^q T_2(t)]}{\ln t},$$
 (A14)

where $T_i(t) = \sum_k 2^{k\hat{\tau}} N_i(k,t)$.

 $^{^{1}}$ As a simple example, we consider the set $1,1/2,1/3,\ldots,n^{-1},\ldots$, and q=0. For this set it can be shown that Eqs. (A1)–(A4) yield $\widetilde{D}_{0}=0$, which is intuitively reasonable for a set that is a countable collection of points. In contrast, Eqs. (1) and (2) yield $D_{0}=1/2$. To see this we first note that, for q=0, Eqs. (1) and (2) give the well-known result $D_{0}=\lim_{\epsilon\to 0}\{[\ln N(\epsilon)]/[\ln(1/\epsilon)]\}$, where $N(\epsilon)$ is the number of ϵ intervals needed to cover the set. Next we observe that the distance between 1/n and 1/(n+1) is approximately $1/n^{2}$ for large n. Thus setting $\epsilon=1/n^{2}$, we need n intervals to cover the first n elements of the set, one interval for each such element. To cover the remaining elements we must cover the interval (1/(n+1),0). This requires $(1/\epsilon)[1/(n+1)]=n^{2}/(n+1) \approx n$ intervals. Thus $N(\epsilon) \approx 2n=2/\epsilon^{1/2}$, yielding $D_{0}=1/2$.

²There is a subtlety here. As *t* increases we are changing *both* the covering *and* the distribution of points. We really should be looking at the limiting distribution and not the time *t* distribution. Since each interval is equally likely to subdivide, the *expected* measure in each of the time *t* intervals is the same, and our computations in this section reflect this approximation.

APPENDIX B: PROOF OF THEOREM FROM APPENDIX A

As we noted in Appendix A, $\tilde{D}_q \leq D_q$ for all $q \geq 0$, so we now show that also $\tilde{D}_q \geq D_q$ for q > 1. Our proof is based on the theorem in [14] that for q > 1, the quantity

$$I_q(s) = \int \left(\int \frac{d\mu(y)}{|x - y|^s} \right)^{q-1} d\mu(x)$$
 (B1)

is finite for $s < D_q$ and infinite for $s > D_q$. To see that $\widetilde{D}_q \ge D_q$, we show that $\widetilde{D}_q \ge s$ whenever $I_q(s)$ is finite. Let $\{S_i\}$ be a disjoint covering of μ , as in Appendix A, and recall that μ_i is the measure of S_i and ϵ_i is its diameter. Then considering only the contribution to $I_q(s)$ from points s and s that are in the same s, we have

by Eq. (A1). This implies that $\Gamma_q((q-1)s, \delta) \leq I_q(s)$ by Eq. (A2), since q > 1, and hence $\Gamma_q((q-1)s) \leq I_q(s)$ by Eq. (A3). Therefore, if $I_q(s)$ is finite, then $\Gamma_q((q-1)s)$ is finite, whence $\tau(q) \geq (q-1)s$, and finally $\tilde{D}_q \geq s$ by Eq. (A4).

 $I_{q}(s) \ge \sum_{i=1}^{N} \int_{S_{i}} \left(\int_{S_{i}} \frac{d\mu(y)}{|x-y|^{s}} \right)^{q-1} d\mu(x)$ $\ge \sum_{i=1}^{N} \int_{S_{i}} \left(\int_{S_{i}} \frac{d\mu(y)}{\epsilon_{i}^{s}} \right)^{q-1} d\mu(x)$ $= \sum_{i=1}^{N} \mu_{i} \left(\frac{\mu_{i}}{\epsilon_{i}^{s}} \right)^{q-1} = \sum_{i=1}^{N} \mu_{i}^{q} / \epsilon_{i}^{(q-1)s}$ $= \Gamma_{q}((q-1)s, \{S_{i}\}, \delta)$

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