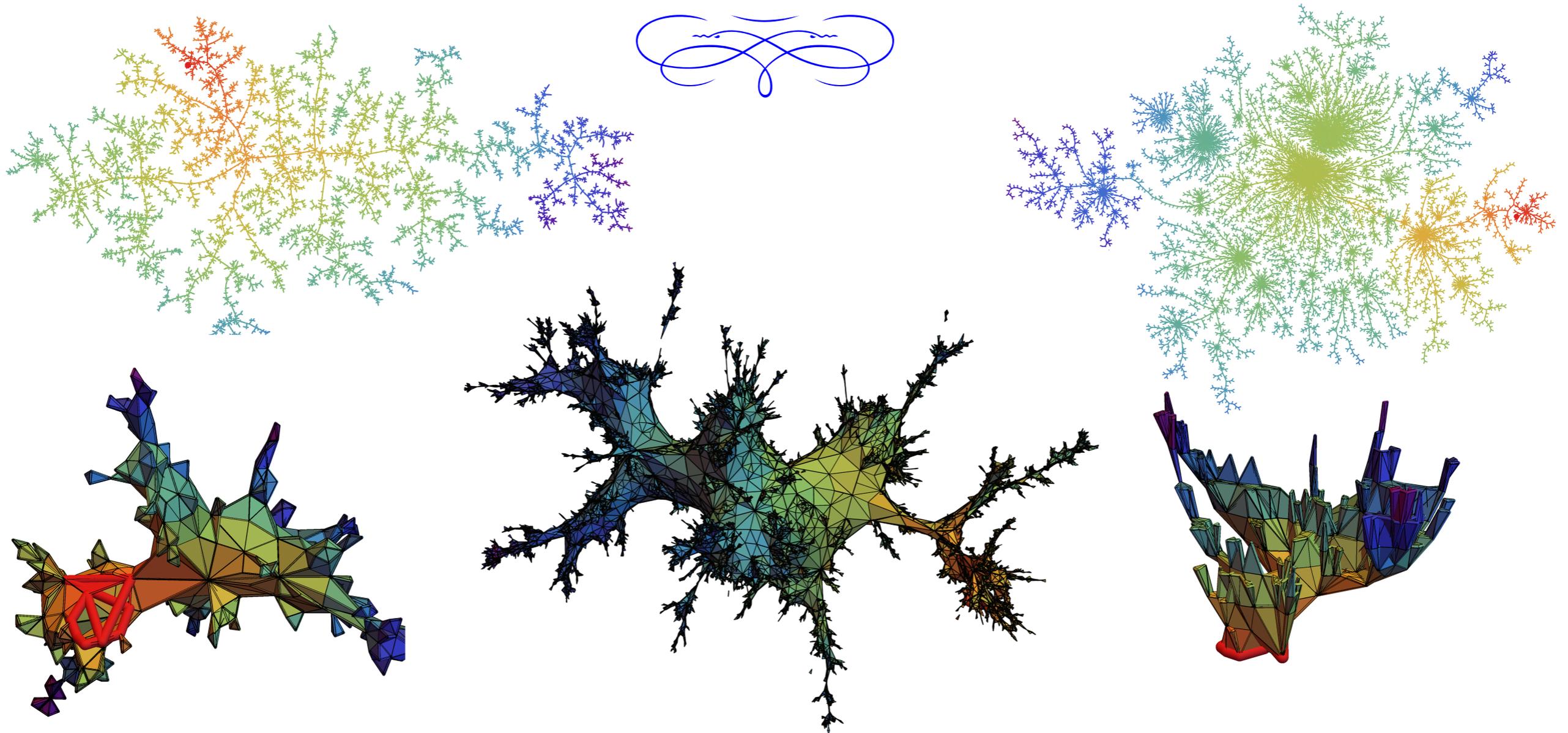


Self-similar growth-fragmentations & random planar maps



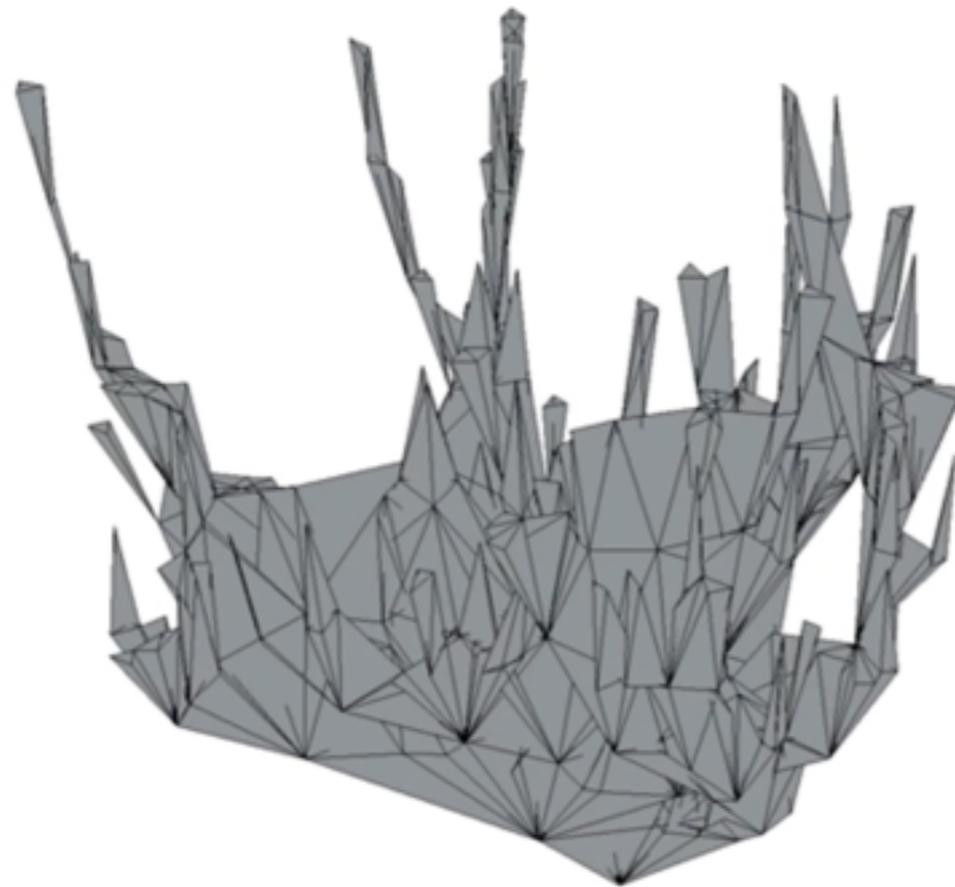
Igor Kortchemski (joint work with J. Bertoin, T. Budd, N. Curien)
CNRS & École polytechnique

Goal

→ **Goal:** study a random surface by studying its level sets.

Goal

↗ **Goal:** study a random surface by studying its level sets.



Goal

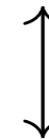
→ **Goal:** study a random surface by studying its level sets.



Goal

↗ **Goal:** study a random surface by studying its level sets.

Random surface



Level sets of
a random surface

Goal

↗ **Goal:** study a random surface by studying its level sets.

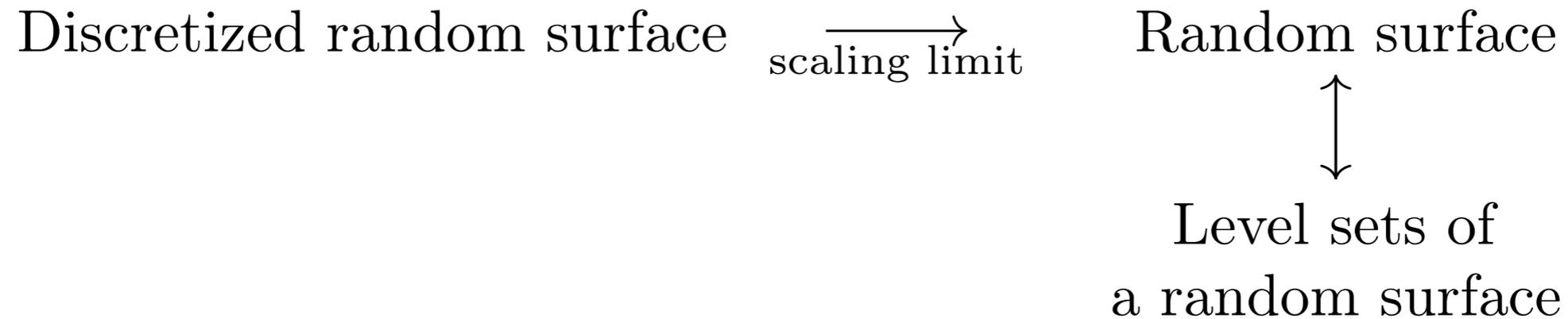
Approach from the discrete side.

Random surface
↕
Level sets of
a random surface

Goal

↗ **Goal:** study a random surface by studying its level sets.

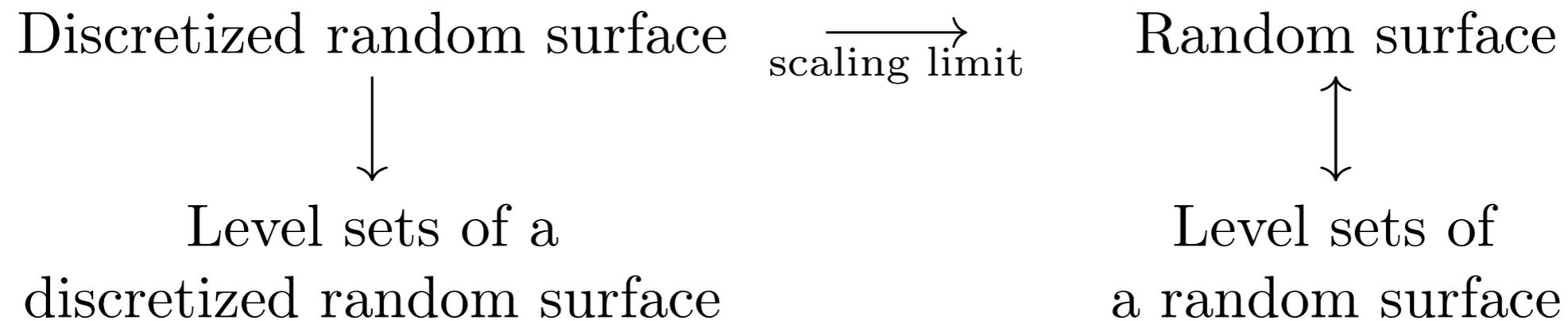
Approach from the discrete side.



Goal

↗ **Goal:** study a random surface by studying its level sets.

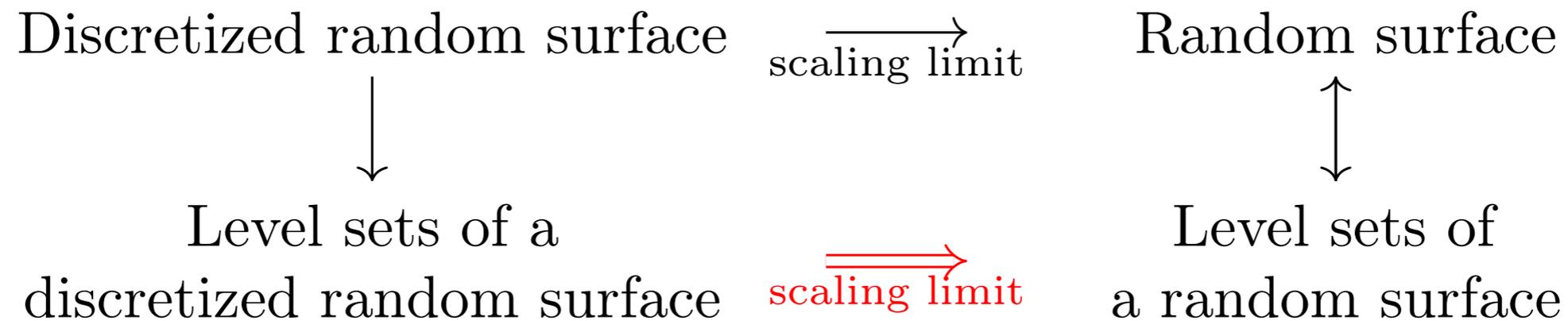
Approach from the discrete side.



Goal

↗ **Goal:** study a random surface by studying its level sets.

Approach from the discrete side.



Goal

↗ **Goal:** study a random surface by studying its level sets.

Approach from the discrete side.

Discretized random surface



Level sets of a
discretized random surface

scaling limit

?

Motivation for studying scaling limits

Let $(X_n)_{n \geq 1}$ be “discrete” objects converging towards a “continuous” object X :

$$X_n \xrightarrow[n \rightarrow \infty]{} X.$$

Motivation for studying scaling limits

Let $(X_n)_{n \geq 1}$ be “discrete” objects converging towards a “continuous” object X :

$$X_n \xrightarrow[n \rightarrow \infty]{} X.$$

Several consequences:

- *From the discrete world to the continuous world:* if a property \mathcal{P} is satisfied by all the X_n and passes to the limit, then X satisfies \mathcal{P} .

Motivation for studying scaling limits

Let $(X_n)_{n \geq 1}$ be “discrete” objects converging towards a “continuous” object X :

$$X_n \xrightarrow[n \rightarrow \infty]{} X.$$

Several consequences:

- *From the discrete world to the continuous world:* if a property \mathcal{P} is satisfied by all the X_n and passes to the limit, then X satisfies \mathcal{P} .
- *From the continuous world to the discrete world:* if a property \mathcal{P} is satisfied by X and passes to the limit, X_n satisfies “approximately” \mathcal{P} for n large.

Motivation for studying scaling limits

Let $(X_n)_{n \geq 1}$ be “discrete” objects converging towards a “continuous” object X :

$$X_n \xrightarrow[n \rightarrow \infty]{} X.$$

Several consequences:

- *From the discrete world to the continuous world:* if a property \mathcal{P} is satisfied by all the X_n and passes to the limit, then X satisfies \mathcal{P} .
- *From the continuous world to the discrete world:* if a property \mathcal{P} is satisfied by X and passes to the limit, X_n satisfies “approximately” \mathcal{P} for n large.
- *Universality:* if $(Y_n)_{n \geq 1}$ is another sequence of objects converging towards X , then X_n and Y_n share approximately the same properties for n large.

Motivation for studying scaling limits

Let $(X_n)_{n \geq 1}$ be “discrete” objects converging towards a “continuous” object X :

$$X_n \xrightarrow[n \rightarrow \infty]{} X.$$

Several consequences:

- *From the discrete world to the continuous world:* if a property \mathcal{P} is satisfied by all the X_n and passes to the limit, then X satisfies \mathcal{P} .
- *From the continuous world to the discrete world:* if a property \mathcal{P} is satisfied by X and passes to the limit, X_n satisfies “approximately” \mathcal{P} for n large.
- *Universality:* if $(Y_n)_{n \geq 1}$ is another sequence of objects converging towards X , then X_n and Y_n share approximately the same properties for n large.

What is the sense of the convergence when the objects are **random**?

Motivation for studying scaling limits

Let $(X_n)_{n \geq 1}$ be “discrete” objects converging towards a “continuous” object X :

$$X_n \xrightarrow[n \rightarrow \infty]{} X.$$

Several consequences:

- *From the discrete world to the continuous world:* if a property \mathcal{P} is satisfied by all the X_n and passes to the limit, then X satisfies \mathcal{P} .
- *From the continuous world to the discrete world:* if a property \mathcal{P} is satisfied by X and passes to the limit, X_n satisfies “approximately” \mathcal{P} for n large.
- *Universality:* if $(Y_n)_{n \geq 1}$ is another sequence of objects converging towards X , then X_n and Y_n share approximately the same properties for n large.

What is the sense of the convergence when the objects are **random**?

 **Convergence in distribution in a certain metric space.**

Outline

I. PLANAR MAPS



II. BIENAYMÉ–GALTON–WATSON TREES

III. RANDOM MAPS AND GROWTH-FRAGMENTATIONS

MOTIVATION



Motivation

What does a typical **random surface** look like?

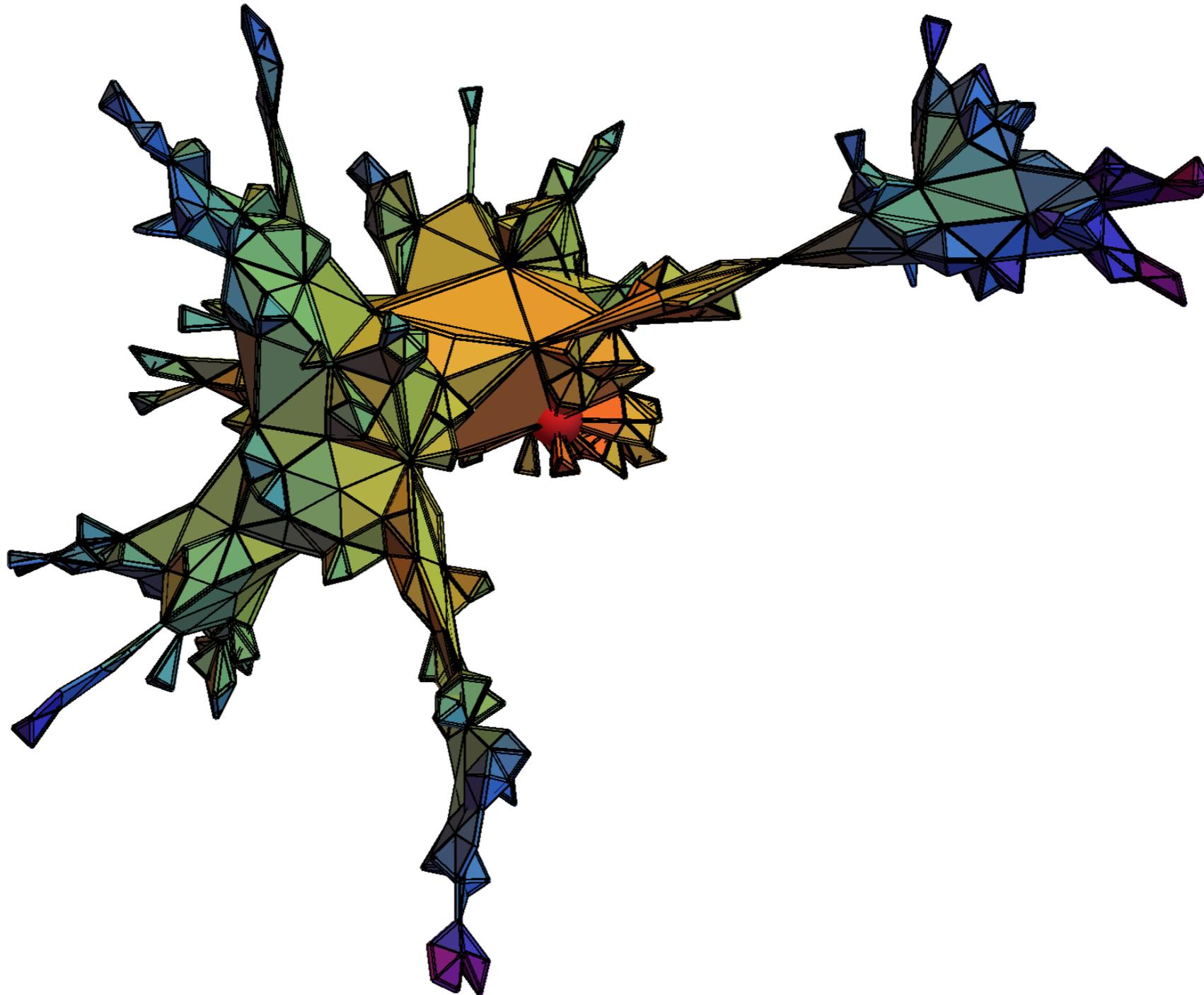
→ **Idea:** construct a **random surface** as a limit of random **discrete surfaces**.

→ **Idea**: construct a **random surface** as a limit of random **discrete surfaces**.

Consider n **triangles**, and glue them together at random to obtain a surface homeomorphic to a **sphere**.

→ **Idea**: construct a **random surface** as a limit of random **discrete surfaces**.

Consider n **triangles**, and glue them together at random to obtain a surface homeomorphic to a **sphere**.



↗ Idea: construct a **random surface** as a limit of random **discrete surfaces**.

Consider n **triangles**, and glue them together at random to obtain a surface homeomorphic to a **sphere**.

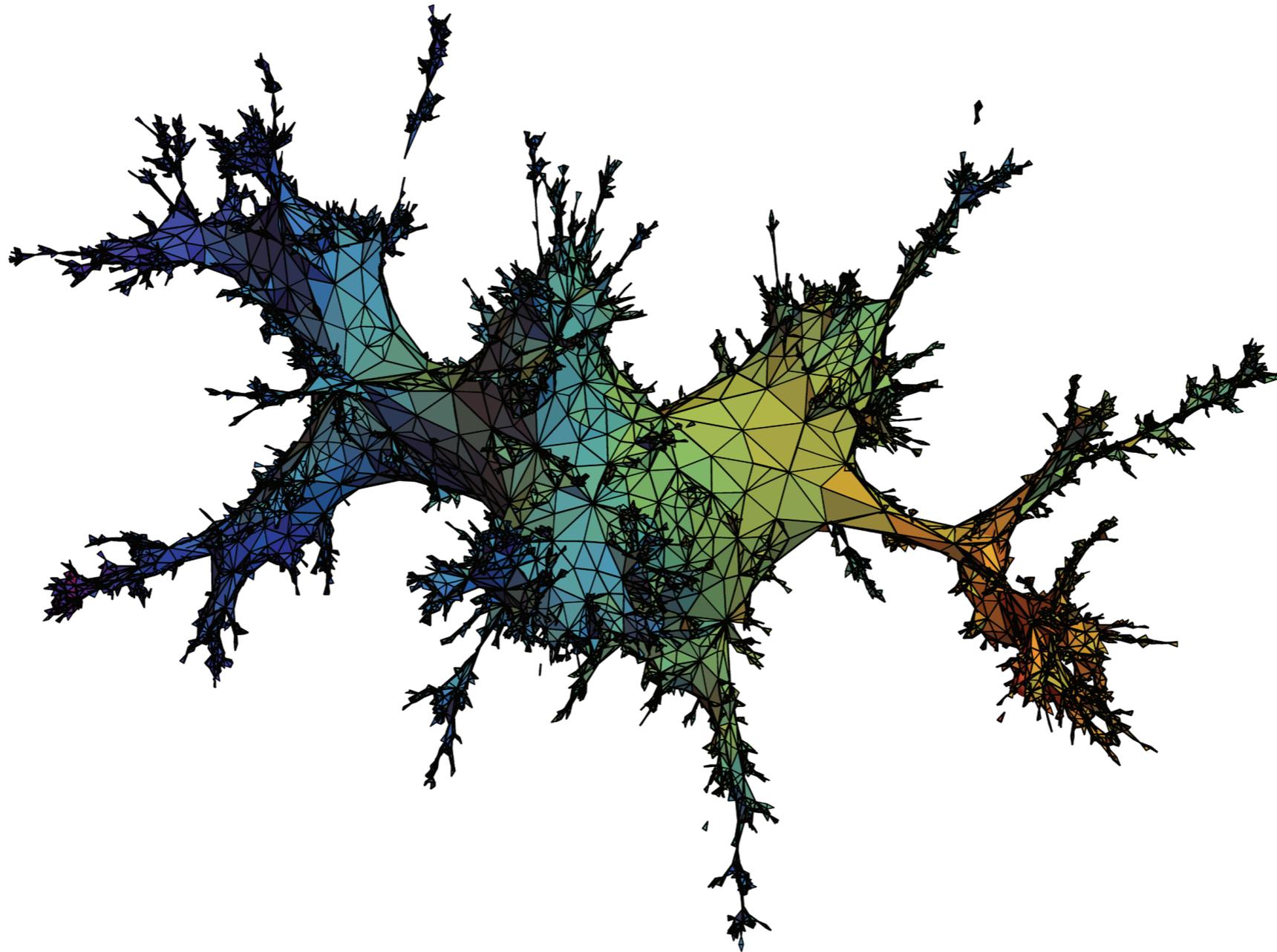


Figure: A large **random triangulation**

The Brownian map

Problem (Schramm, ICM '06): Let T_n be a random uniform triangulation of the sphere with n triangles.

The Brownian map

Problem (Schramm, ICM '06): Let T_n be a **random uniform triangulation** of the sphere with n triangles. View T_n as a compact metric space, by equipping its vertices with the graph distance.

The Brownian map

Problem (Schramm, ICM '06): Let T_n be a **random uniform triangulation** of the sphere with n triangles. View T_n as a compact metric space, by equipping its vertices with the graph distance. Show that $n^{-1/4} \cdot T_n$ converges to a random compact metric space homeomorphic to the sphere (**the Brownian map**)

The Brownian map

Problem (Schramm, ICM '06): Let T_n be a **random uniform triangulation** of the sphere with n triangles. View T_n as a compact metric space, by equipping its vertices with the graph distance. Show that $n^{-1/4} \cdot T_n$ converges to a random compact metric space homeomorphic to the sphere (**the Brownian map**), in distribution for the **Gromov–Hausdorff** topology.

The Brownian map

Problem (Schramm, ICM '06): Let T_n be a **random uniform triangulation** of the sphere with n triangles. View T_n as a compact metric space, by equipping its vertices with the graph distance. Show that $n^{-1/4} \cdot T_n$ converges to a random compact metric space homeomorphic to the sphere (**the Brownian map**), in distribution for the **Gromov–Hausdorff** topology.

