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Ínría

A generator of random convex polygons in a disc

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Un générateur de polygones convexes aléatoires dans un disque

Résumé : Nous proposons un algorithme qui génère un polygone aléatoire défini par l'enveloppe convexe de n points aléatoires indpéndants et uniformément distribués dans le disque, sans avoir à générer explicitement tous les points.

Mots-clés : Enveloppe convexe, polygone aléatoire, générateur

1 Introduction

Let \mathcal{D} be a disc in \mathbb{R}^2 with radius R centered at \mathfrak{o} , and (x_1, \ldots, x_n) a sample of n points uniformly and independently distributed in \mathcal{D} . Let's define the polygon P_n as the convex hull of (x_1, \ldots, x_n) , and $f_0(P_n)$ its number of vertices. This kind of polygon has been well studied, and using [Bá92, Rei10] one can easily see that

$$\mathbb{E}f_0(P_n) = c \ n^{\frac{1}{3}} + o(n^{\frac{1}{3}}) \tag{1}$$

where $c = 2^{\frac{4}{3}} 3^{-\frac{1}{3}} \Gamma(\frac{5}{3}) \pi^{\frac{5}{3}} \approx 3.383228964$ and Γ denotes the usual Gamma function.

To generate such a polygon, one can explicitly generate n points uniformly in \mathcal{D} and compute the convex hull. For a very large number of points, it could be interesting to generate fewer points to get the same polygon, for example to evaluate the constant that are not explicitly known for the asymptotic distribution of the perimeter, or other parameters such has the higher moments of the extremal points.

In this note, we propose an algorithm that generates far fewer points at random in order to get P_n , so that the time and the memory needed is reduced for n large.

2 Algorithm

We start with random polygon P_i in \mathcal{D} where *i* is very small, and we increase the number of points until we get P_n .

The idea is that given the convex hull of small number of points, the number of points generated in \mathcal{D} that are deeply inside (and so won't change the convex hull) is a random variable that we can easily simulate, so that we need to generate only the small number of points that are close to the convex hull.

The outline of the algorithm is the following :

- Generate a small number of points in \mathcal{D} and compute its convex hull;
- Compute the radius of the largest disc centered at o inscribed in the convex hull;
- Choose a number of points to simulate at this step;
- Simulate the number of points that fall in this inscribed disc at this step;
- Generate the rest of the random points in the annulus defined by these two discs and update the convex hull.
- continue this process until the sum of the simulated and generated points is equal to n.

To simplify the notations, we assume \mathcal{D} to have radius 1.

Notations

- n is the total number of points simulated and m_i is the total number of points from step 1 to step i;
- k_i is the number of generated points at step *i*, and $k = \sum_i k_i$ is the total number of generated points;

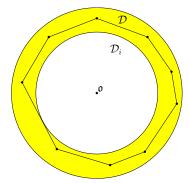


Figure 1: At step i, we simulate the number of points that falls in \mathcal{D}_i and we generate points uniformly in the yellow annulus

- h_i is size of the convex hull at step *i*, and *h* is the size of the final convex hull.
- p_i is the probability to fall in the annulus at step *i*.

Initialization First, we have to generate a random polytope P_i with a small number of points in \mathcal{D} such that \mathfrak{o} is inside the random polytope. This is not too much to ask, as (see [AS08])

$$\mathbb{P}(\mathfrak{o} \notin P_i) = 2^{-(i-1)}i \tag{2}$$

We initialize the random polygon by generating 100 points in the disc. As the probability that $\mathfrak{o} \notin P_{100}$ is lower than 1.6×10^{-28} , it's very unlikely that this is not enough. Otherwise, we add another sample of 100 points, until $\mathfrak{o} \in P_{100,j}$ for some j.

Simulation of Points At the beginning of each step *i* of the loop, we are given a polygon P_{m_i} , which is the convex hull of m_i points. Let s_i the number of new points simulated at step *i*. We choose $s_i = m_i$ if $m_i < n \log^{-2} n$, and $s_i = n \log^{-2} n$ otherwise. Let \mathcal{D}_i be the largest inscribed disc in P_{m_i} centered in \mathfrak{o} , and r_i its radius.

Using a simulation of a binomial variable of parameter s_i and $1 - r_i^2$, we can evaluate the number of points that falls in $\mathcal{D} \setminus \mathcal{D}_i$. As the points in \mathcal{D}_i won't change the convex hull, we don't need to generate them. Then, we generate the rest of the points uniformly in the annulus $\mathcal{D} \setminus \mathcal{D}_i$, and we update the convex hull using a Graham scan.

Generation of random points To generate random points in an annulus with radii r_i and 1, one need to generate the polar angles uniformly in $[-\pi, \pi)$ and the squared radii uniformly in $[r_i^2, 1)$.

As we want to perform a Graham scan in linear time, the points have to be sorted by their polar angles. This can be done in expected linear time and size, using a bucket sort, as the angles are uniformly chosen, see [CSRL01, 8.4].