Solved by **Le Gall** (as well as for other families of maps including quadrangulations) in 2011, and independently by **Miermont** in 2011 for quadrangulations.

The Brownian map

Problem (Schramm, ICM '06): Let T_n be a **random uniform triangulation** of the sphere with n triangles. View T_n as a compact metric space, by equipping its vertices with the graph distance. Show that $n^{-1/4} \cdot T_n$ converges to a random compact metric space homeomorphic to the sphere (**the Brownian map**), in distribution for the **Gromov–Hausdorff** topology.

Solved by **Le Gall** (as well as for other families of maps including quadrangulations) in 2011, and independently by **Miermont** in 2011 for quadrangulations.

Since, convergence to the **Brownian map** has been established for many different models of random maps (**Beltran & Le Gall, Addario-Berry & Albenque, Bettinelli, Bettinelli & Jacob & Miermont, Abraham, Bettinelli & Miermont, Baur & Miermont & Ray**)

The Brownian map

Problem (Schramm, ICM '06): Let T_n be a **random uniform triangulation** of the sphere with n triangles. View T_n as a compact metric space, by equipping its vertices with the graph distance. Show that $n^{-1/4} \cdot T_n$ converges to a random compact metric space homeomorphic to the sphere (**the Brownian map**), in distribution for the **Gromov–Hausdorff** topology.

Solved by **Le Gall** (as well as for other families of maps including quadrangulations) in 2011, and independently by **Miermont** in 2011 for quadrangulations.

Since, convergence to the **Brownian map** has been established for many different models of random maps (**Beltran & Le Gall, Addario-Berry & Albenque, Bettinelli, Bettinelli & Jacob & Miermont, Abraham, Bettinelli & Miermont, Baur & Miermont & Ray**), using different techniques, such as bijections with labeled trees (**Cori–Vauquelin–Schaeffer, Bouttier–Di Francesco–Guitter**).

↗ Other motivations:

- connections with 2D Liouville Quantum Gravity ([David](#), [Duplantier](#), [Garban](#), [Kupianen](#), [Maillard](#), [Miller](#), [Rhodes](#), [Sheffield](#), [Vargas](#), [Zeitouni](#)).
- study of random planar maps decorated with statistical physics models ([Angel](#), [Berestycki](#), [Borot](#), [Bouttier](#), [Guitter](#), [Chen](#), [Curien](#), [Gwynne](#), [K.](#), [Kassel](#), [Laslier](#), [Mao](#), [Ray](#), [Richier](#), [Sheffield](#), [Sun](#), [Wilson](#)).

LEVEL SETS OF THE BROWNIAN MAP

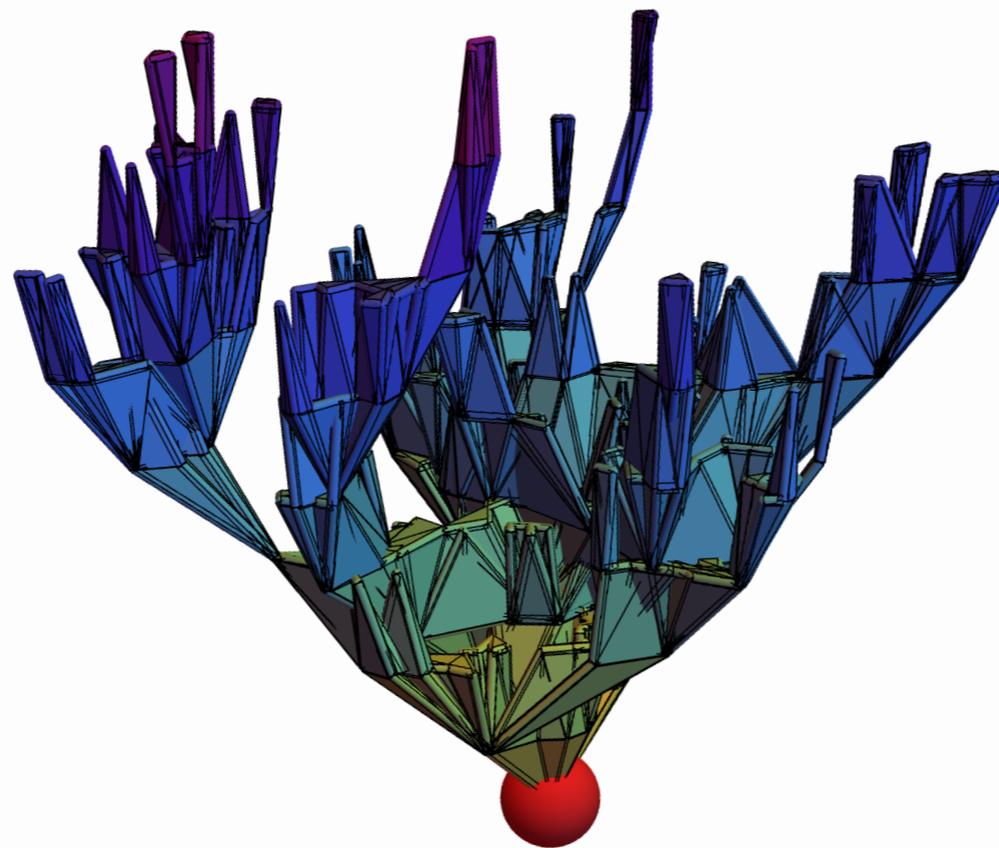


Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x .

Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x .



Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x . Now, for every $h > 0$, remove all the points which are not in the ball of radius h centered at the root, and look at the lengths of the cycles as h grows (level set process).



Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x . Now, for every $h > 0$, remove all the points which are not in the ball of radius h centered at the root, and look at the lengths of the cycles as h grows (level set process).

↗ **Questions** (related to the “breadth-first search” of the **Brownian map** of **Miller & Sheffield**):

Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x . Now, for every $h > 0$, remove all the points which are not in the ball of radius h centered at the root, and look at the lengths of the cycles as h grows (level set process).

↗ **Questions** (related to the “breadth-first search” of the **Brownian map** of **Miller & Sheffield**):

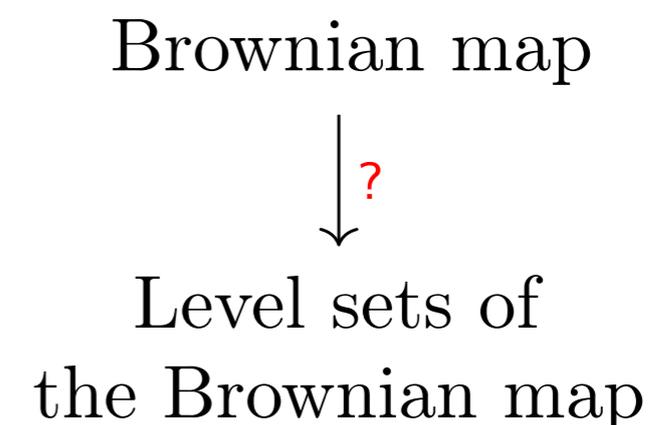
- What is the law of the level set process of the **Brownian map** as h grows?

Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x . Now, for every $h > 0$, remove all the points which are not in the ball of radius h centered at the root, and look at the lengths of the cycles as h grows (level set process).

↗ **Questions** (related to the “breadth-first search” of the **Brownian map** of **Miller & Sheffield**):

- What is the law of the level set process of the **Brownian map** as h grows?

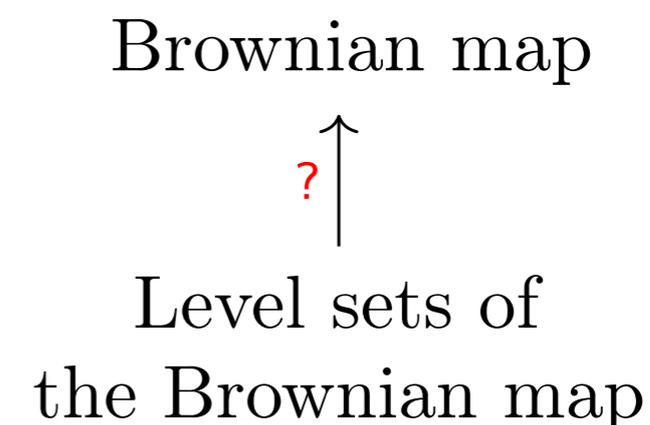


Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x . Now, for every $h > 0$, remove all the points which are not in the ball of radius h centered at the root, and look at the lengths of the cycles as h grows (level set process).

↗ **Questions** (related to the “breadth-first search” of the **Brownian map** of **Miller & Sheffield**):

- What is the law of the level set process of the **Brownian map** as h grows?
- Can one reconstruct the **Brownian map** from the level set processes?



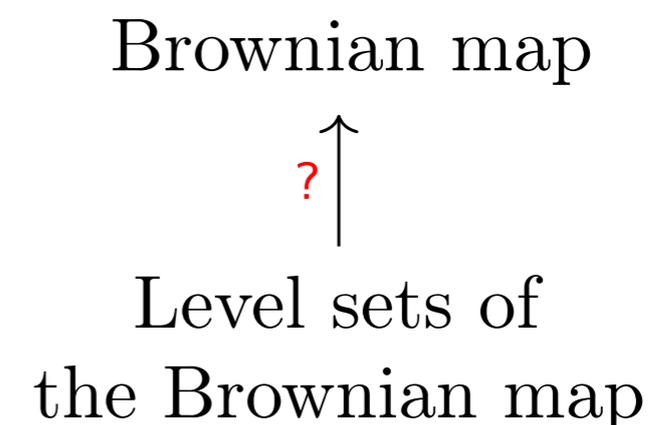
Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x . Now, for every $h > 0$, remove all the points which are not in the ball of radius h centered at the root, and look at the lengths of the cycles as h grows (level set process).

↗ **Questions** (related to the “breadth-first search” of the **Brownian map** of **Miller & Sheffield**):

- What is the law of the level set process of the **Brownian map** as h grows?
- Can one reconstruct the **Brownian map** from the level set processes?

↗ **Our result:** scaling limit of the level set process of **random triangulations** (**discrete maps**).



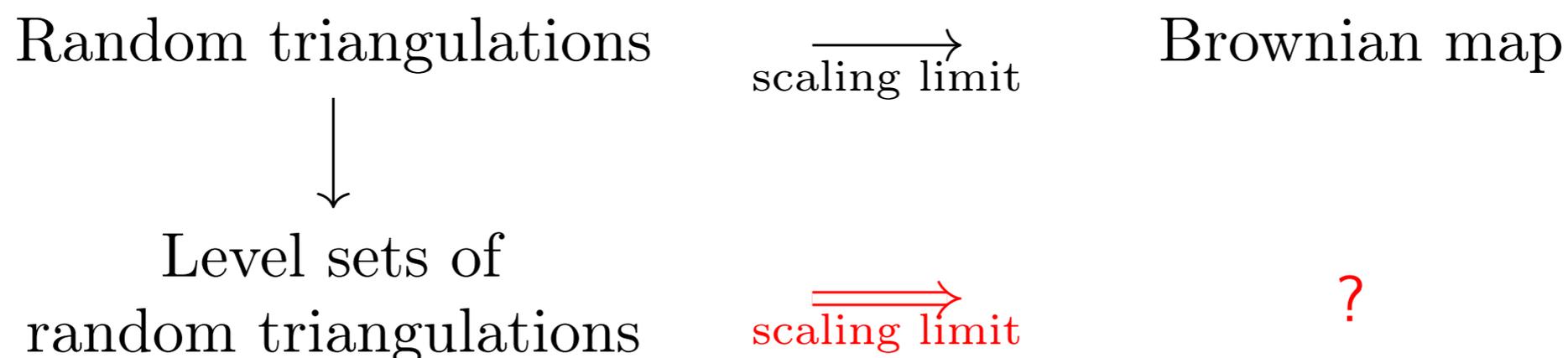
Level sets of the Brownian Map

Imagine the **Brownian map** in such a way that every point at metric distance x from the root is at height x . Now, for every $h > 0$, remove all the points which are not in the ball of radius h centered at the root, and look at the lengths of the cycles as h grows (level set process).

↗ **Questions** (related to the “breadth-first search” of the **Brownian map** of **Miller & Sheffield**):

- What is the law of the level set process of the **Brownian map** as h grows?
- Can one reconstruct the **Brownian map** from the level set processes?

↗ **Our result:** scaling limit of the level set process of **random triangulations** (**discrete maps**).



TRIANGULATIONS



Definitions

A **map** is a finite connected graph properly embedded in the sphere (up to continuous orientation preserving deformations).

Definitions

A **map** is a finite connected graph properly embedded in the sphere (up to continuous orientation preserving deformations).

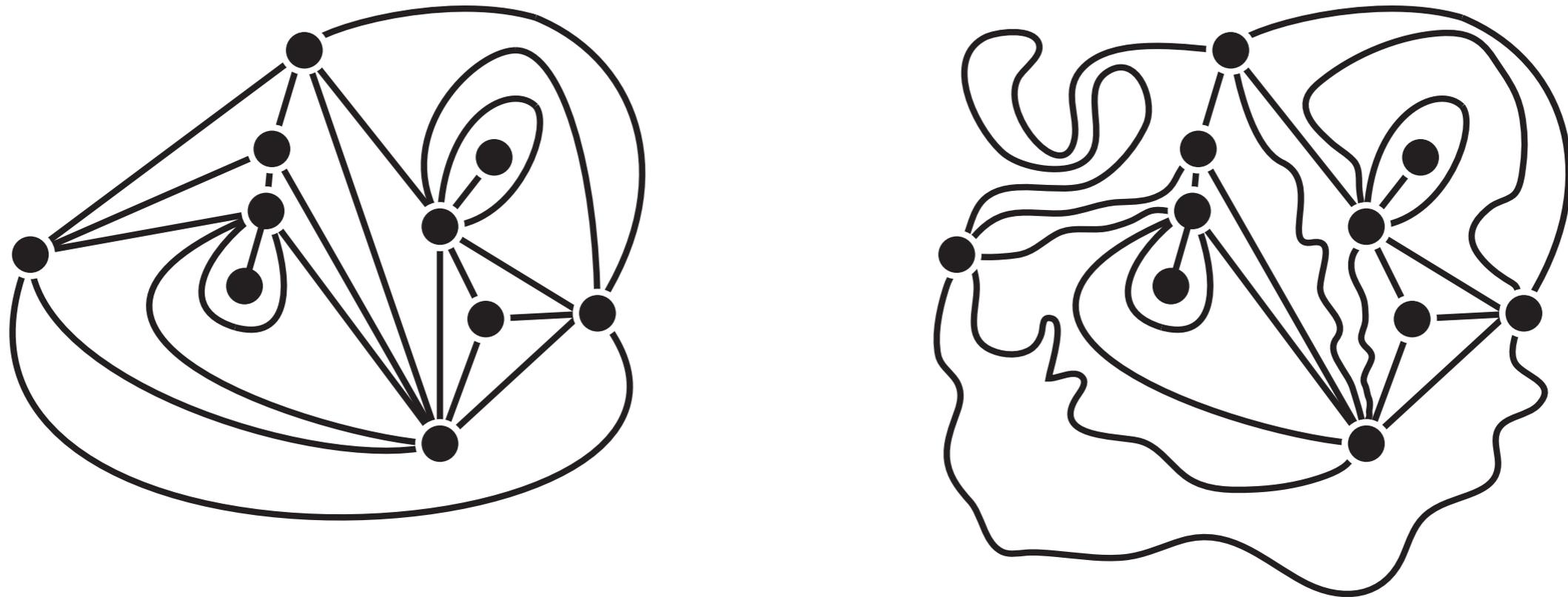


Figure: Two identical maps.

Definitions

A **map** is a finite connected graph properly embedded in the sphere (up to continuous orientation preserving deformations). A map is a **triangulation** when all the faces are triangles.

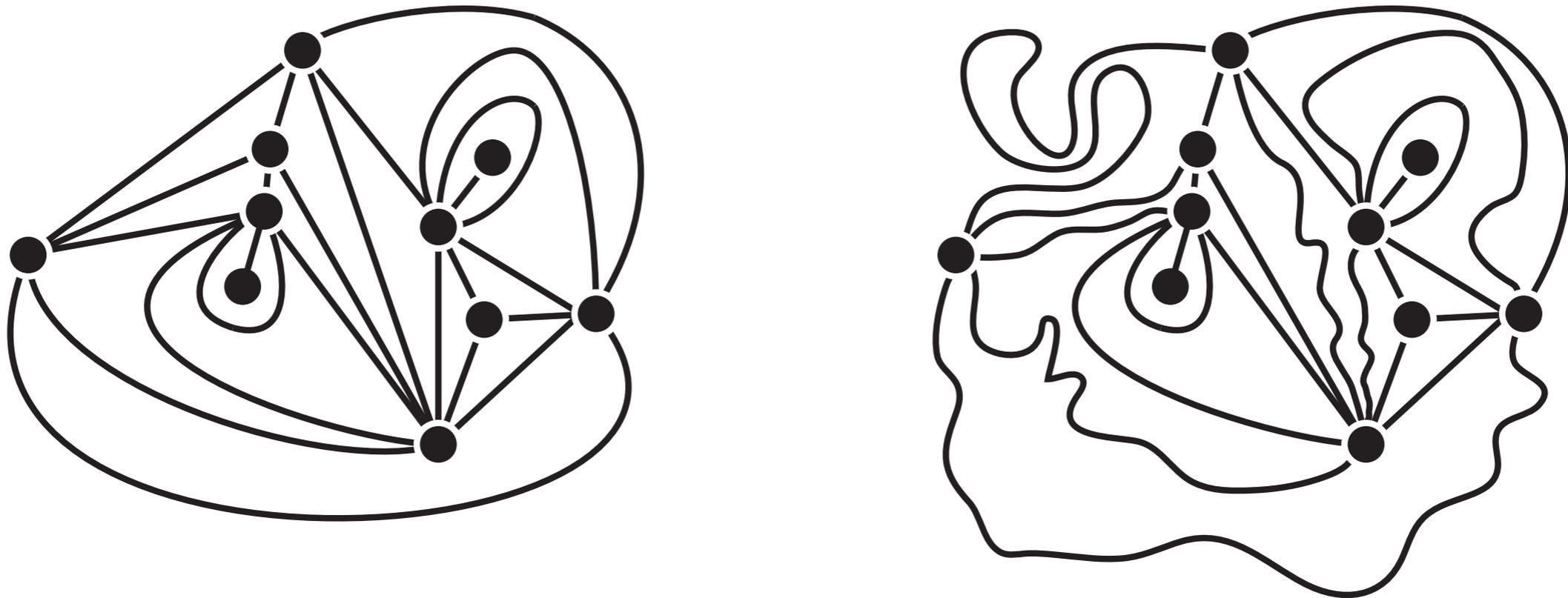


Figure: Two identical triangulations.

Definitions

A **map** is a finite connected graph properly embedded in the sphere (up to continuous orientation preserving deformations). A map is a **triangulation** when all the faces are triangles. A map is **rooted** when an oriented edge is distinguished.

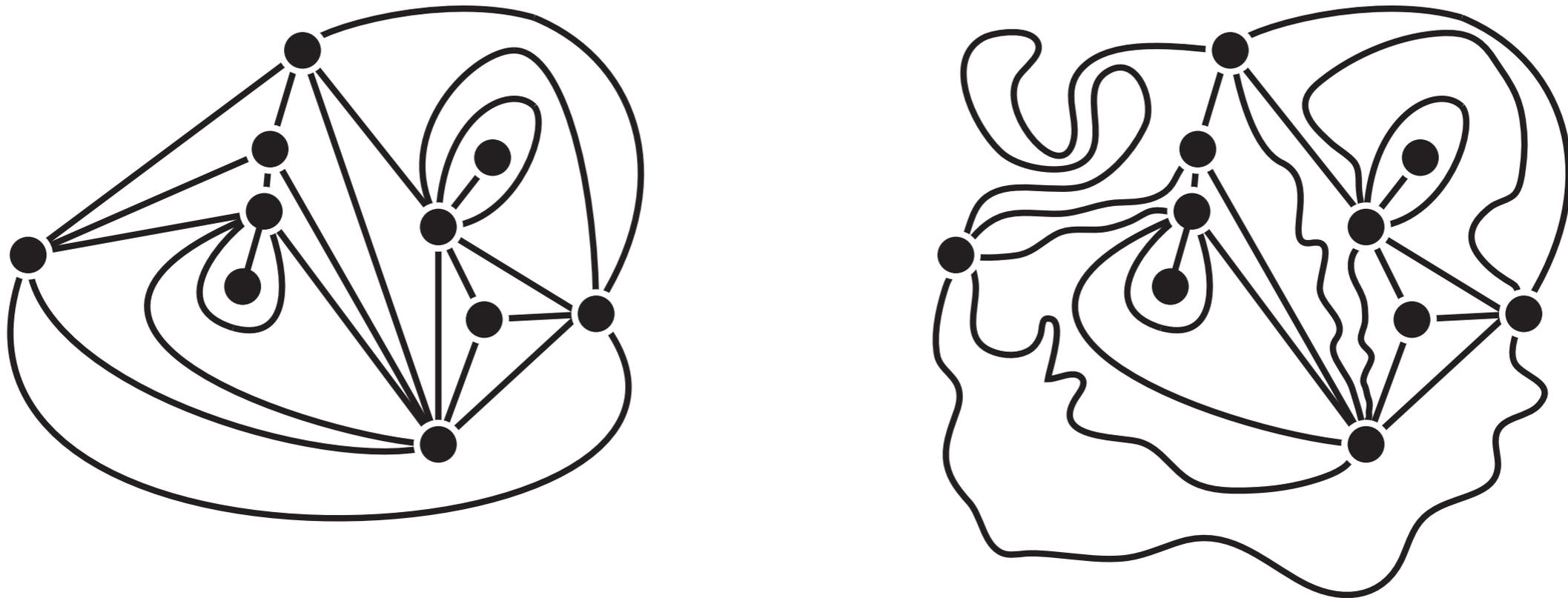


Figure: Two identical triangulations.

Definitions

A **map** is a finite connected graph properly embedded in the sphere (up to continuous orientation preserving deformations). A map is a **triangulation** when all the faces are triangles. A map is **rooted** when an oriented edge is distinguished.

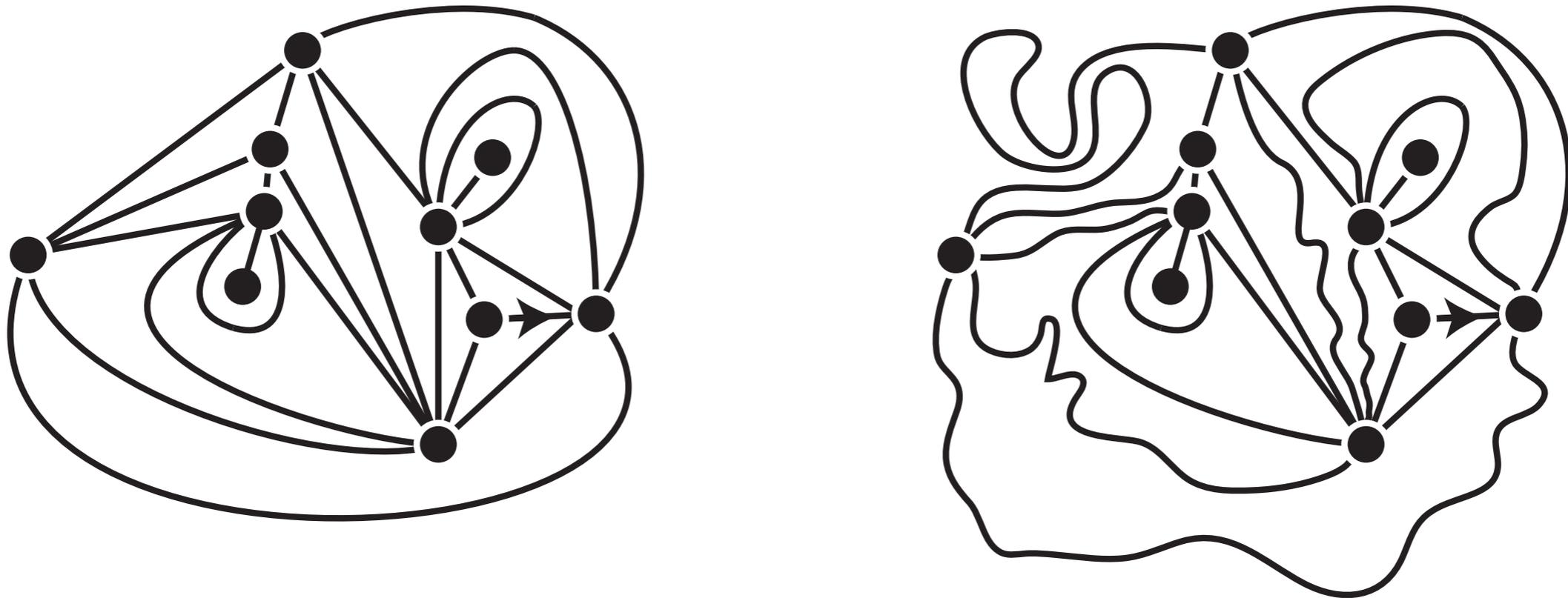


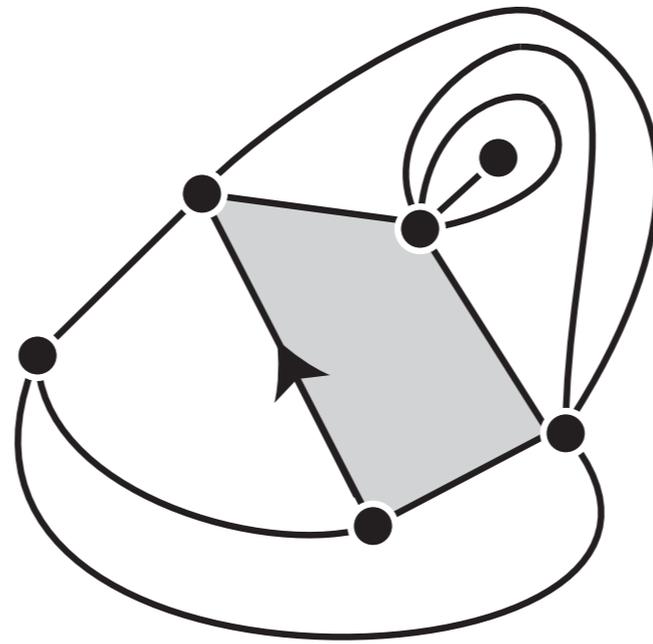
Figure: Two identical rooted triangulations.

TRIANGULATIONS WITH A BOUNDARY



Definitions

A **triangulation with a boundary** is a map where all the faces are triangles, except maybe the one on the right of the root edge which is called the external face.



Definitions

A **triangulation with a boundary** is a map where all the faces are triangles, except maybe the one on the right of the root edge which is called the external face.

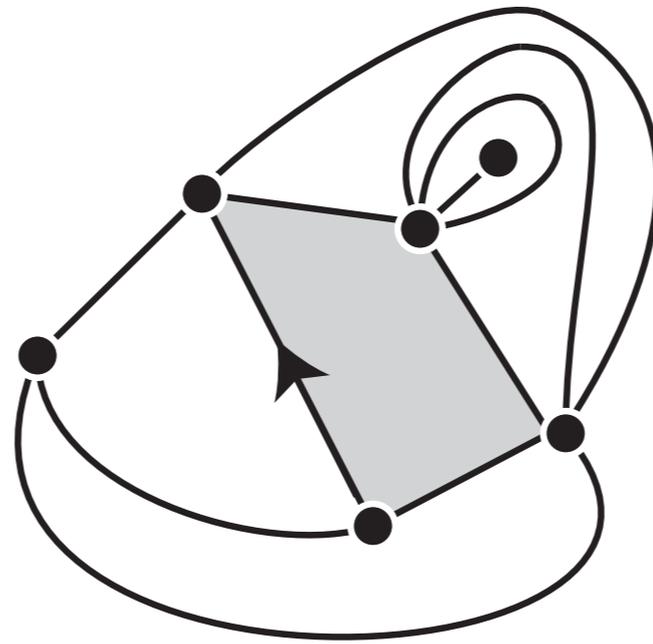


Figure: A triangulation with a boundary with two internal vertices (not adjacent to the external face).

Definitions

A **triangulation with a boundary** is a map where all the faces are triangles, except maybe the one on the right of the root edge which is called the external face.

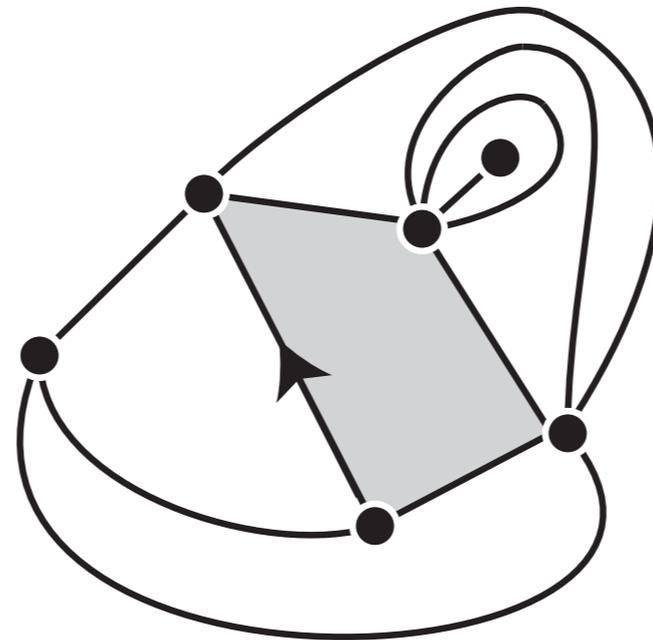


Figure: A triangulation with a boundary with two internal vertices (not adjacent to the external face).

A **triangulation of the p -gon** is a triangulation whose boundary is **simple** and has length p .

Definitions

A **triangulation with a boundary** is a map where all the faces are triangles, except maybe the one on the right of the root edge which is called the external face.

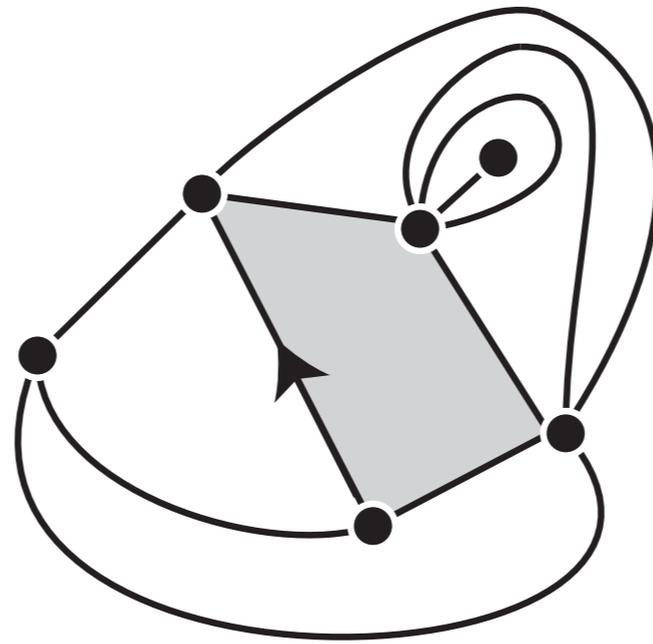


Figure: A triangulation of the 4-gon with two internal vertices (not adjacent to the external face).

A **triangulation of the p -gon** is a triangulation whose boundary is **simple** and has length p .

 what probability measure of planar maps?

Outline

I. PLANAR MAPS

II. BIENAYMÉ–GALTON–WATSON TREES



III. RANDOM MAPS AND GROWTH-FRAGMENTATIONS

Plane trees

We only consider **rooted plane trees**.

Plane trees

We only consider **rooted plane trees**.



Figure: Two different plane trees.

Plane trees

We only consider **rooted plane trees**.



Figure: Two different plane trees.

→ Natural question: what does a large “typical” **plane rooted tree** look like?

Plane trees

We only consider **rooted plane trees**.

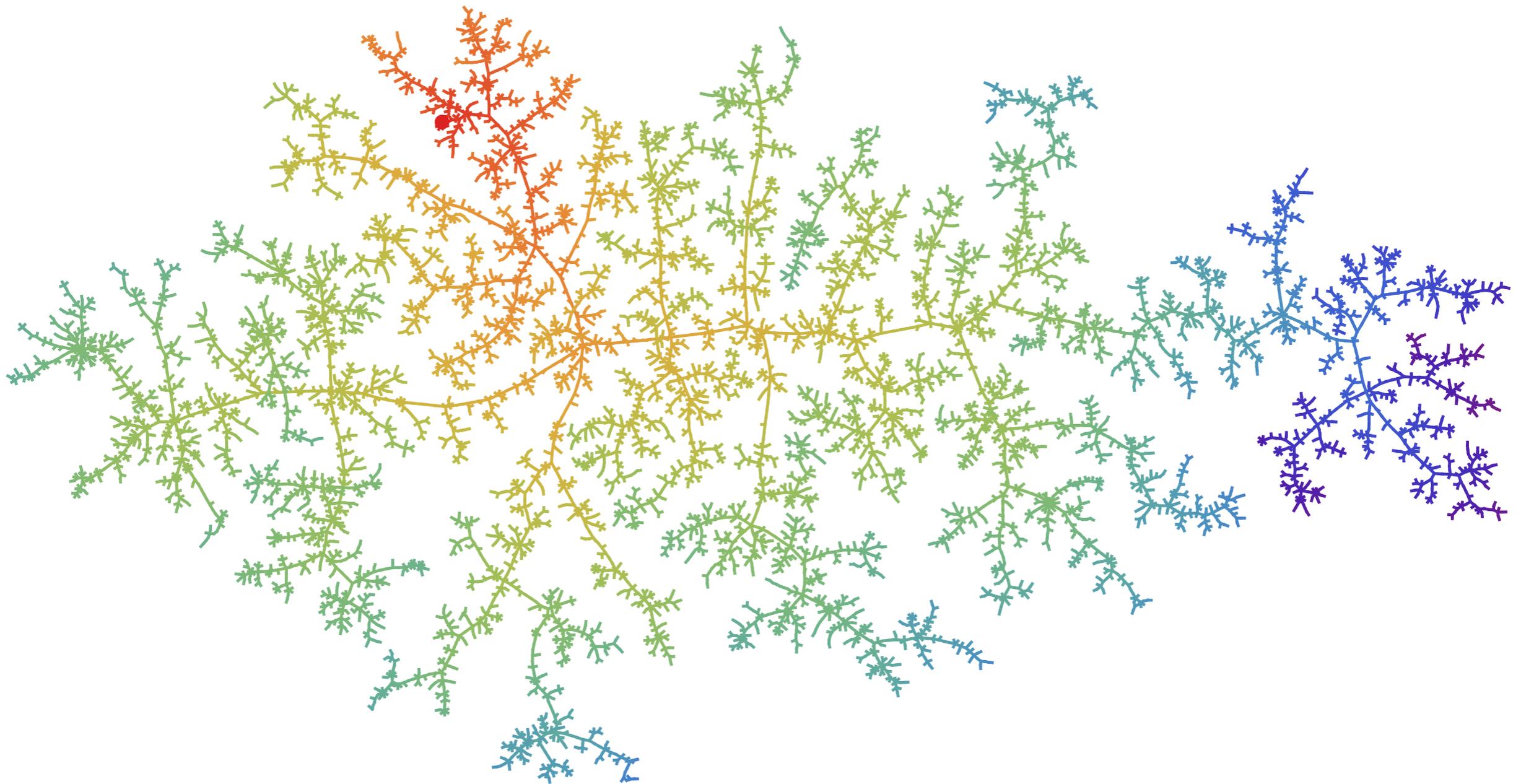


Figure: Two different plane trees.

→ Natural question: what does a large “typical” **plane rooted tree** look like?

→ Let t_n be a large **random plane tree**, chosen uniformly at random among all **rooted plane trees** with n vertices.

A simulation of a large random tree



Uniform plane trees

↗ To study a **uniform plane rooted tree** with n vertices, a key fact is that they can be seen as a **BGW** tree conditioned to have n vertices, with offspring distribution $\mu(i) = \frac{1}{2^{i+1}}$ for $i \geq 0$.

Uniform plane trees

↗ To study a **uniform plane rooted tree** with n vertices, a key fact is that they can be seen as a **BGW** tree conditioned to have n vertices, with offspring distribution $\mu(i) = \frac{1}{2^{i+1}}$ for $i \geq 0$.

Reason: a tree with n vertices then has probability 2^{-2n-1} .

Uniform plane trees

→ To study a **uniform plane rooted tree** with n vertices, a key fact is that they can be seen as a **BGW** tree conditioned to have n vertices, with offspring distribution $\mu(i) = \frac{1}{2^{i+1}}$ for $i \geq 0$.

Reason: a tree with n vertices then has probability 2^{-2n-1} .

→ Where does this geometric distribution come from?

Uniform plane trees

↪ To study a **uniform plane rooted tree** with n vertices, a key fact is that they can be seen as a **BGW** tree conditioned to have n vertices, with offspring distribution $\mu(i) = \frac{1}{2^{i+1}}$ for $i \geq 0$.

Reason: a tree with n vertices then has probability 2^{-2n-1} .

↪ Where does this geometric distribution come from?

One looks for a random tree \mathcal{T} such that for every tree τ

$$\mathbb{P}(\mathcal{T} = \tau) = \frac{\chi^{\text{size of } \tau}}{W(\chi)}$$

Uniform plane trees

↪ To study a **uniform plane rooted tree** with n vertices, a key fact is that they can be seen as a **BGW** tree conditioned to have n vertices, with offspring distribution $\mu(i) = \frac{1}{2^{i+1}}$ for $i \geq 0$.

Reason: a tree with n vertices then has probability 2^{-2n-1} .

↪ Where does this geometric distribution come from?

One looks for a random tree \mathcal{T} such that for every tree τ

$$\mathbb{P}(\mathcal{T} = \tau) = \frac{x^{\text{size of } \tau}}{W(x)}, \quad W(x) = \sum_{n \geq 1} \frac{1}{n} \binom{2n-2}{n-1} x^n = \frac{1 - \sqrt{1-4x}}{2}.$$

Uniform plane trees

↪ To study a **uniform plane rooted tree** with n vertices, a key fact is that they can be seen as a **BGW** tree conditioned to have n vertices, with offspring distribution $\mu(i) = \frac{1}{2^{i+1}}$ for $i \geq 0$.

Reason: a tree with n vertices then has probability 2^{-2n-1} .

↪ Where does this geometric distribution come from?

One looks for a random tree \mathcal{T} such that for every tree τ

$$\mathbb{P}(\mathcal{T} = \tau) = \frac{x^{\text{size of } \tau}}{W(x)}, \quad W(x) = \sum_{n \geq 1} \frac{1}{n} \binom{2n-2}{n-1} x^n = \frac{1 - \sqrt{1-4x}}{2}.$$

The radius of convergence is $1/4$, and by taking $x = 1/4$, one gets a **BGW** tree with offspring distribution μ .

Simply generated trees

In particular, **uniform plane trees** are particular cases of so-called simply generated (or Boltzmann) trees:

Simply generated trees

In particular, **uniform plane trees** are particular cases of so-called simply generated (or Boltzmann) trees:

Given a sequence $w = (w(i); i \geq 0)$ of nonnegative real numbers, with every $\tau \in \mathbb{T}$, associate a weight $\Omega^w(\tau)$:

$$\Omega^w(\tau) = \prod_{u \in \tau} w(\text{number of children of } u).$$

Simply generated trees

In particular, **uniform plane trees** are particular cases of so-called simply generated (or Boltzmann) trees:

Given a sequence $w = (w(i); i \geq 0)$ of nonnegative real numbers, with every $\tau \in \mathbb{T}$, associate a weight $\Omega^w(\tau)$:

$$\Omega^w(\tau) = \prod_{u \in \tau} w(\text{number of children of } u).$$

Then, if \mathbb{T}_n is the set of all trees with n vertices, for every $\tau \in \mathbb{T}_n$, set

$$\mathbb{P}_n^w(\tau) = \frac{\Omega^w(\tau)}{\sum_{T \in \mathbb{T}_n} \Omega^w(T)}.$$

SCALING LIMITS OF LARGE SIMPLY GENERATED TREES



Large simply generated trees

→ If the weight sequence is sufficiently regular, the scaling limit of **simply generated trees** is the **Brownian tree** (Aldous).

Large simply generated trees

↗ If the weight sequence is sufficiently regular, the scaling limit of **simply generated trees** is the **Brownian tree** (Aldous).

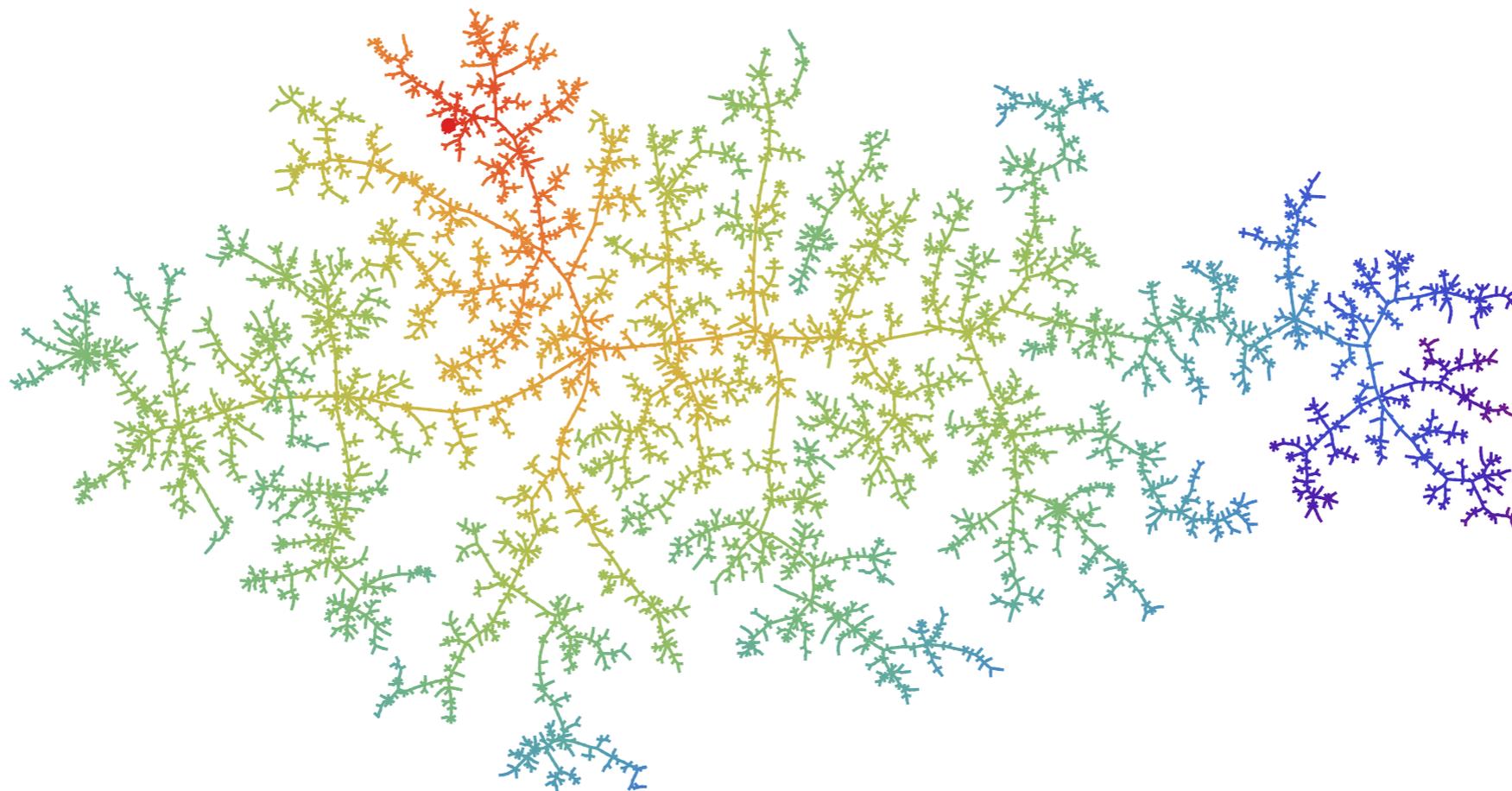


Figure: A non isometric embedding of a realization of the Brownian tree.

Large simply generated trees

- If the weight sequence is sufficiently regular, the scaling limit of **simply generated trees** is the **Brownian tree** (Aldous).
- If the weight sequence has a heavy tail behavior, the scaling limit of **simply generated trees** is a **stable tree** (Duquesne, Le Gall, Le Jan).

Large simply generated trees

- ↗ If the weight sequence is sufficiently regular, the scaling limit of **simply generated trees** is the **Brownian tree** (Aldous).
- ↗ If the weight sequence has a heavy tail behavior, the scaling limit of **simply generated trees** is a **stable tree** (Duquesne, Le Gall, Le Jan).

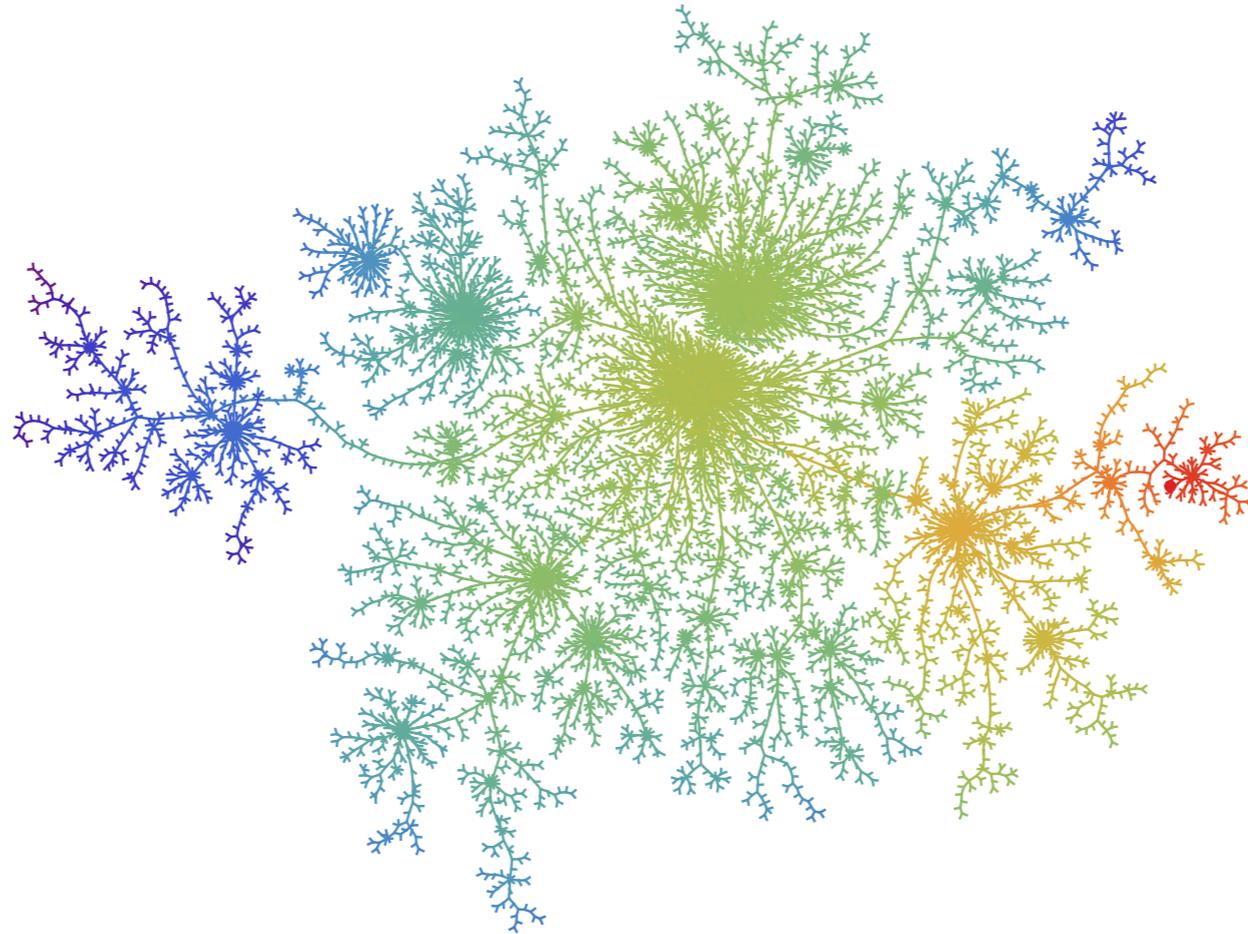


Figure: A non isometric embedding of a realization of a **stable tree** with index 1.2.

Outline

I. PLANAR MAPS

II. BIENAYMÉ–GALTON–WATSON TREES

III. SCALING LIMITS OF LEVEL SETS OF RANDOM MAPS



Random maps

 What probability distribution on plane triangulations?

Random maps

↗ What probability distribution on plane triangulations?

For **BGW** trees: how to force a **BGW tree** to be large?

Random maps

↗ What probability distribution on plane triangulations?

For **BGW** trees: how to force a **BGW tree** to be large? One way is to condition it to have size p .

Random maps

↗ What probability distribution on plane triangulations?

For **BGW** trees: how to force a **BGW tree** to be large? One way is to condition it to have size p . Another way is to consider a forest of p **BGW trees**.

Random maps

↗ What probability distribution on plane triangulations?

For **BGW** trees: how to force a **BGW tree** to be large? One way is to condition it to have size p . Another way is to consider a forest of p **BGW trees**.

↗ Similarly, for **planar triangulations** we will take a Boltzmann distribution on **planar triangulations** with a large boundary p .

Definitions

Let $\mathcal{T}_{n,p}$ denote the set of all **triangulations of the p -gon** with n internal vertices.

Definitions

Let $\mathcal{T}_{n,p}$ denote the set of all **triangulations of the p -gon** with n internal vertices. We have (**Krikun**)

$$\#\mathcal{T}_{n,p} = 4^{n-1} \frac{p (2p)! (2p + 3n - 5)!!}{(p!)^2 n! (2p + n - 1)!!}$$

Definitions

Let $\mathcal{T}_{n,p}$ denote the set of all **triangulations of the p -gon** with n internal vertices. We have (**Krikun**)

$$\#\mathcal{T}_{n,p} = 4^{n-1} \frac{p (2p)! (2p + 3n - 5)!!}{(p!)^2 n! (2p + n - 1)!!} \underset{n \rightarrow \infty}{\sim} C(p) (12\sqrt{3})^n n^{-5/2}.$$

Definitions

Let $\mathcal{T}_{n,p}$ denote the set of all **triangulations of the p -gon** with n internal vertices. We have (**Krikun**)

$$\#\mathcal{T}_{n,p} = 4^{n-1} \frac{p (2p)! (2p + 3n - 5)!!}{(p!)^2 n! (2p + n - 1)!!} \underset{n \rightarrow \infty}{\sim} C(p) (12\sqrt{3})^n n^{-5/2}.$$

Therefore, the radius of convergence of $\sum_{n \geq 0} \#\mathcal{T}_{n,p} z^n$ is $(12\sqrt{3})^{-1}$.