Full algorithm

Data: integer n**Result**: Convex hull of n uniformly chosen points in the disc \mathcal{D} Simulated Points $\leftarrow 0$; do Generate $\min(100, n - Simulated_Points)$ points in the disc of radius 1; Simulated Points \leftarrow Simulated Points+min(100, n - Simulated Points); while \mathfrak{o} is not in the convex hull and Simulated Points < n; while Simulated Points < n do Compute inscribed radius; $p \leftarrow inscribed \ radius^2;$ if Simulated Points $< n \log^{-2} n$ then $k \leftarrow Simulated Points;$ else $k \leftarrow \min(\lfloor n \log^{-2} n \rfloor, n - Simulated_points);$ end $X \leftarrow$ Simulation of Binomial variable with parameters k, 1-p;Generate X points uniformly and sorted in the annulus of radii *inscribed* radius, 1; Simulated Points \leftarrow Simulated Points +k; Merge the list of the convex hull and the new points; Update the convex hull with a Graham scan on the list; end return Convex hull

Algorithm 1: Algorithm of the Generator of Random Polygon in a disc

3 Complexity

Clearly the size complexity is $\max_i(h_i + k_i)$ and the time complexity is $\sum_i(h_i + k_i)$ since the Graham scan and the points generation are linear in the number of points [Gra72].

For the initialization, as the probability that $\mathfrak{o} \notin P_n$ decreases exponentially, it is very unlikely that more than one loop is necessary. Let's call \mathfrak{p} the minimal number of points such that $\mathfrak{o} \in P_{\mathfrak{p}}$. Using formula (2), the expectation of \mathfrak{p} is very small :

$$\mathbb{E}(\mathfrak{p}) = \sum_{j=1}^{\infty} j \mathbb{P}(\mathfrak{p} = j) = \sum_{j=1}^{\infty} \mathbb{P}(\mathfrak{p} \ge j)$$
$$= \sum_{i=1}^{3} \mathbb{P}(\mathfrak{p} \ge j) + \sum_{j=4}^{\infty} \mathbb{P}(\mathfrak{o} \not\in P_{j-1})$$
$$= 3 + \sum_{j=3}^{\infty} \mathbb{P}(\mathfrak{o} \not\in P_j)$$
$$= 3 + \sum_{i=3}^{\infty} j 2^{-(j-1)} = 3 + 2 = 5.$$
(3)

Thus, the expected size and time complexity of the initialization is O(1).

RR n° 8467

For i big enough, we have [Bár89]:

$$\mathbb{E} d_H(P_{m_i}, \mathcal{D}) = \Theta\left(\frac{\log m_i}{m_i}\right)^{\frac{2}{3}}$$
(4)

where d_H , the Hausdorff distance, is the maximal distance between a point in P_i and the boundary of \mathcal{D} .

Recall that \mathcal{D}_i is the annulus with radii $r_i = 1 - d_H(P_{m_i}, \mathcal{D})$ and 1 and let p_i be the probability that a random point in \mathcal{D} falls in \mathcal{D}_i Using (4), there exist a constant $c_0 > 0$ such that, for i large $\mathbb{E}(p_i) = 1 - r_i^2 < 2(1 - r_i) < c_0 \left(\frac{\log m_i}{m_i}\right)^{\frac{2}{3}}$ Let' s call i_{τ} the last step *i* where $m_i < \frac{n}{\log^2 n}$.

At each step $i \leq i_{\tau}$, k_i is a binomial variable with parameter p_i and $m_i < \frac{n}{\log^2 n}$. Thus, for i large enough,

$$\mathbb{E}(k_i) = \mathbb{E}(\mathbb{E}(k_i|p_i)) = m_i \mathbb{E}p_i = O(m_i^{\frac{1}{3}} \log^{\frac{2}{3}}(m_i)) = O\left(\left(\frac{n}{\log^2 n}\right)^{\frac{1}{3}} \log^{\frac{2}{3}}n\right) = O(n^{\frac{1}{3}}).$$
(5)

As we choose $m_{i+1} = 2m_i$, i_{τ} is bounded by $\log_2(n)$.

For $i > i_{\tau}$, k_i is a binomial variable with parameter p_i and $\frac{n}{\log^2 n}$, so using (5) and the fact that $m_i > \frac{n}{\log^2 n}$, we get $\mathbb{E}k_i = O(n^{\frac{1}{3}})$. As we simulate at each step $i > i_{\tau}$ at least $\frac{n}{\log^2(n)}$, the number of step after i_{τ} is bounded by $\log^2(n)$.