Definitions

Let $\mathcal{T}_{n,p}$ denote the set of all **triangulations of the p -gon** with n internal vertices. We have (**Krikun**)

$$\#\mathcal{T}_{n,p} = 4^{n-1} \frac{p (2p)! (2p + 3n - 5)!!}{(p!)^2 n! (2p + n - 1)!!} \underset{n \rightarrow \infty}{\sim} C(p) (12\sqrt{3})^n n^{-5/2}.$$

Therefore, the radius of convergence of $\sum_{n \geq 0} \#\mathcal{T}_{n,p} z^n$ is $(12\sqrt{3})^{-1}$.

Set

$$Z(p) = \sum_{n=0}^{\infty} \left(\frac{1}{12\sqrt{3}} \right)^n \#\mathcal{T}_{n,p} < \infty.$$

Definitions

Let $\mathcal{T}_{n,p}$ denote the set of all **triangulations of the p -gon** with n internal vertices. We have (Krikun)

$$\#\mathcal{T}_{n,p} = 4^{n-1} \frac{p (2p)! (2p + 3n - 5)!!}{(p!)^2 n! (2p + n - 1)!!} \underset{n \rightarrow \infty}{\sim} C(p) (12\sqrt{3})^n n^{-5/2}.$$

Therefore, the radius of convergence of $\sum_{n \geq 0} \#\mathcal{T}_{n,p} z^n$ is $(12\sqrt{3})^{-1}$.
Set

$$Z(p) = \sum_{n=0}^{\infty} \left(\frac{1}{12\sqrt{3}} \right)^n \#\mathcal{T}_{n,p} < \infty.$$

A **triangulation of the p -gon** chosen at random with probability

$$(12\sqrt{3})^{-\#(\text{internal vertices})} Z(p)^{-1}$$

is called a **Boltzmann triangulation of the p -gon**.

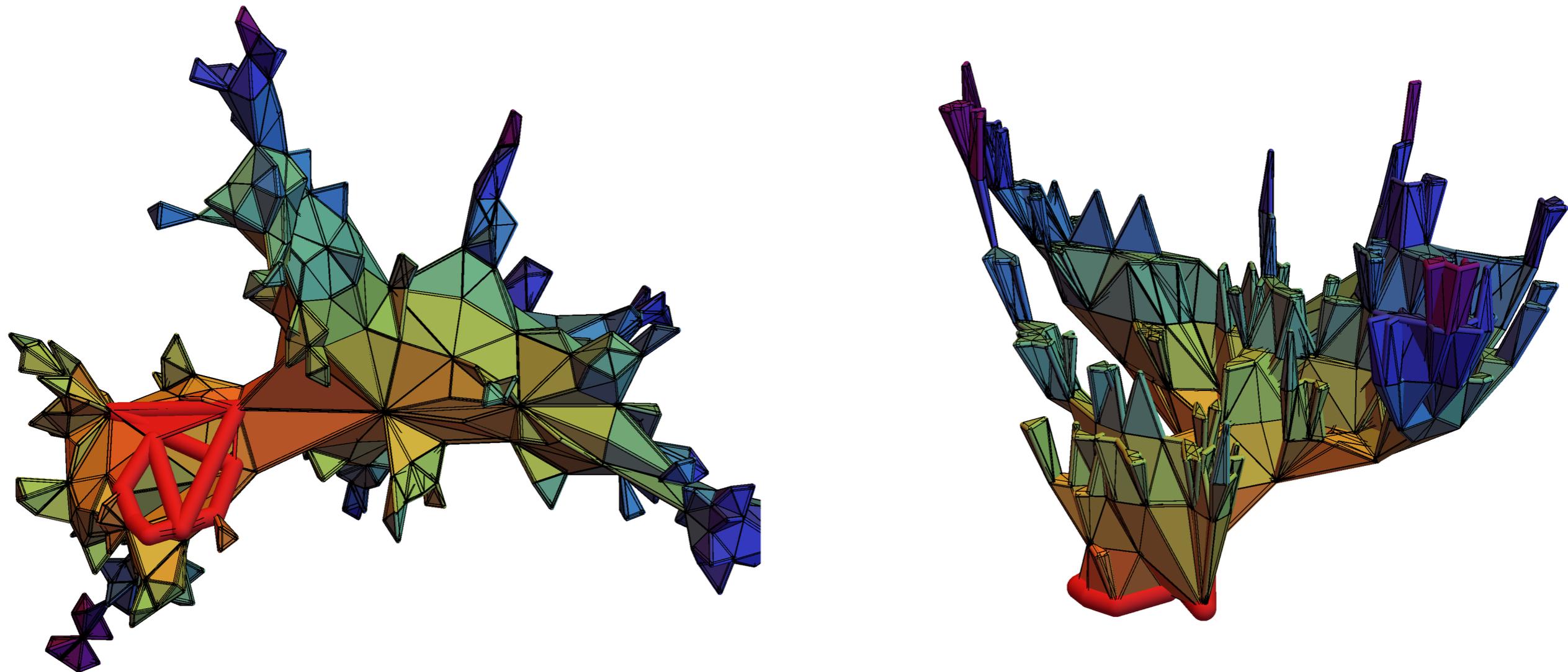


Figure: A Boltzmann triangulation of the 9-gon.

LEVEL SETS OF BOLTZMANN TRIANGULATIONS WITH A BOUNDARY



Large Boltzmann triangulations with a boundary

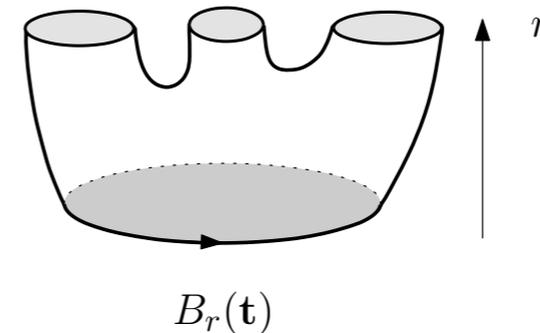
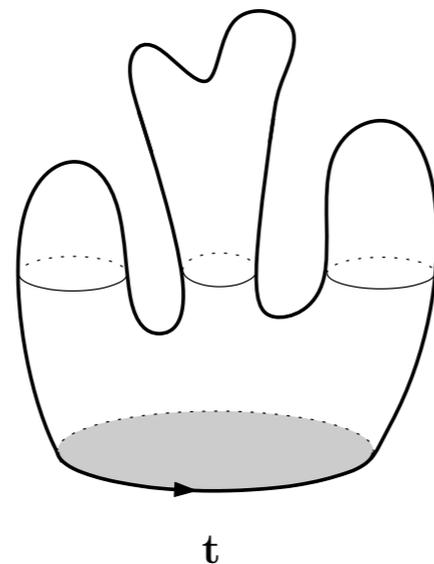
Let $T^{(p)}$ be a random Boltzmann triangulation of the p -gon

Large Boltzmann triangulations with a boundary

Let $\mathcal{T}^{(p)}$ be a random **Boltzmann triangulation** of the p -gon, let $B_r(\mathcal{T}^{(p)})$ be the map made of the vertices with distance at most r from the boundary

Large Boltzmann triangulations with a boundary

Let $T^{(p)}$ be a random Boltzmann triangulation of the p -gon, let $B_r(T^{(p)})$ be the map made of the vertices with distance at most r from the boundary

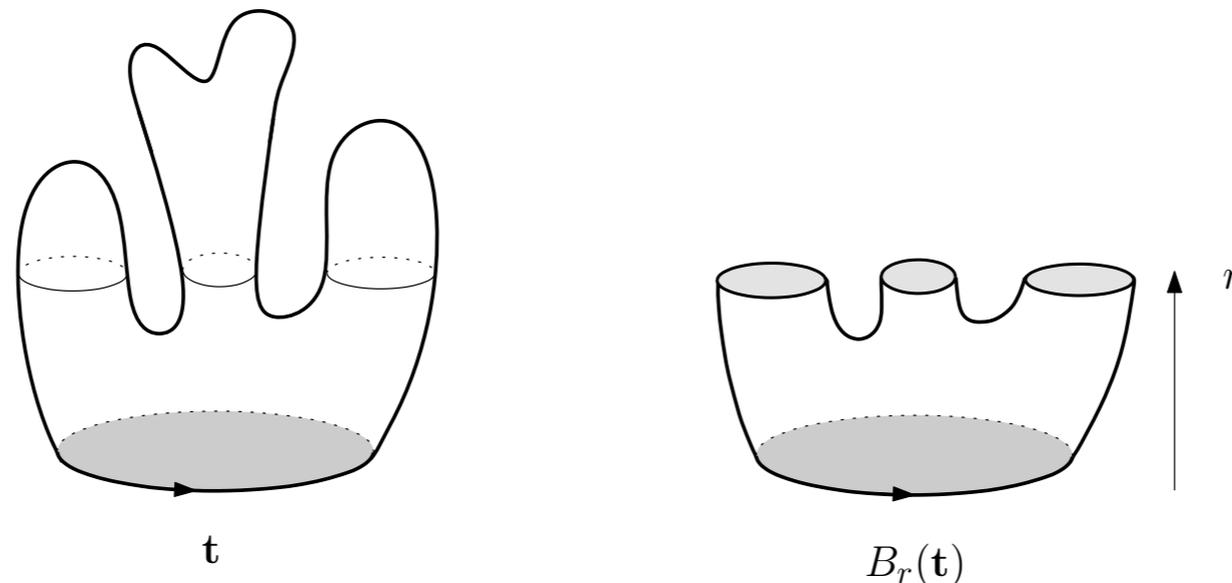


Large Boltzmann triangulations with a boundary

Let $\mathcal{T}^{(p)}$ be a random **Boltzmann triangulation** of the p -gon, let $B_r(\mathcal{T}^{(p)})$ be the map made of the vertices with distance at most r from the boundary, and

$$\mathbb{L}^{(p)}(r) := \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right).$$

be lengths (or perimeters) of the cycles of $B_r(\mathcal{T}^{(p)})$, ranked in decreasing order.

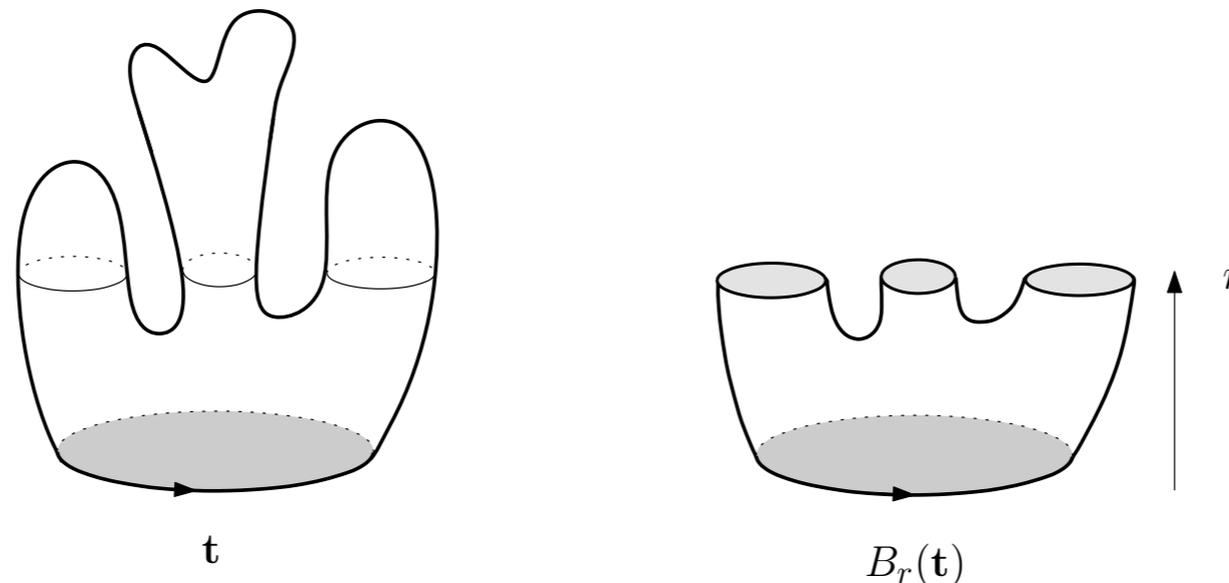


Large Boltzmann triangulations with a boundary

Let $\mathcal{T}^{(p)}$ be a random **Boltzmann triangulation** of the p -gon, let $B_r(\mathcal{T}^{(p)})$ be the map made of the vertices with distance at most r from the boundary, and

$$\mathbb{L}^{(p)}(r) := \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right).$$

be lengths (or perimeters) of the cycles of $B_r(\mathcal{T}^{(p)})$, ranked in decreasing order.



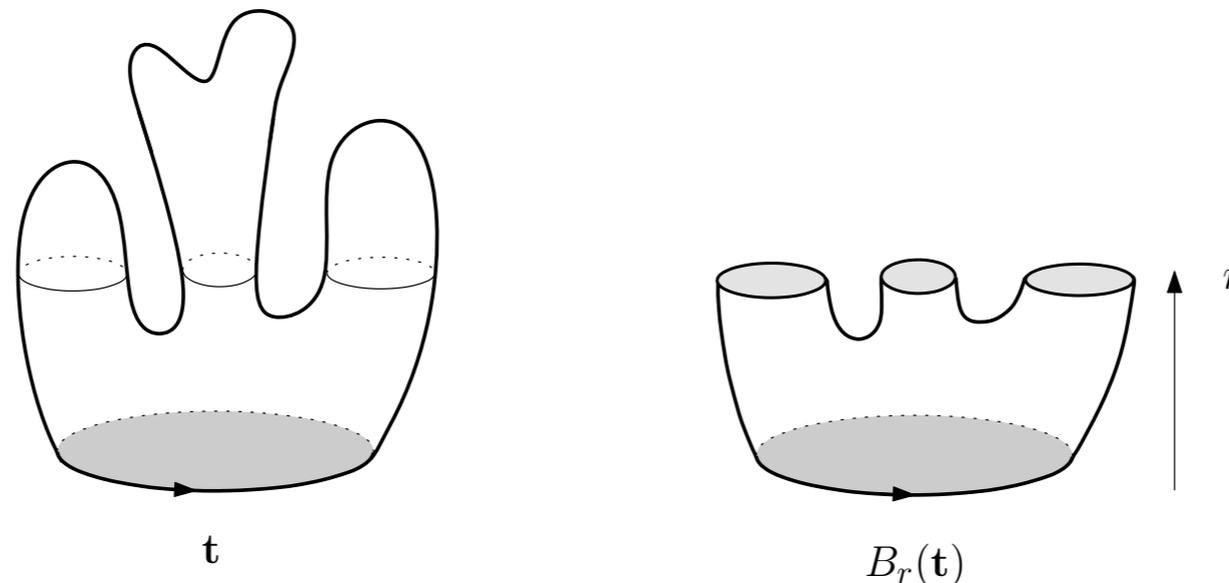
↪ **Goal:** obtain a functional invariance principle of $(\mathbb{L}^{(p)}(r); r \geq 0)$.

Large Boltzmann triangulations with a boundary

Let $\mathcal{T}^{(p)}$ be a random **Boltzmann triangulation** of the p -gon, let $B_r(\mathcal{T}^{(p)})$ be the map made of the vertices with distance at most r from the boundary, and

$$\mathbb{L}^{(p)}(r) := \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right).$$

be lengths (or perimeters) of the cycles of $B_r(\mathcal{T}^{(p)})$, ranked in decreasing order.



→ **Goal:** obtain a functional invariance principle of $(\mathbb{L}^{(p)}(r); r \geq 0)$. In this spirit, a “breadth-first search” of the **Brownian map** is given by [Miller & Sheffield](#).

Simulation



The theorem

Recall that $\mathbb{L}^{(p)}(r) = \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right)$ are the lengths of the cycles of $B_r(\mathbb{T}^{(p)})$ ranked in decreasing order.

The theorem

Recall that $\mathbb{L}^{(p)}(r) = \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right)$ are the lengths of the cycles of $B_r(\mathbb{T}^{(p)})$ ranked in decreasing order.

Theorem (Bertoin, Curien, K. '15).

We have

$$\left(\frac{1}{p} \cdot \mathbb{L}^{(p)}(t\sqrt{p}); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(\mathbb{X} \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

The theorem

Recall that $\mathbb{L}^{(p)}(r) = \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right)$ are the lengths of the cycles of $B_r(\mathbb{T}^{(p)})$ ranked in decreasing order.

Theorem (Bertoin, Curien, K. '15).

We have

$$\left(\frac{1}{p} \cdot \mathbb{L}^{(p)}(t\sqrt{p}); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(\mathbb{X} \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

in distribution in ℓ_3^\downarrow , where $\mathbb{X} = (\mathbb{X}(t); t \geq 0)$ is a càdlàg process with values in ℓ_3^\downarrow

The theorem

Recall that $\mathbb{L}^{(p)}(r) = \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right)$ are the lengths of the cycles of $B_r(\mathbb{T}^{(p)})$ ranked in decreasing order.

Theorem (Bertoin, Curien, K. '15).

We have

$$\left(\frac{1}{p} \cdot \mathbb{L}^{(p)}(t\sqrt{p}); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(\mathbb{X} \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

in distribution in ℓ_3^\downarrow , where $\mathbb{X} = (\mathbb{X}(t); t \geq 0)$ is a càdlàg process with values in ℓ_3^\downarrow , which is a *self-similar growth-fragmentation process* (Bertoin '15).

THE MAIN TOOL: A PEELING EXPLORATION



Geometry of random maps

Several techniques to study **random maps**:

Geometry of random maps

Several techniques to study **random maps**:

– **bijective techniques**,

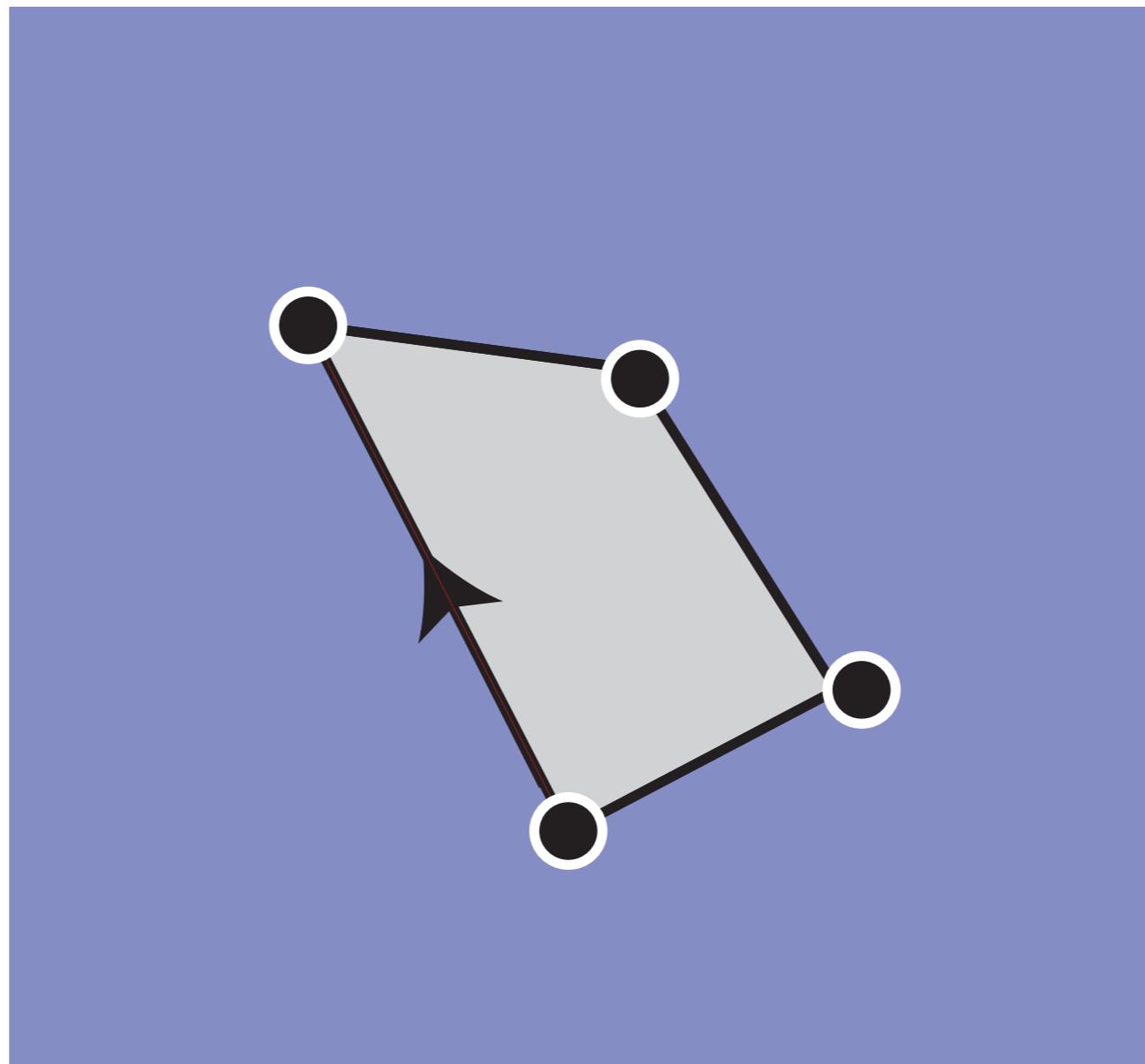
Geometry of random maps

Several techniques to study **random maps**:

- **bijection techniques**,
- **peeling**, which is a Markovian way to iteratively explore a random map ([Watabiki '95](#), [Angel '03](#), [Budd '14](#)).

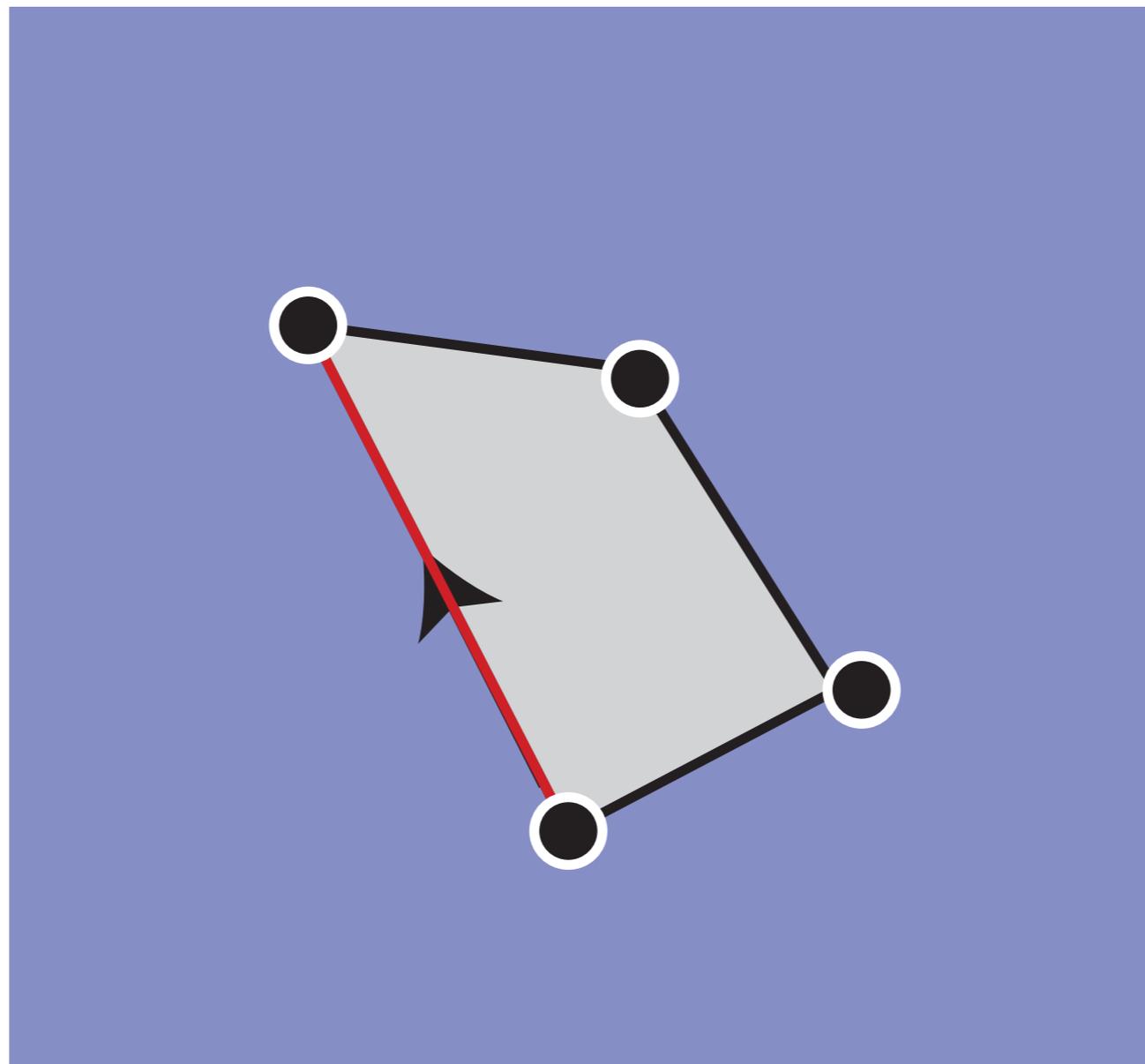
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



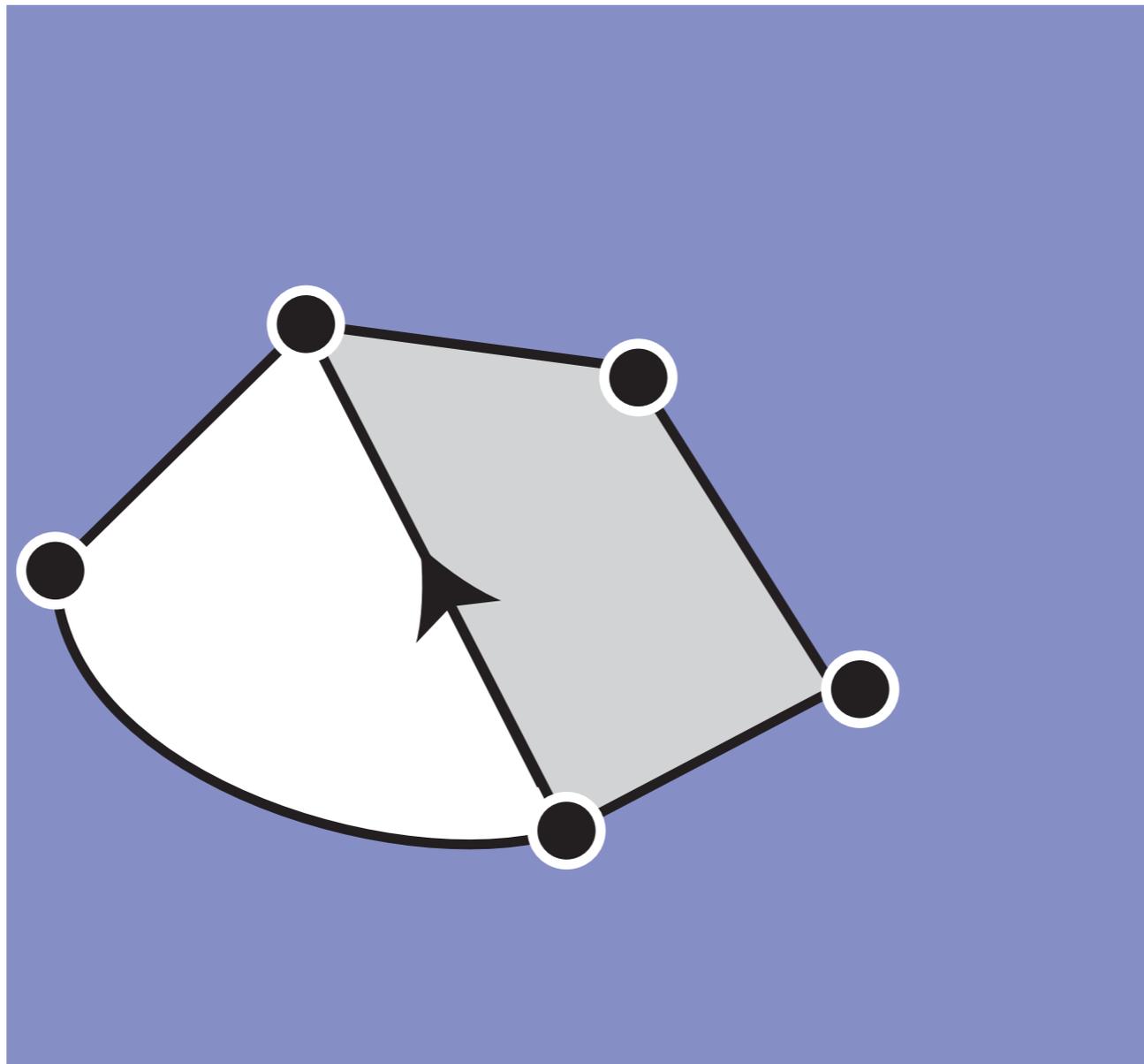
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



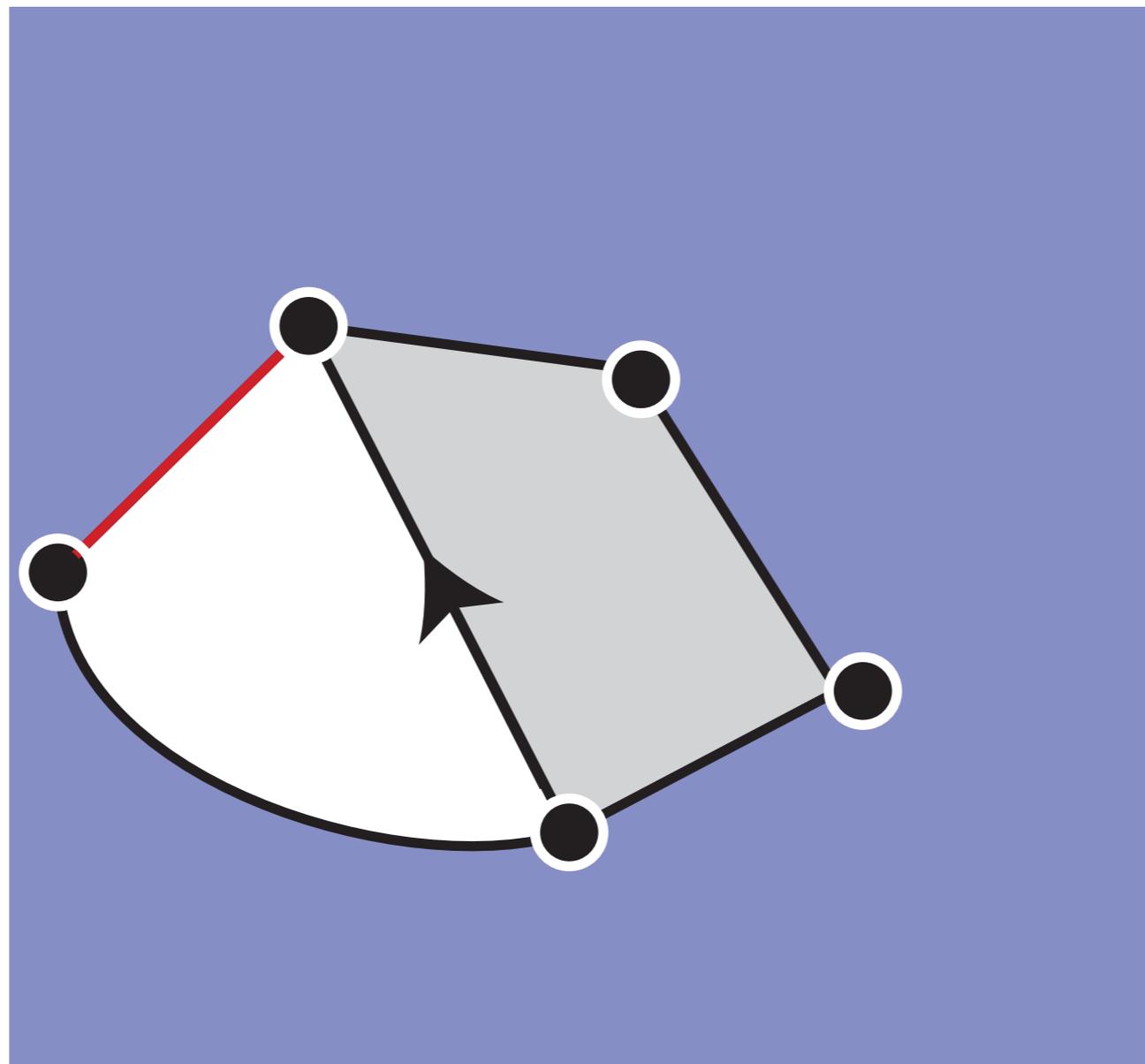
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



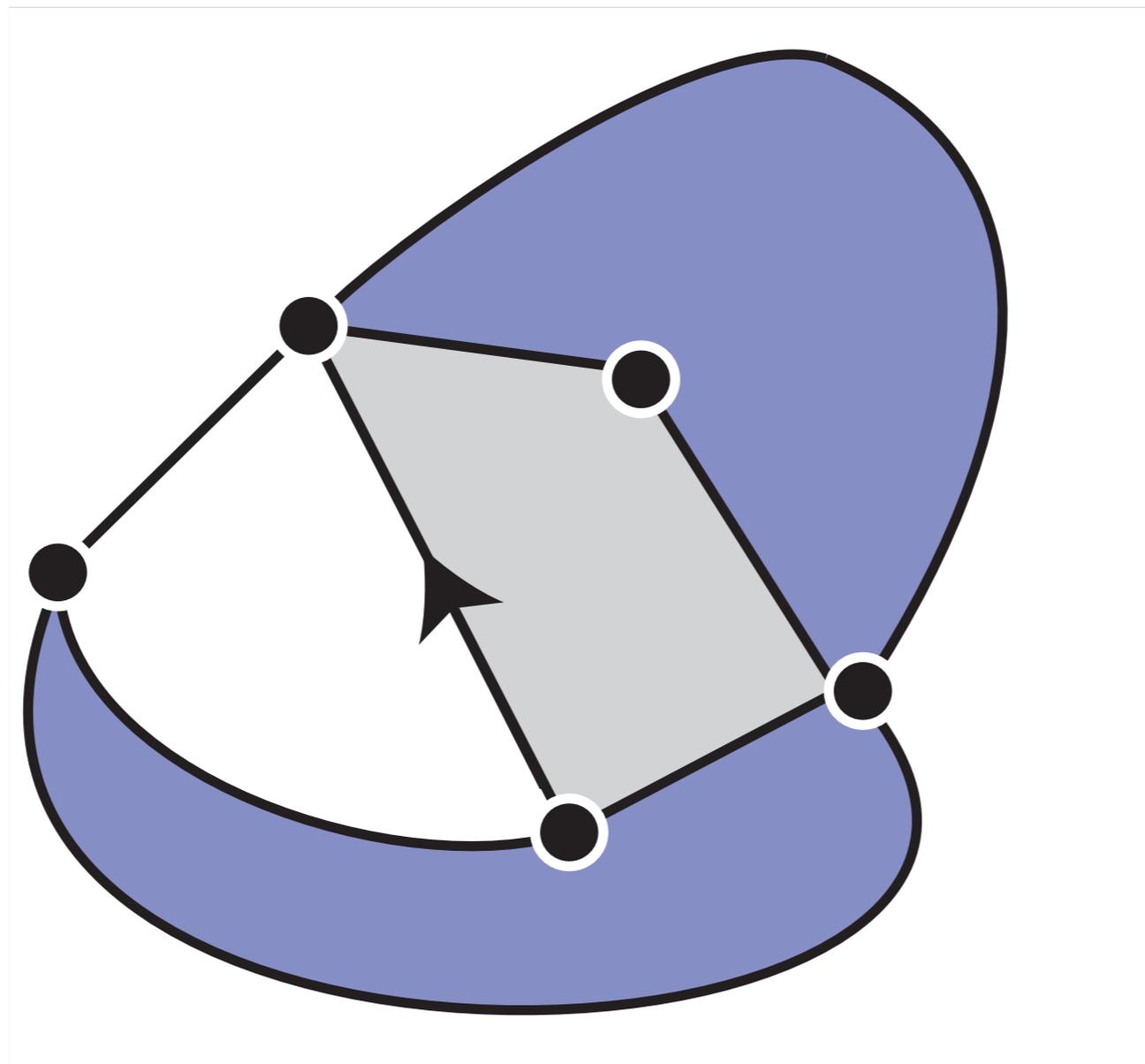
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



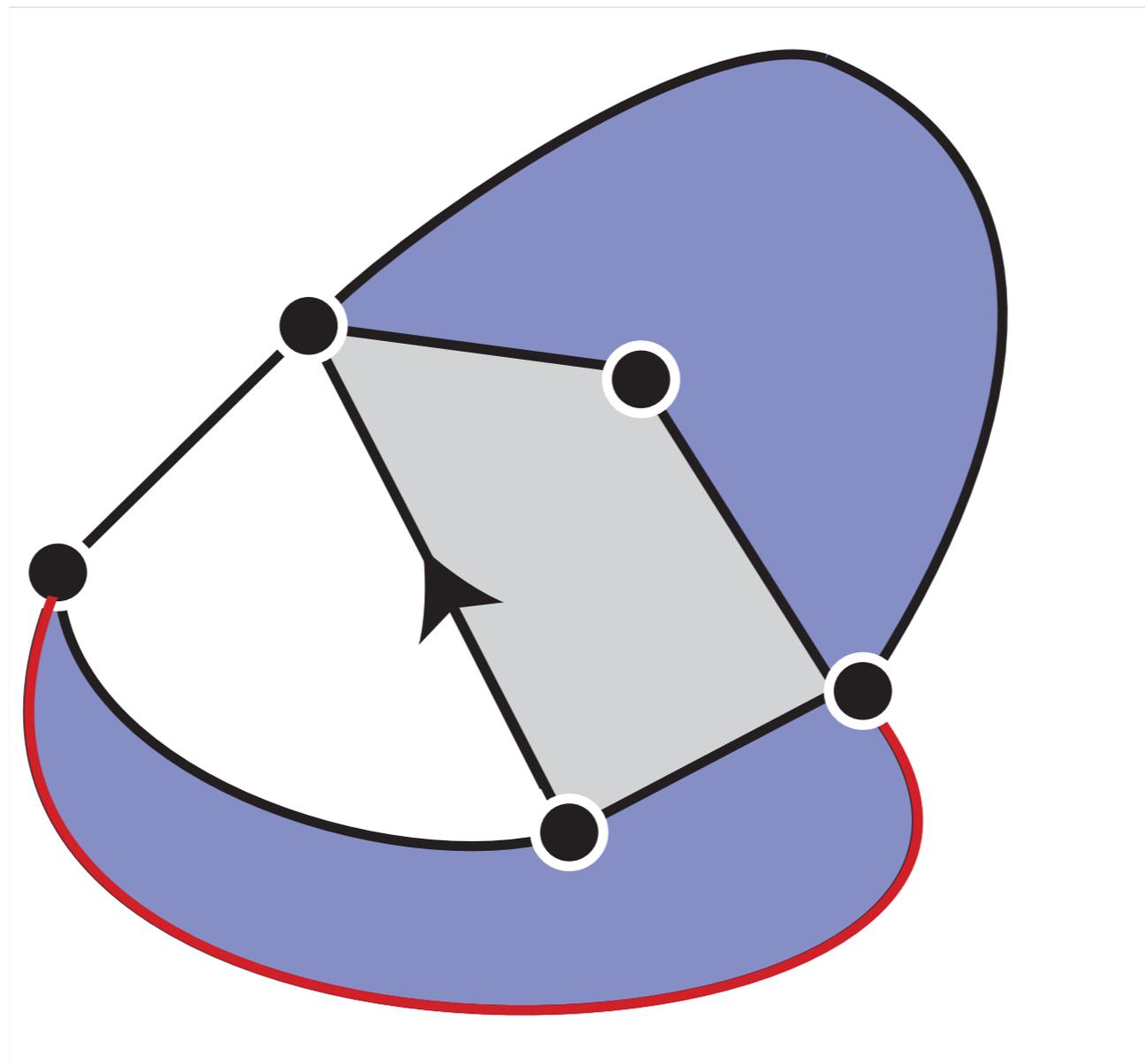
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



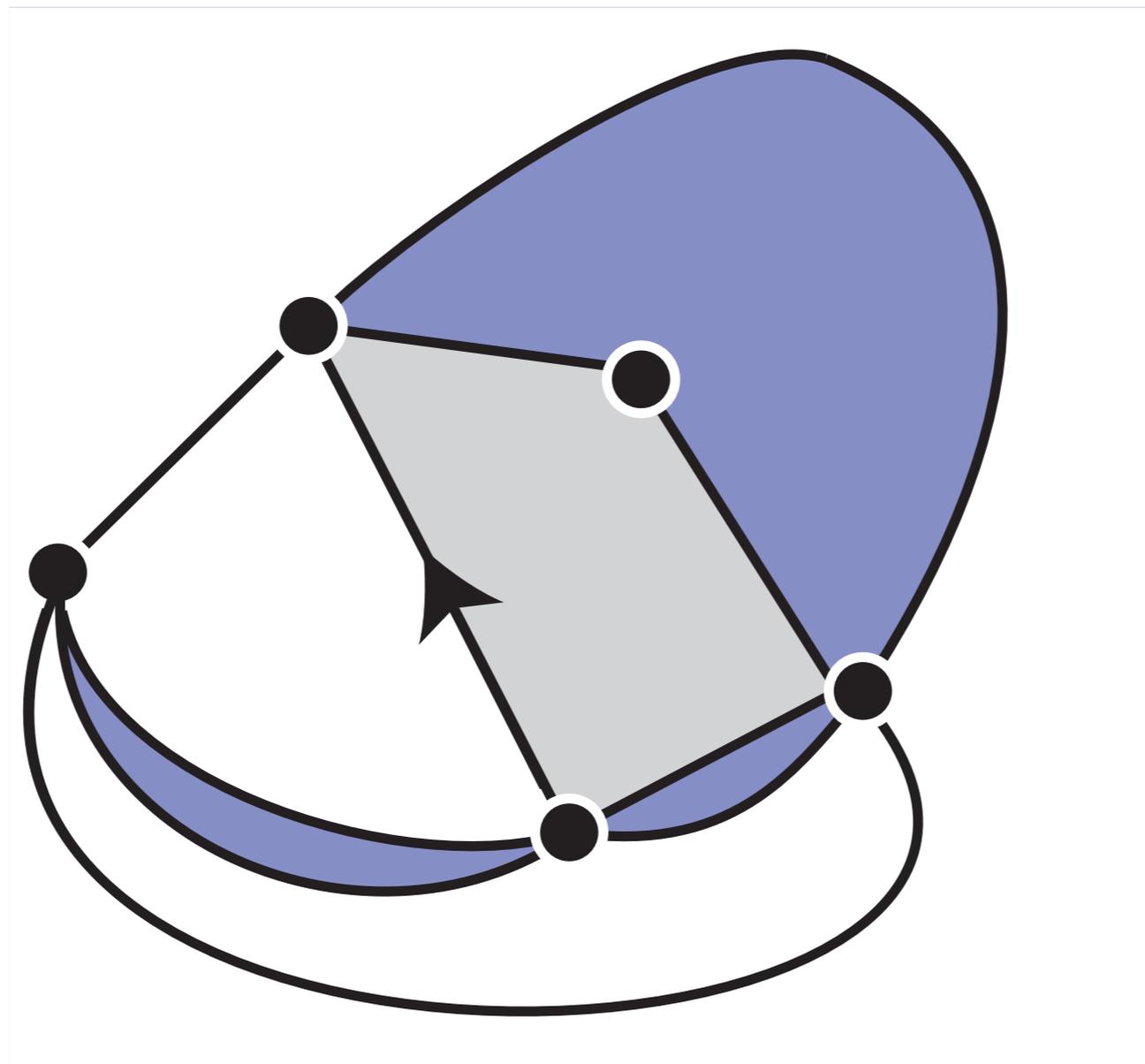
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



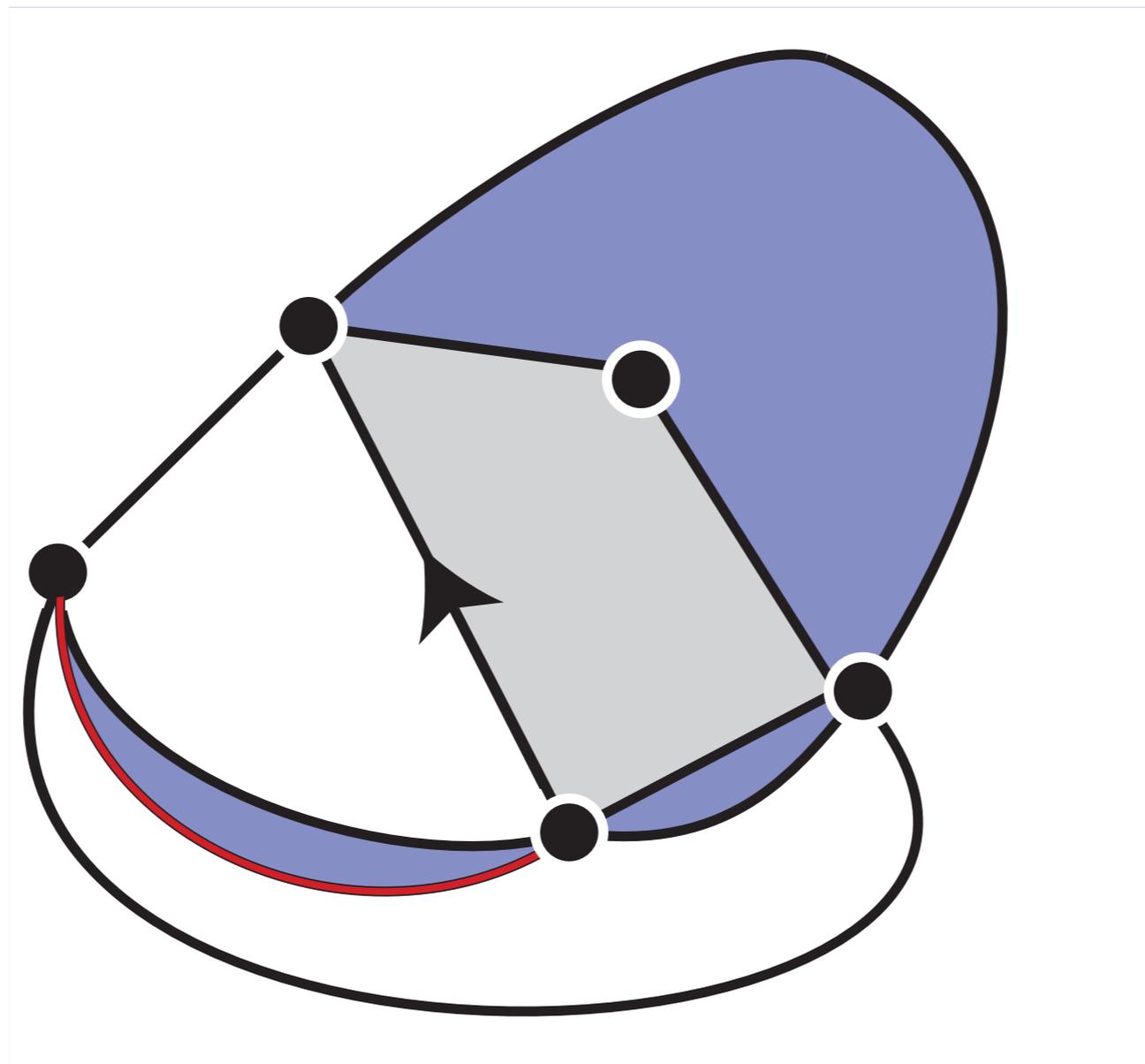
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



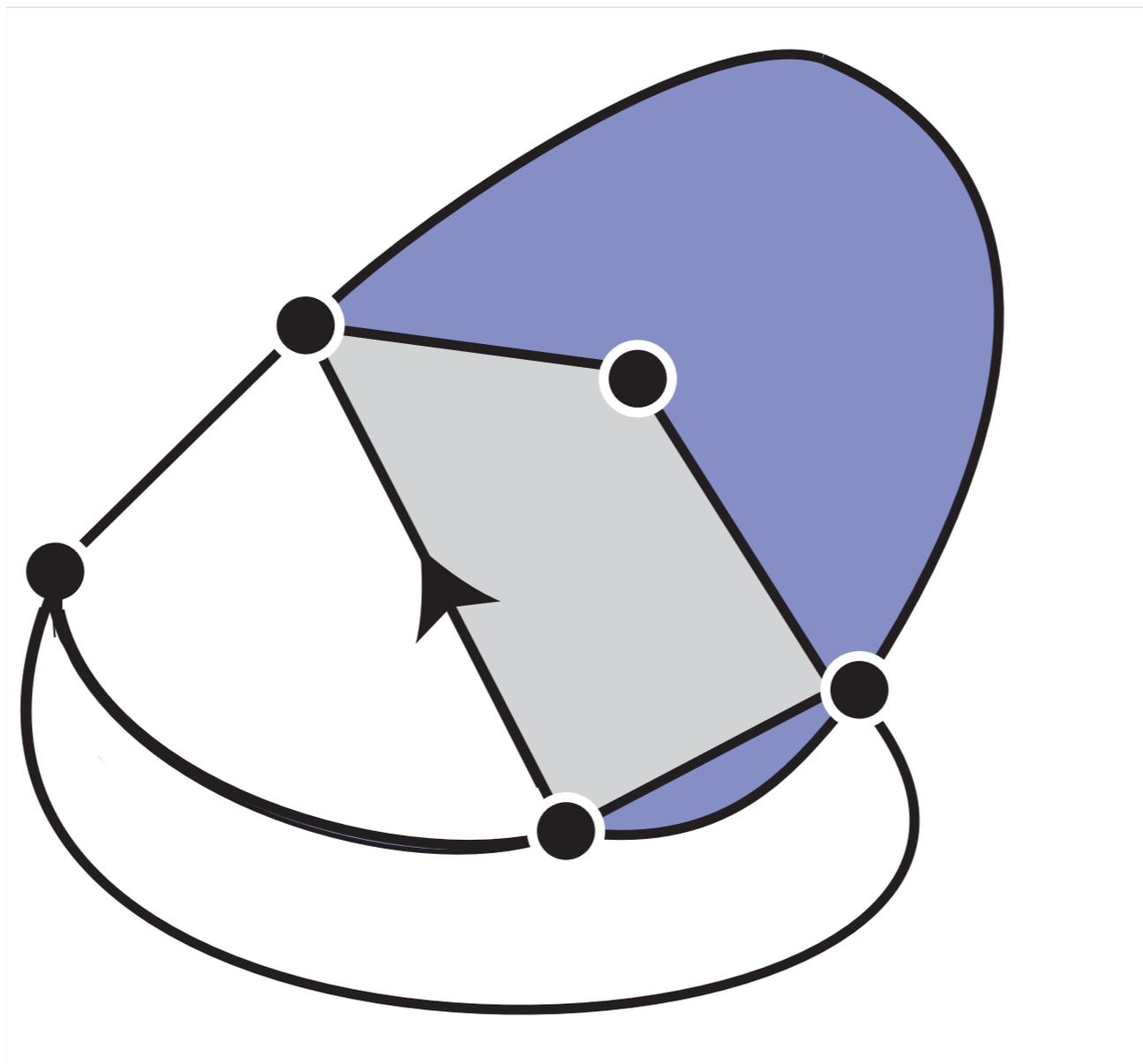
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