Now,

$$\mathbb{E}k = \sum_{i=1}^{i_{\tau}} \mathbb{E}k_i + \sum_{i>i_{\tau}} \mathbb{E}k_i$$

$$\leq O\left(\log^{\frac{2}{3}} n\left(\sum_{i=1}^{\log_2 n} m_i^{\frac{1}{3}}\right)\right) + \log^2 n \ O(n^{\frac{1}{3}})$$

$$\leq O\left(\log^{\frac{2}{3}} n\left(\sum_{i=1}^{\log_2 n} (2^i)^{\frac{1}{3}}\right)\right) + O(n^{\frac{1}{3}}\log^2 n) = O(n^{\frac{1}{3}}\log^2 n)$$
(6)

At each step *i*, the expected size of the convex hull is $O(m_i^{\frac{1}{3}}) = O(n^{\frac{1}{3}})$, and $\mathbb{E}k_i = O(n^{\frac{1}{3}})$. Thus, the expected size complexity is $O(n^{\frac{1}{3}})$.

As our points are sorted according to their polar angle, computing the convex hull with a Graham scan is done in linear time $(O(n^{\frac{1}{3}} \log^2 n))$, the generation of the k points and the computation of the largest annulus as well $(O(n^{\frac{1}{3}}))$. Thus, the expected time complexity is $O(n^{\frac{1}{3}}\log^2 n)$.

Experiments 4

This algorithm has been implemented in C++, and will be submitted for integration in the CGAL library.

As expected, the distribution of $\mathbb{E}f_0(P_n)$ is asymptotically the same as the theoretical one, see equation (1), with the same constant :

This estimation has been done one 1000 experiments for each value of $\log_{10} n$.

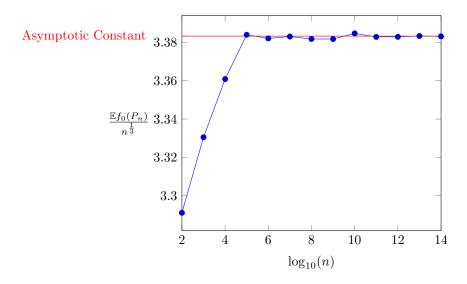


Figure 2: Average size of P_n divided by $n^{\frac{1}{3}}$

Complexities To evaluate the size complexity, we just compute the largest list of points used in the loop, see Figure 3. The time complexity can be evaluated by estimating the total number of points generated, see Figure 4.

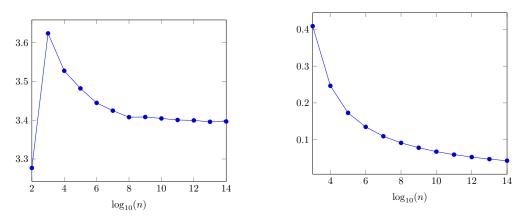


Figure 3: Average maximal size divided by $n^{\frac{1}{3}}$ Figure 4: Average number of generated points divided by $n^{\frac{1}{3}} \log^2 n$

As a first application, we can estimate the distribution of the smallest and the largest edge of a random polytope, see Figure 5 and Figure 6. The average have been done on 100 experiments for each value of $\log_{10} n$. The expected minimal edge seems to be $O(n^{-\frac{1}{3}})$, and the maximal $O\left(\frac{\log^2 n}{n}\right)^{\frac{1}{3}}$.

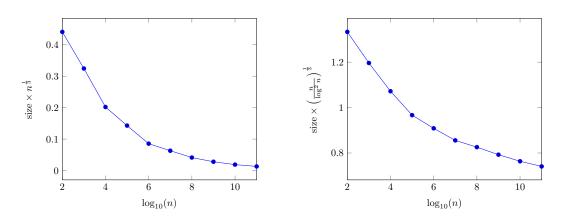


Figure 5: Average length of the minimal edge Figure 6: Average length of the maximal edge

5 Conclusion

We propose an algorithm that generates random polygons given by the convex hull of random points, without generating all the points. There is no theoretical obstacle to generalize to higher dimension. The theoretical results used for the evaluation of the complexity are known in arbitrary dimension [Bár89, Rei10], so the analysis can be done as well.

We can reduce the expected time complexity by a logarithmic factor if we allow to increase the expected size complexity by a logarithmic factor. Instead of bounding the simulated points at step i by $\frac{n}{\log^2 n}$ when m_i becomes bigger than $\frac{n}{\log^2 n}$, we can choose to always simulate m_i points. In this case,

$$\mathbb{E}k_i = O(m_i^{\frac{1}{3}} \log^{\frac{2}{3}} m_i) = O(n^{\frac{1}{3}} \log^{\frac{2}{3}} n)$$
(7)

so the expected size complexity becomes $O(n^{\frac{1}{3}} \log^{\frac{2}{3}} n)$. On the other hand, the expected time complexity is reduced to $O(n^{\frac{1}{3}} \log^{\frac{2}{3}} n)$, as

$$\mathbb{E}k = \sum_{i} \mathbb{E}k_{i}$$
$$= O\left(\log^{\frac{2}{3}} n\left(\sum_{i=1}^{\log_{2} n} m_{i}^{\frac{1}{3}}\right)\right) = O\left(n^{\frac{1}{3}}\log^{\frac{2}{3}} n\right).$$

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