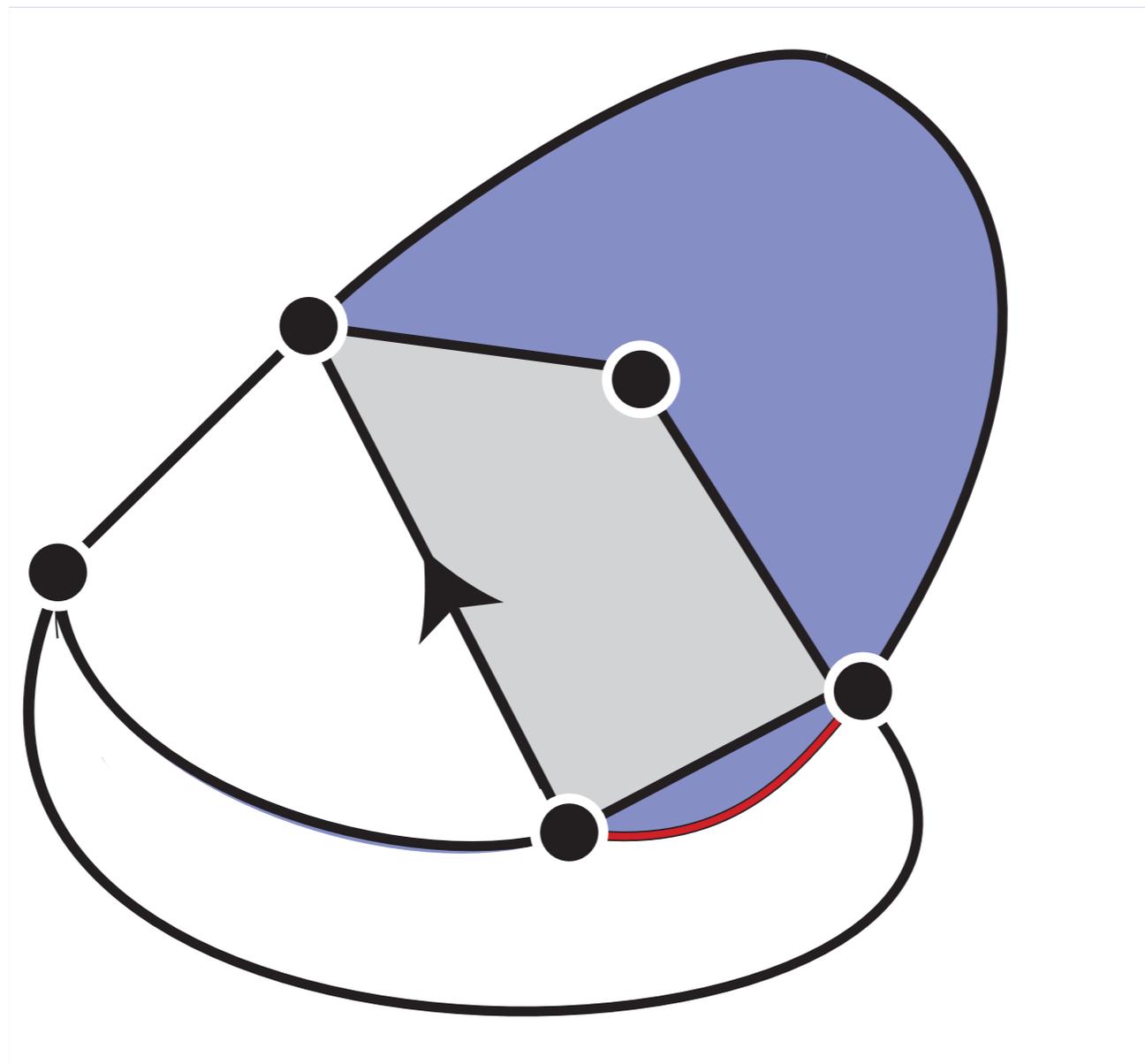
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



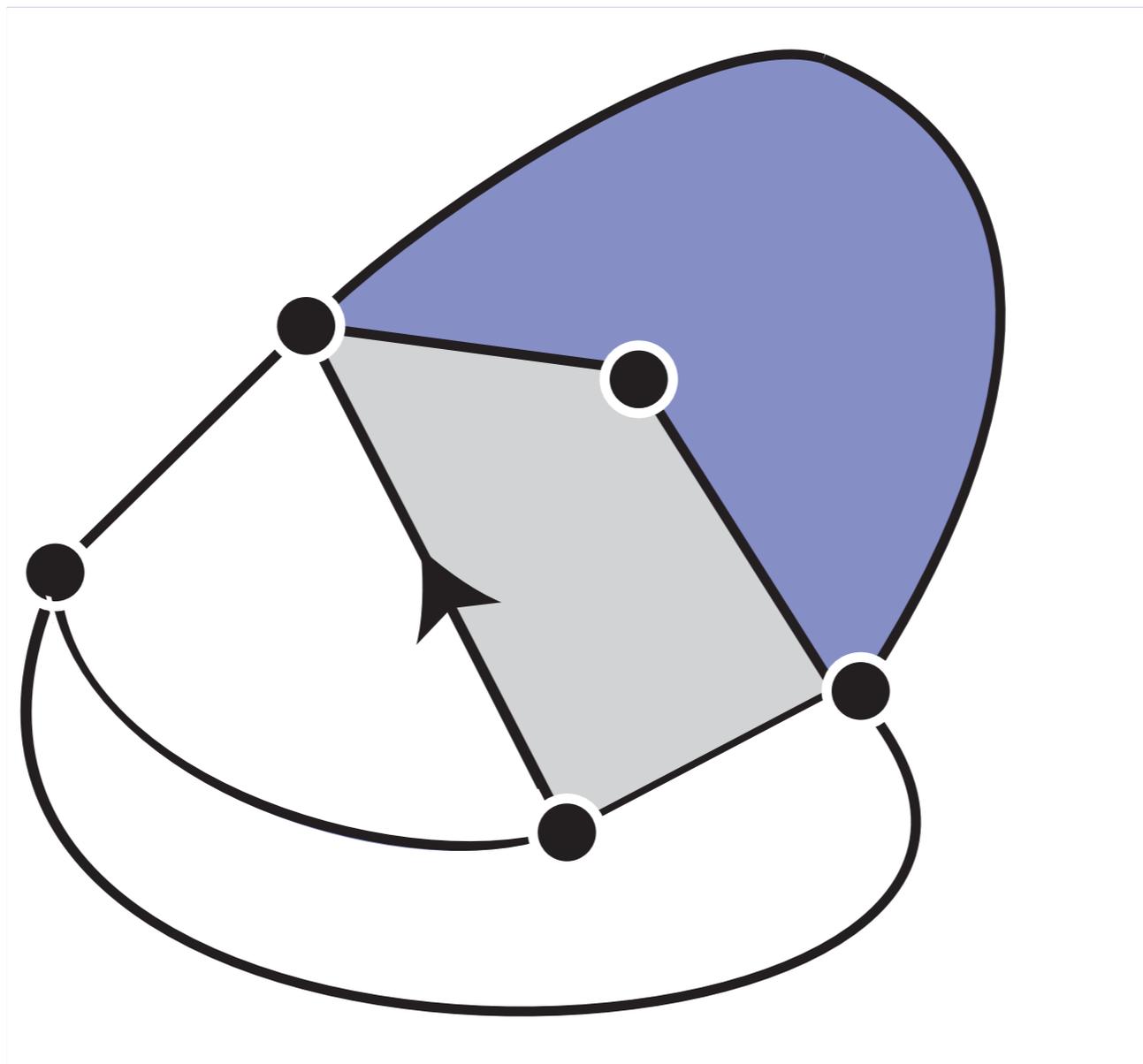
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



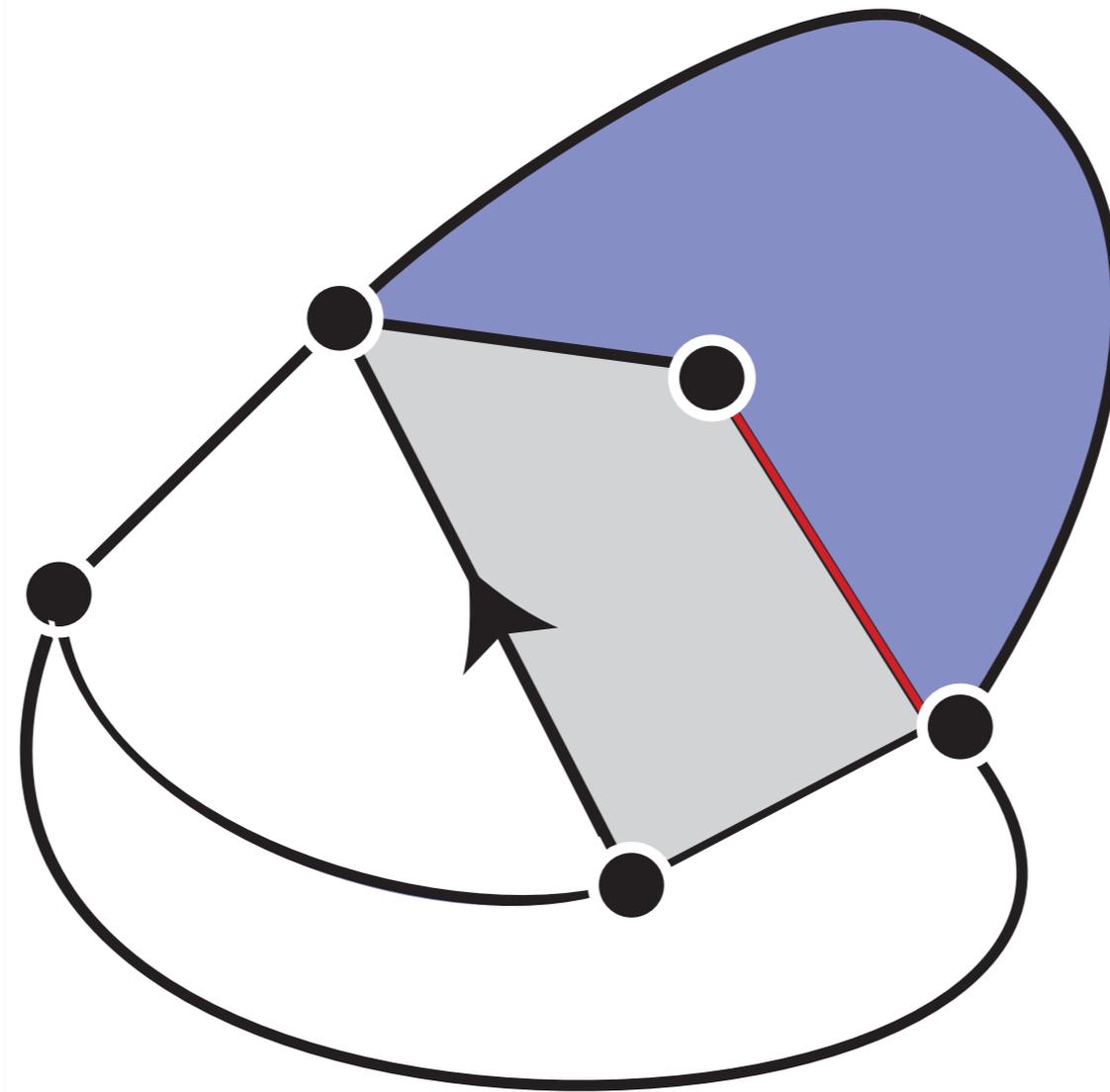
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



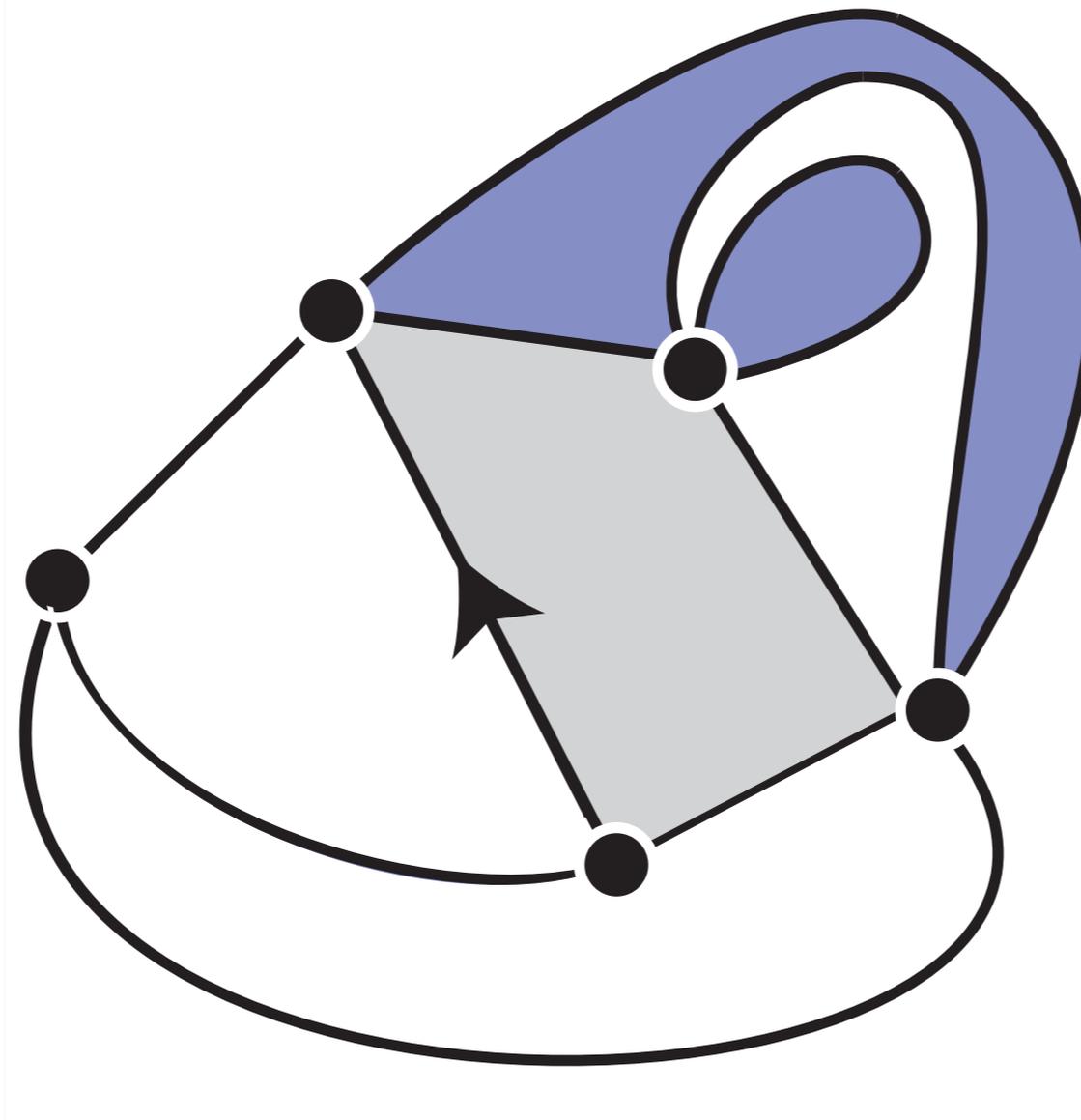
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



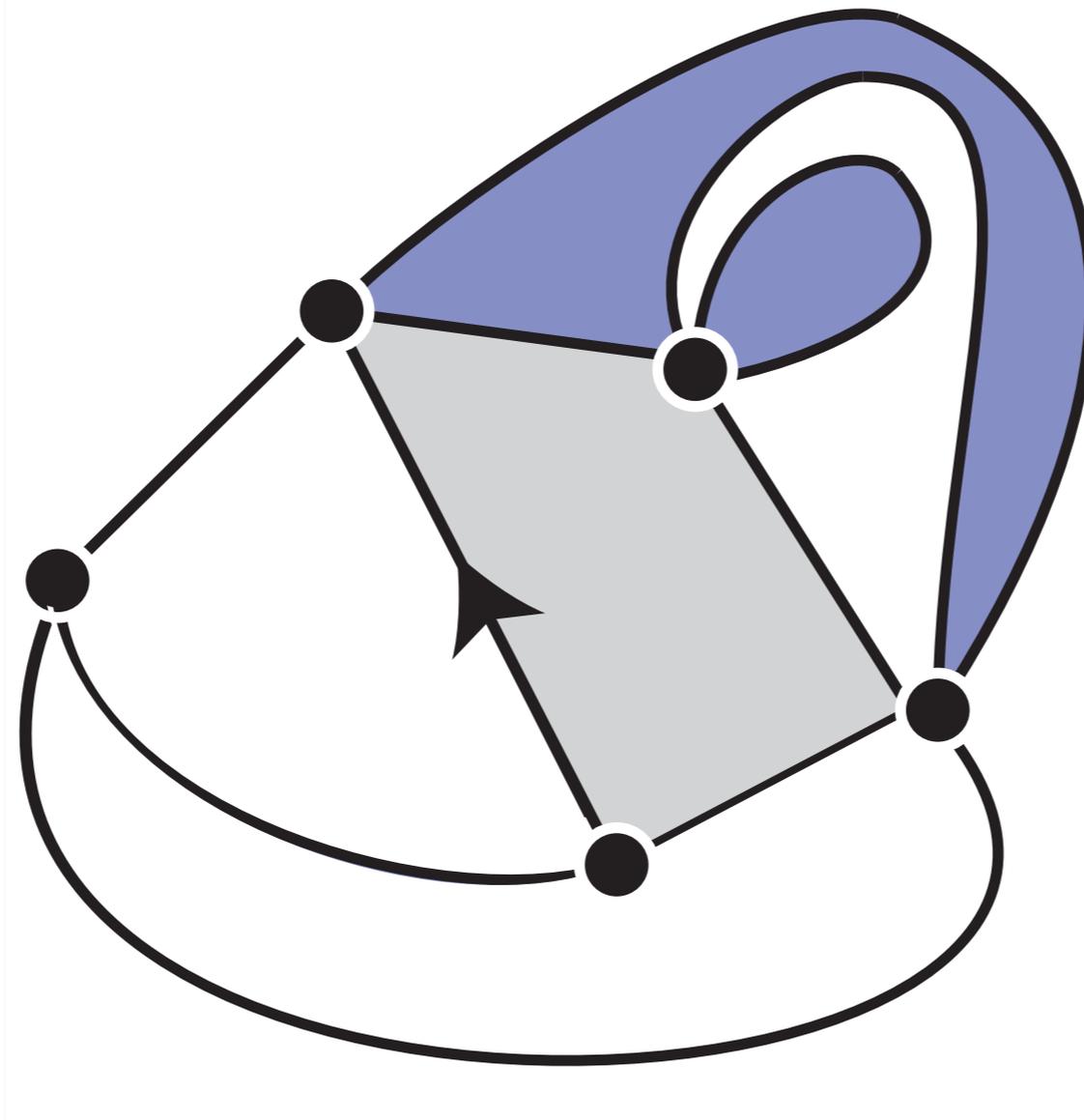
Branching peeling

Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



Branching peeling

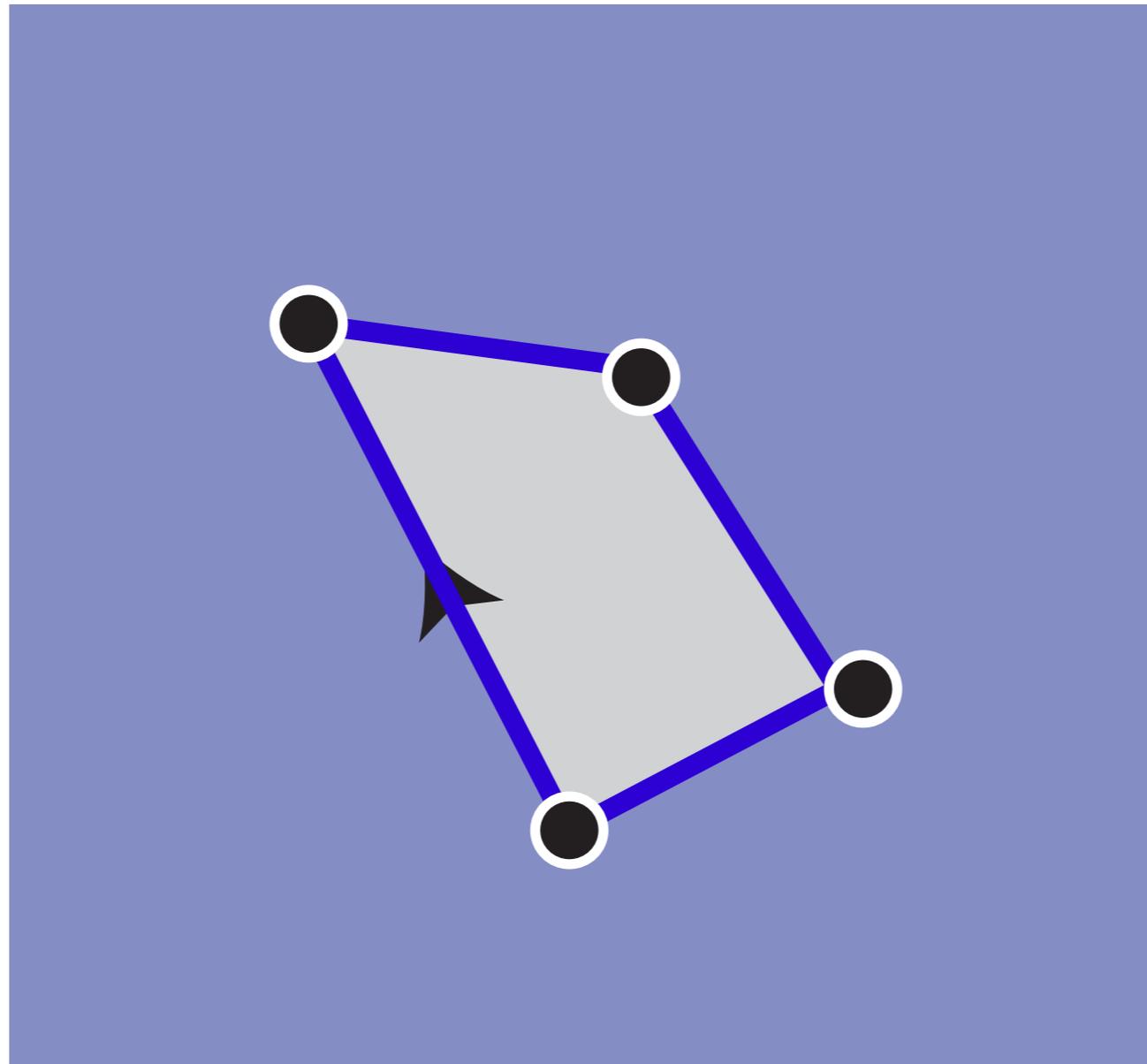
Intuitively, a **branching peeling** of a **triangulation with a boundary t** is an iterative exploration of t starting from the boundary and by discovering a new triangle at each step by *peeling an edge* using a deterministic algorithm \mathcal{A} .



And so on...

Following the locally largest cycle

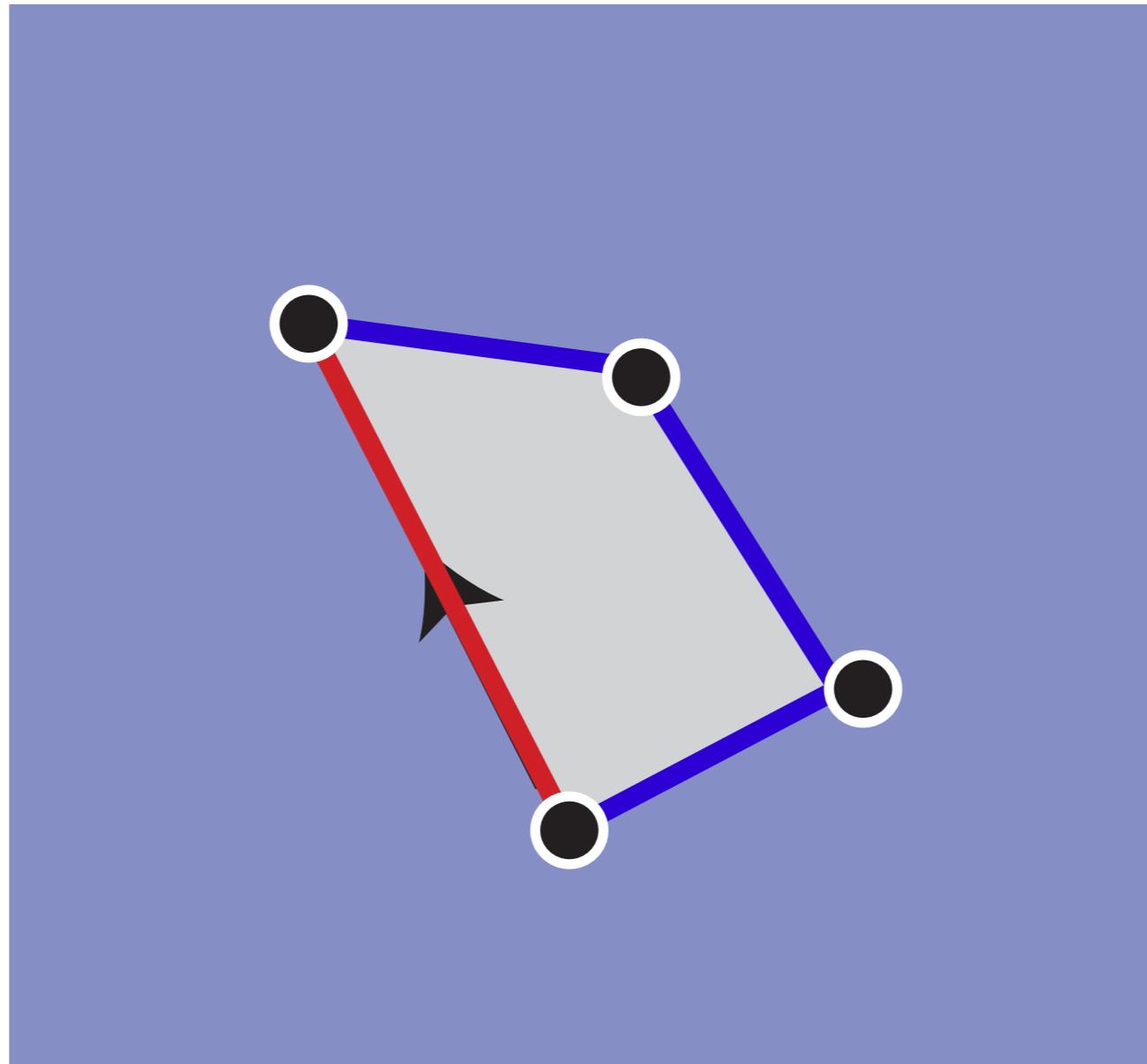
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4$$

Following the locally largest cycle

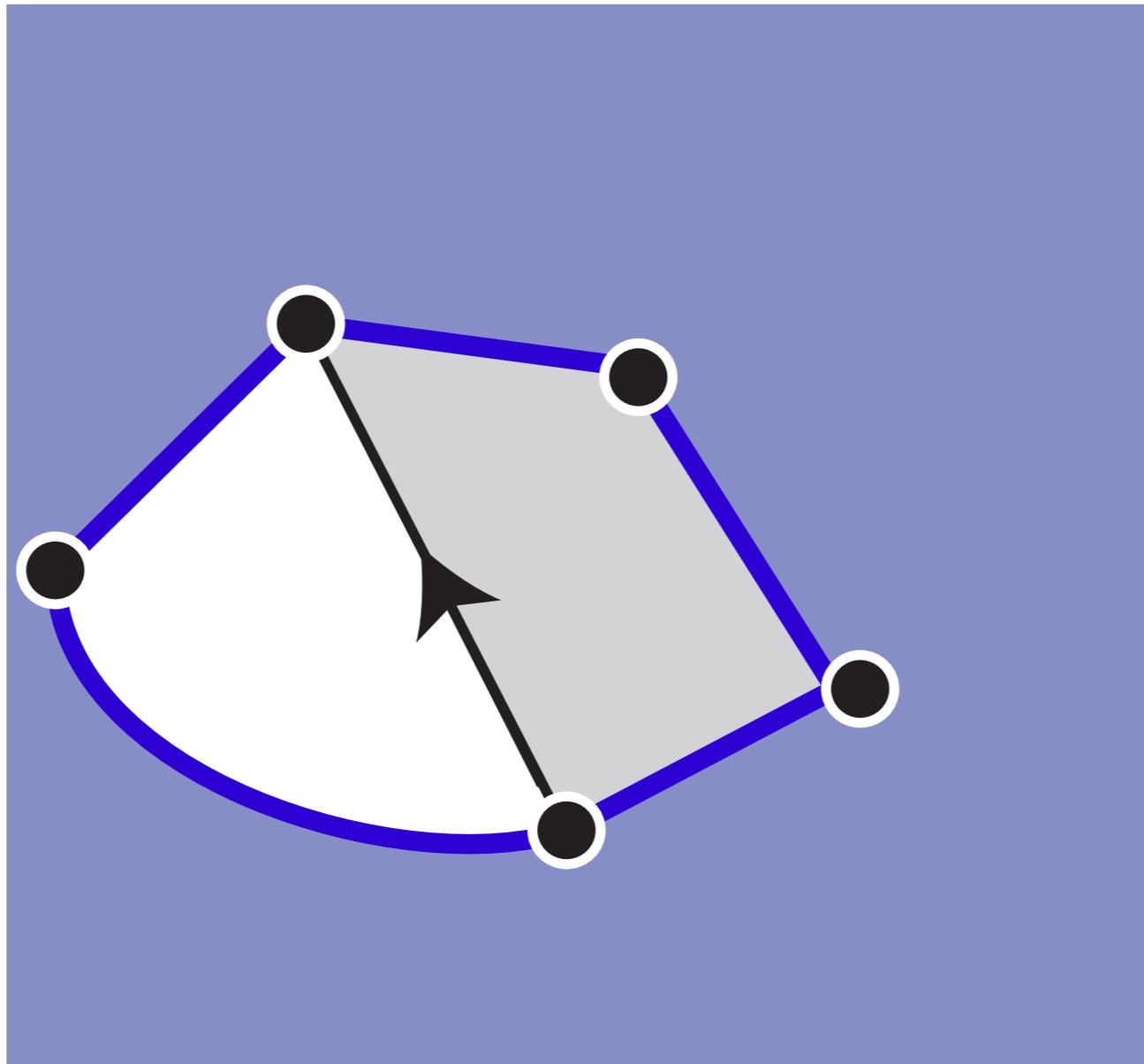
 **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4$$

Following the locally largest cycle

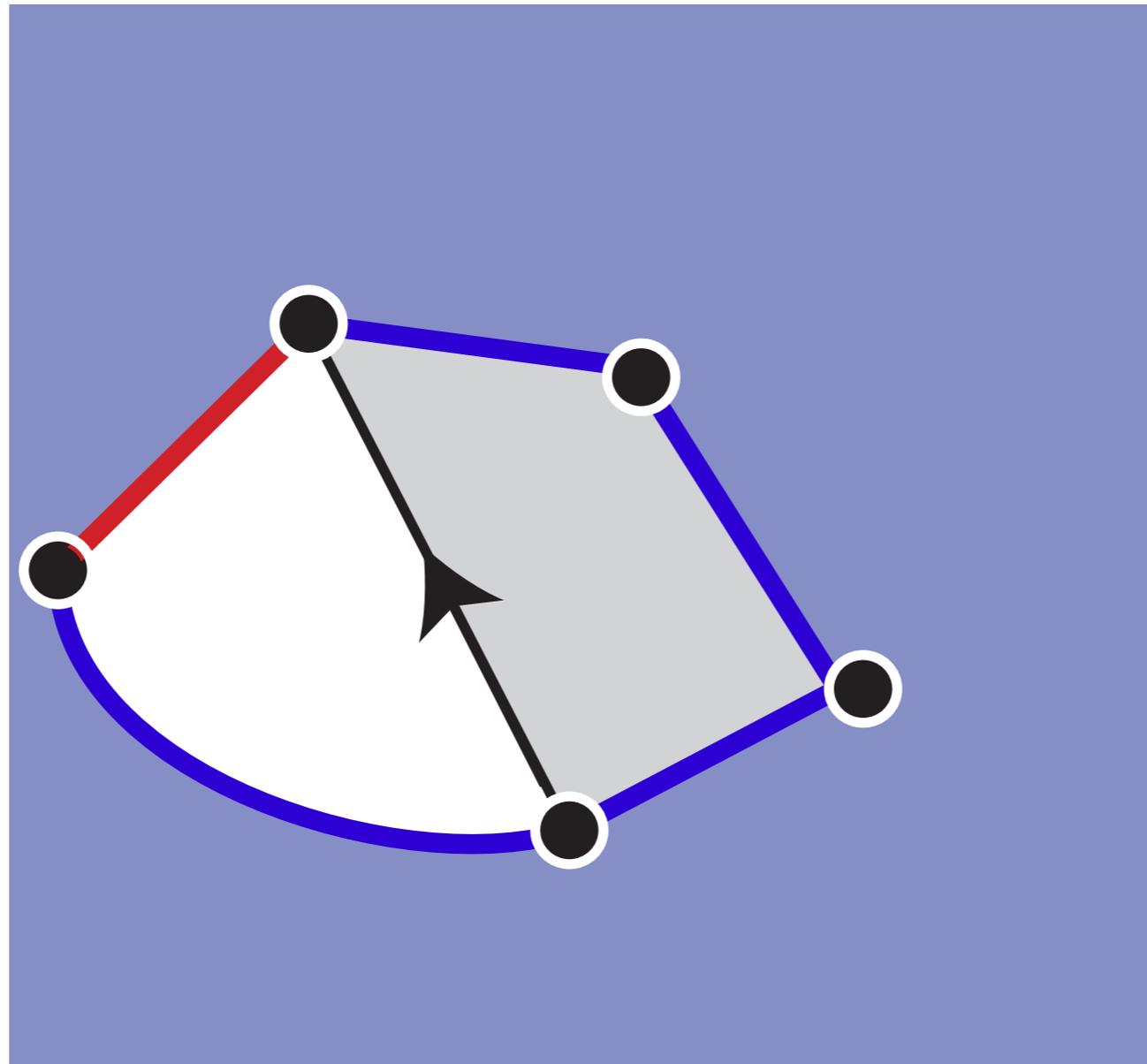
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5$$

Following the locally largest cycle

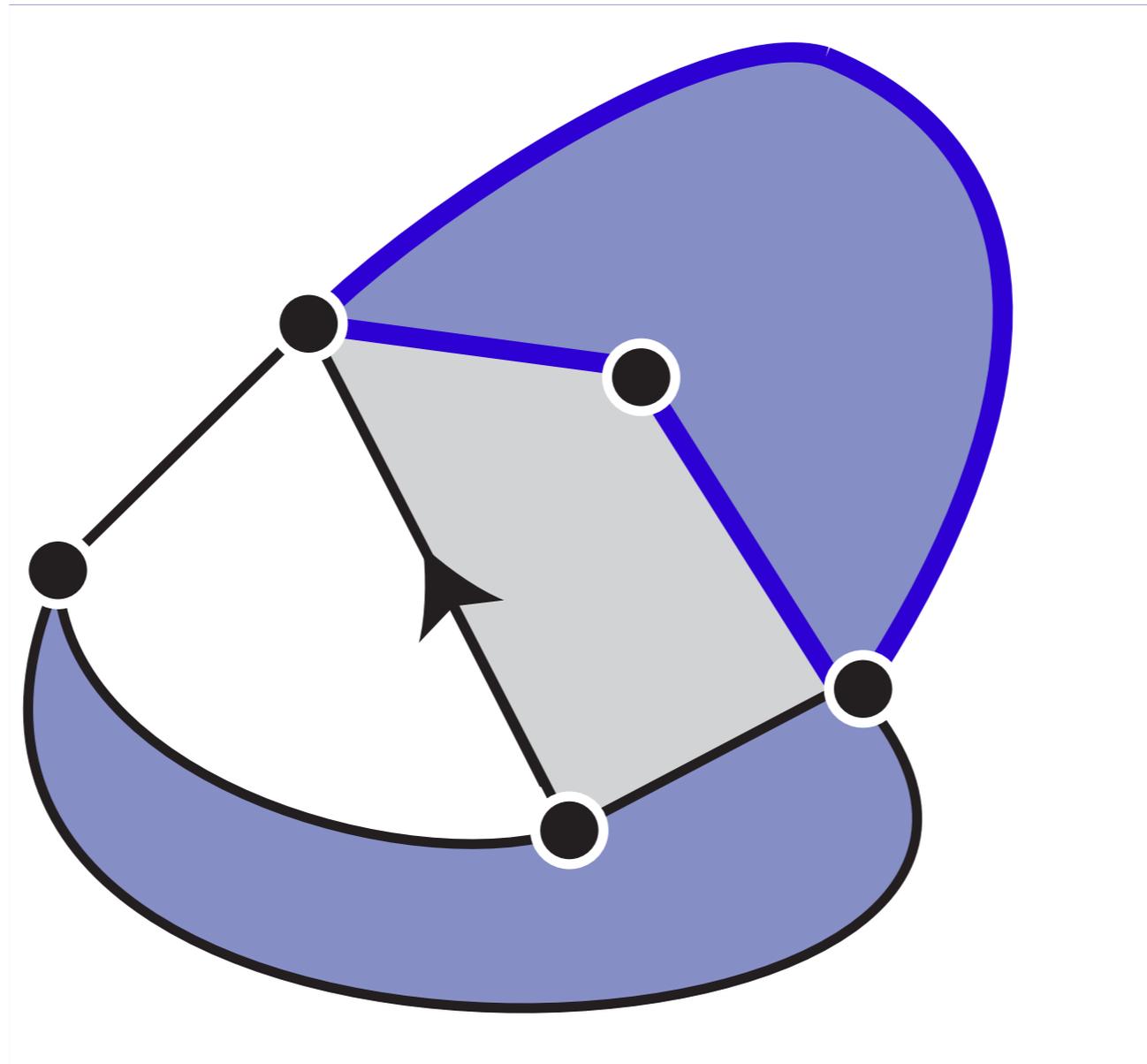
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5$$

Following the locally largest cycle

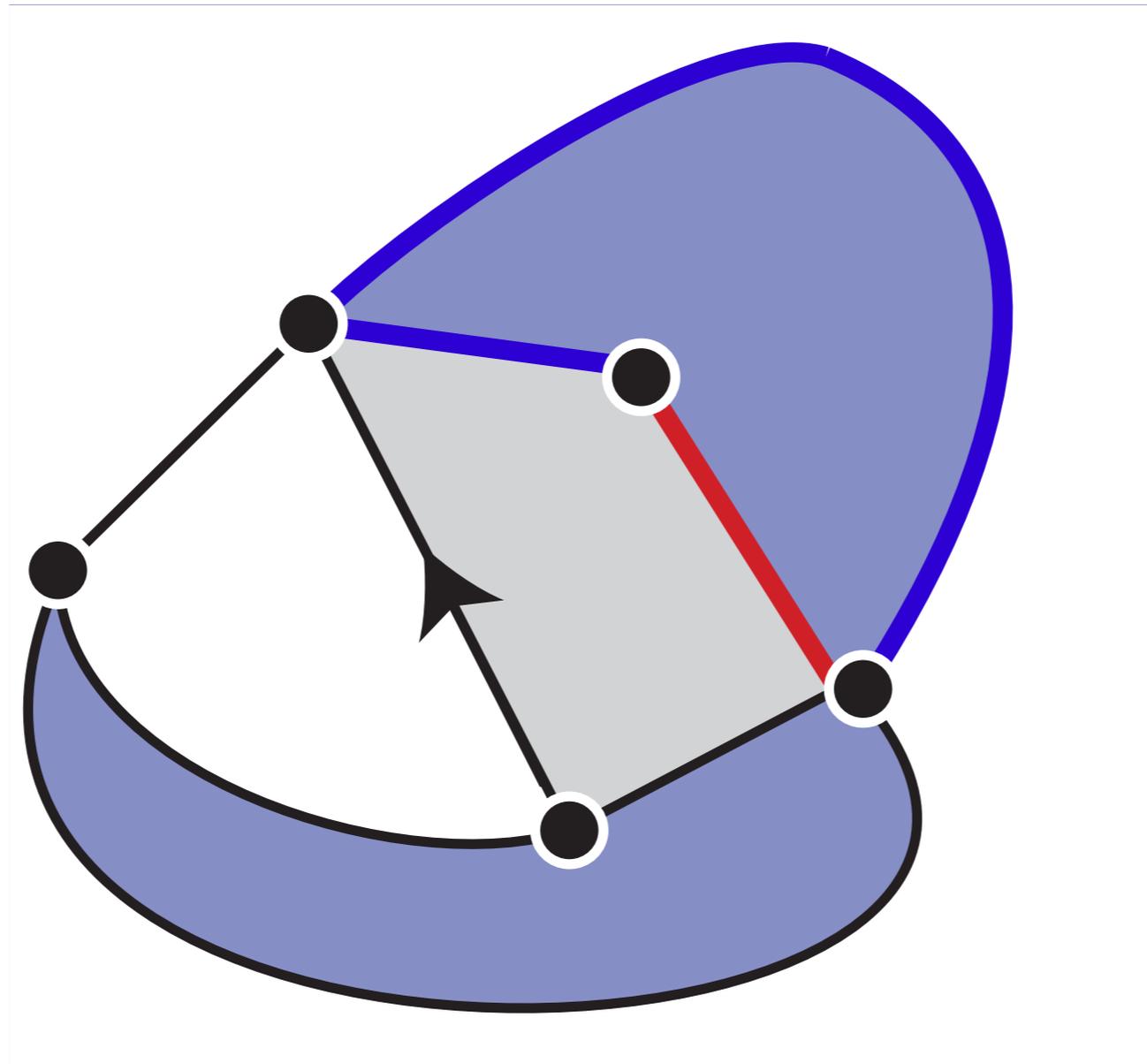
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5, \tilde{L}^{(4)}(2) = 3$$

Following the locally largest cycle

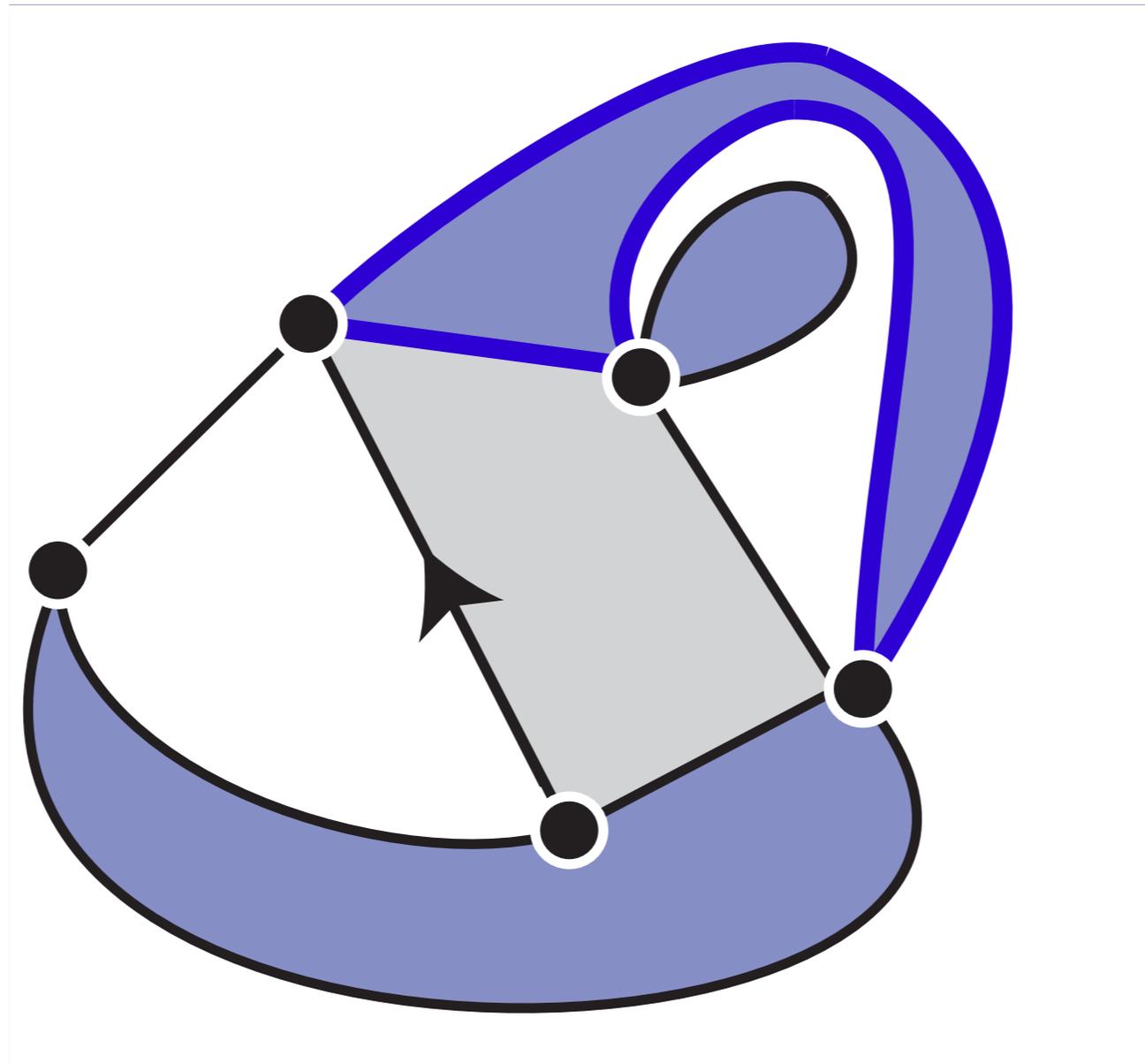
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5, \tilde{L}^{(4)}(2) = 3$$

Following the locally largest cycle

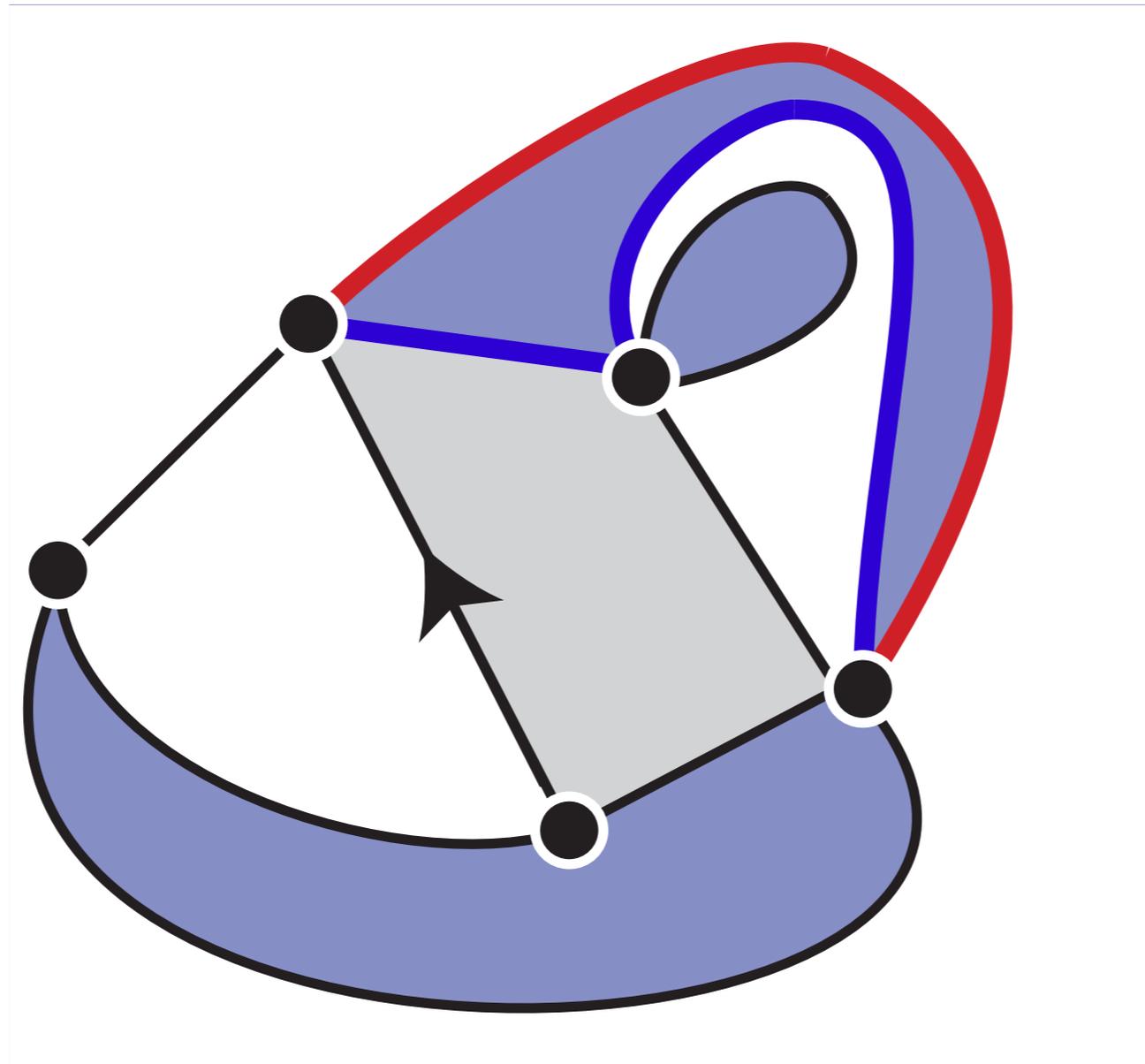
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5, \tilde{L}^{(4)}(2) = 3, \tilde{L}^{(4)}(3) = 3$$

Following the locally largest cycle

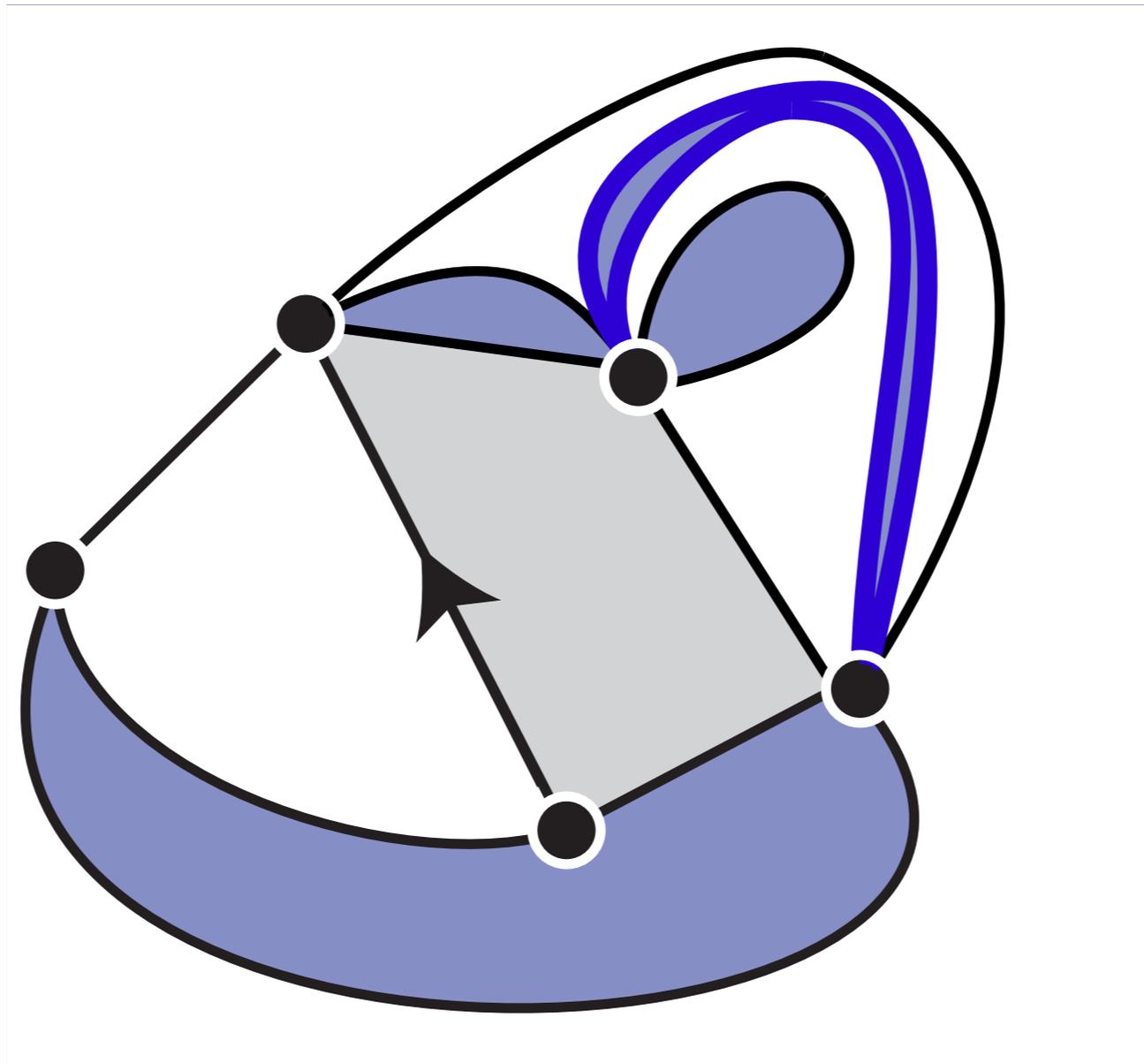
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5, \tilde{L}^{(4)}(2) = 3, \tilde{L}^{(4)}(3) = 3$$

Following the locally largest cycle

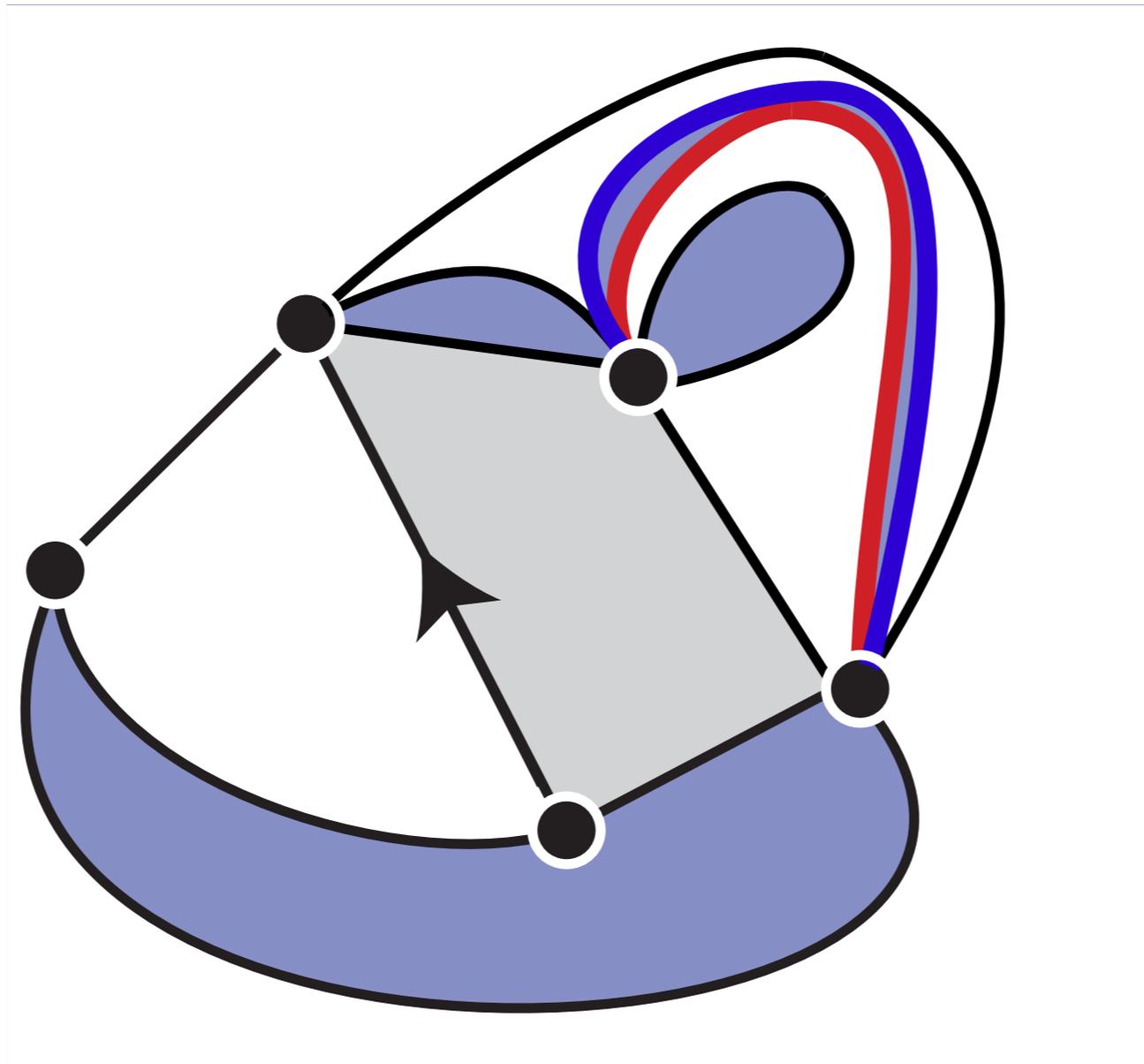
↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5, \tilde{L}^{(4)}(2) = 3, \tilde{L}^{(4)}(3) = 3, \tilde{L}^{(4)}(4) = 2$$

Following the locally largest cycle

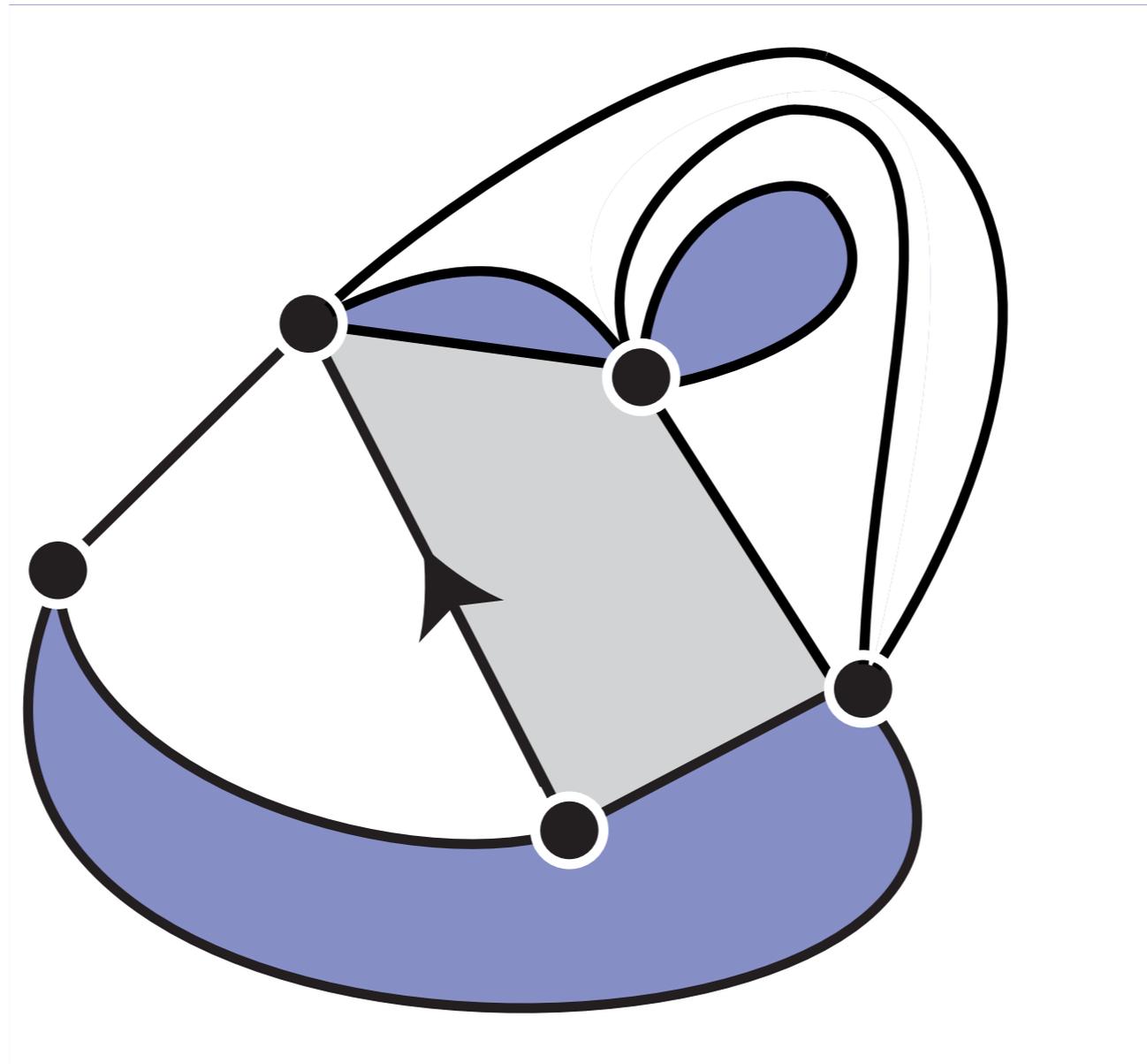
↗ Idea: at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5, \tilde{L}^{(4)}(2) = 3, \tilde{L}^{(4)}(3) = 3, \tilde{L}^{(4)}(4) = 2$$

Following the locally largest cycle

↗ **Idea:** at each peeling step, peel along the current locally largest cycle. Let $\tilde{L}^{(p)}(i)$ its length after i peeling steps.



$$\tilde{L}^{(4)}(0) = 4, \tilde{L}^{(4)}(1) = 5, \tilde{L}^{(4)}(2) = 3, \tilde{L}^{(4)}(3) = 3, \tilde{L}^{(4)}(4) = 2, \tilde{L}^{(4)}(5) = 0.$$

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathsf{T}^{(p)}$.

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathcal{T}^{(p)}$.

↪ **Key point:** $(\tilde{L}^{(p)}(i); i \geq 0)$ is a Markov chain starting at p , absorbed at 0 and with explicit transitions.

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathcal{T}^{(p)}$.

→ **Key point:** $(\tilde{L}^{(p)}(i); i \geq 0)$ is a Markov chain starting at p , absorbed at 0 and with explicit transitions. In addition, the triangulations filling-in the holes of non-explored regions are independent **Boltzmann triangulations with a boundary**.

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathcal{T}^{(p)}$.

↗ **Key point:** $(\tilde{L}^{(p)}(i); i \geq 0)$ is a Markov chain starting at p , absorbed at 0 and with explicit transitions. In addition, the triangulations filling-in the holes of non-explored regions are independent **Boltzmann triangulations with a boundary**.

If $L_{\text{height}}^{(p)}(r)$ is the length of the locally largest cycle at **height** r

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathcal{T}^{(p)}$.

↗ **Key point:** $(\tilde{L}^{(p)}(i); i \geq 0)$ is a Markov chain starting at p , absorbed at 0 and with explicit transitions. In addition, the triangulations filling-in the holes of non-explored regions are independent **Boltzmann triangulations with a boundary**.

If $L_{\text{height}}^{(p)}(r)$ is the length of the locally largest cycle at **height** r , using **Bertoin & K. '14** and **Curien & Le Gall '14**, we get that

Proposition (Bertoin, Curien & K. '15).

We have

$$\left(\frac{1}{p} L_{\text{height}}^{(p)}(\lfloor \sqrt{p} \cdot t \rfloor); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(X \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathcal{T}^{(p)}$.

↗ **Key point:** $(\tilde{L}^{(p)}(i); i \geq 0)$ is a Markov chain starting at p , absorbed at 0 and with explicit transitions. In addition, the triangulations filling-in the holes of non-explored regions are independent **Boltzmann triangulations with a boundary**.

If $L_{\text{height}}^{(p)}(r)$ is the length of the locally largest cycle at **height** r , using **Bertoin & K. '14** and **Curien & Le Gall '14**, we get that

Proposition (Bertoin, Curien & K. '15).

We have

$$\left(\frac{1}{p} L_{\text{height}}^{(p)}(\lfloor \sqrt{p} \cdot t \rfloor); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(X \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

where X is a càdlàg self-similar Markov process with index $-1/2$

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathcal{T}^{(p)}$.

↗ **Key point:** $(\tilde{L}^{(p)}(i); i \geq 0)$ is a Markov chain starting at p , absorbed at 0 and with explicit transitions. In addition, the triangulations filling-in the holes of non-explored regions are independent **Boltzmann triangulations with a boundary**.

If $L_{\text{height}}^{(p)}(r)$ is the length of the locally largest cycle at **height** r , using **Bertoin & K. '14** and **Curien & Le Gall '14**, we get that

Proposition (Bertoin, Curien & K. '15).

We have

$$\left(\frac{1}{p} L_{\text{height}}^{(p)}(\lfloor \sqrt{p} \cdot t \rfloor); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(X \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

where X is a càdlàg self-similar Markov process with index $-1/2$ (i.e. $t \mapsto c \cdot X(c^{-1/2}t)$ has the same law as X started at c), with $X(0) = 1$ and only negative jumps

Scaling limit of the locally largest cycle

Recall that $\tilde{L}^{(p)}(i)$ is the length of the locally largest cycle after i peeling steps of $\mathcal{T}^{(p)}$.

↗ **Key point:** $(\tilde{L}^{(p)}(i); i \geq 0)$ is a Markov chain starting at p , absorbed at 0 and with explicit transitions. In addition, the triangulations filling-in the holes of non-explored regions are independent **Boltzmann triangulations with a boundary**.

If $L_{\text{height}}^{(p)}(r)$ is the length of the locally largest cycle at **height** r , using **Bertoin & K. '14** and **Curien & Le Gall '14**, we get that

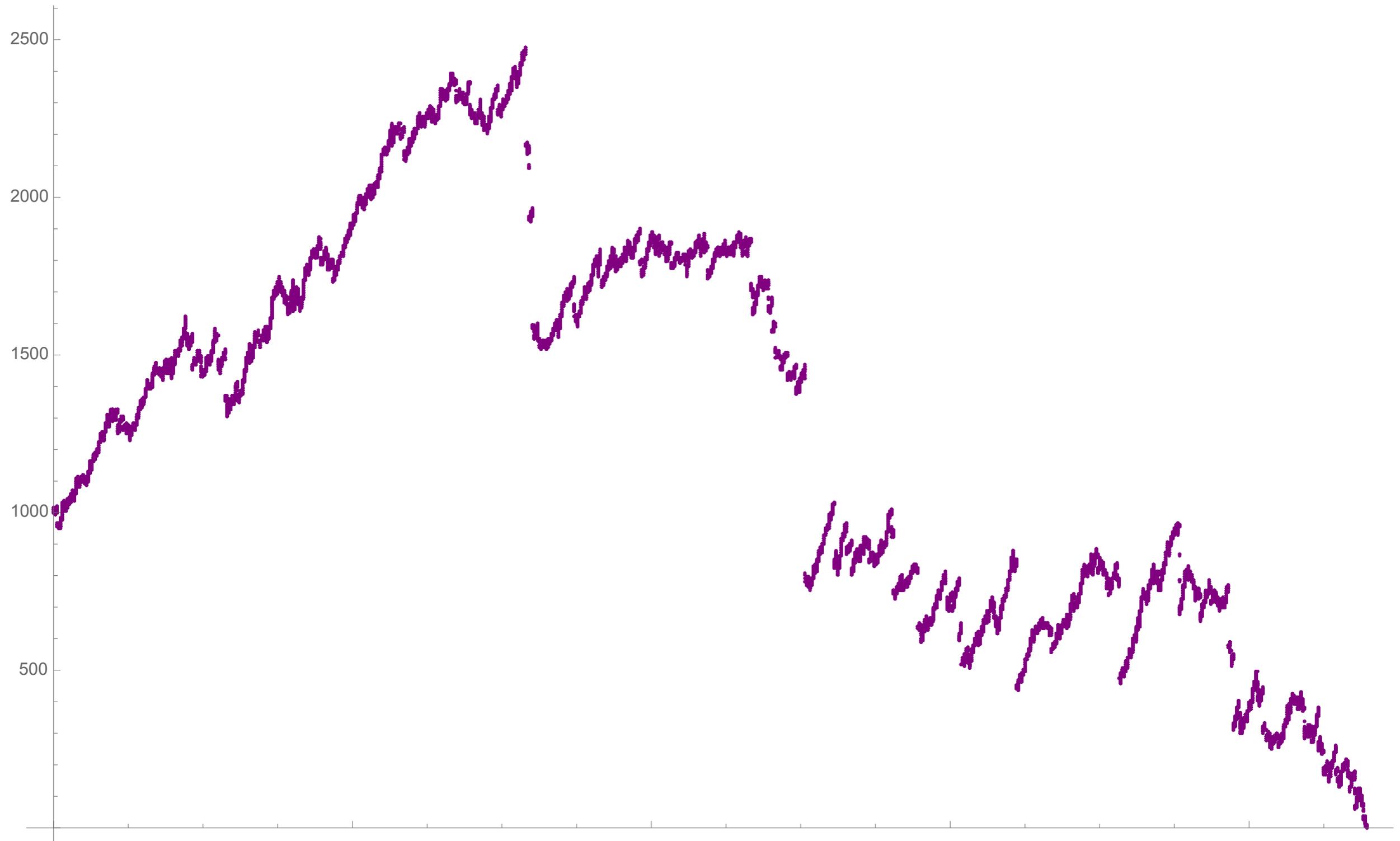
Proposition (Bertoin, Curien & K. '15).

We have

$$\left(\frac{1}{p} L_{\text{height}}^{(p)}(\lfloor \sqrt{p} \cdot t \rfloor); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(X \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

where X is a càdlàg self-similar Markov process with index $-1/2$ and absorbed at 0.

A simulation of X



The self-similar Markov process X

Let ξ be a spectrally negative Lévy process with Laplace exponent

$$\Psi(q) = -\frac{8}{3}q + \int_{1/2}^1 (x^q - 1 + q(1-x)) (x(1-x))^{-5/2} dx,$$

so that $\mathbb{E}[\exp(q\xi(t))] = \exp(t\Psi(q))$ for every $t \geq 0$, $q \geq 0$ and $\xi(t) \rightarrow -\infty$ when $t \rightarrow \infty$.

The self-similar Markov process X

Let ξ be a spectrally negative Lévy process with Laplace exponent

$$\Psi(q) = -\frac{8}{3}q + \int_{1/2}^1 (x^q - 1 + q(1-x)) (x(1-x))^{-5/2} dx,$$

so that $\mathbb{E}[\exp(q\xi(t))] = \exp(t\Psi(q))$ for every $t \geq 0$, $q \geq 0$ and $\xi(t) \rightarrow -\infty$ when $t \rightarrow \infty$.

Then

$$X(t) = \exp(\xi(\tau(t))), \quad t \geq 0$$

The self-similar Markov process X

Let ξ be a spectrally negative Lévy process with Laplace exponent

$$\Psi(q) = -\frac{8}{3}q + \int_{1/2}^1 (x^q - 1 + q(1-x)) (x(1-x))^{-5/2} dx,$$

so that $\mathbb{E}[\exp(q\xi(t))] = \exp(t\Psi(q))$ for every $t \geq 0$, $q \geq 0$ and $\xi(t) \rightarrow -\infty$ when $t \rightarrow \infty$.

Set

$$\tau(t) = \inf \left\{ u \geq 0; \int_0^u e^{\xi(s)/2} ds > t \right\}, \quad t \geq 0$$

with the convention $\inf \emptyset = \infty$, i.e. $\tau(t) = \infty$ when $t \geq \int_0^\infty e^{\xi(s)/2} ds$.

Then

$$X(t) = \exp(\xi(\tau(t))), \quad t \geq 0$$

The self-similar Markov process X

Let ξ be a spectrally negative Lévy process with Laplace exponent

$$\Psi(q) = -\frac{8}{3}q + \int_{1/2}^1 (x^q - 1 + q(1-x)) (x(1-x))^{-5/2} dx,$$

so that $\mathbb{E}[\exp(q\xi(t))] = \exp(t\Psi(q))$ for every $t \geq 0$, $q \geq 0$ and $\xi(t) \rightarrow -\infty$ when $t \rightarrow \infty$.

Set

$$\tau(t) = \inf \left\{ u \geq 0; \int_0^u e^{\xi(s)/2} ds > t \right\}, \quad t \geq 0$$

with the convention $\inf \emptyset = \infty$, i.e. $\tau(t) = \infty$ when $t \geq \int_0^\infty e^{\xi(s)/2} ds$.

Then

$$X(t) = \exp(\xi(\tau(t))), \quad t \geq 0$$

with the convention $\exp(\xi(\infty)) = 0$.

**DESCRIPTION OF THE LIMITING PROCESS: A
GROWTH-FRAGMENTATION PROCESS**



Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations.

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

Growth-fragmentations: genealogical vision

- We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .
- Start at time 0 with one cell of size 1, whose size evolves according to X .

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

- Start at time 0 with one cell of size 1, whose size evolves according to X . Interpret each (negative) jump of X as the division of a cell, that is if $\Delta X(t) = X(t) - X(t-) = -y < 0$, the cell divides at time t into a mother cell (with size $X(t)$) and one daughter cell (of size y).

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

– Start at time 0 with one cell of size 1, whose size evolves according to X .

Interpret each (negative) jump of X as the division of a cell, that is if

$\Delta X(t) = X(t) - X(t-) = -y < 0$, the cell divides at time t into a mother cell (with size $X(t)$) and one daughter cell (of size y).

 After the division, the size of the daughter cell evolves as an independent version of X (started from y)

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

– Start at time 0 with one cell of size 1, whose size evolves according to X .

Interpret each (negative) jump of X as the division of a cell, that is if $\Delta X(t) = X(t) - X(t-) = -y < 0$, the cell divides at time t into a mother cell (with size $X(t)$) and one daughter cell (of size y).

 After the division, the size of the daughter cell evolves as an independent version of X (started from y), independently of all the other evolutions.

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

– Start at time 0 with one cell of size 1, whose size evolves according to X .

Interpret each (negative) jump of X as the division of a cell, that is if $\Delta X(t) = X(t) - X(t-) = -y < 0$, the cell divides at time t into a mother cell (with size $X(t)$) and one daughter cell (of size y).

 After the division, the size of the daughter cell evolves as an independent version of X (started from y), independently of all the other evolutions.

And so one for the daughters, great grand-daughters, and so on...

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

– Start at time 0 with one cell of size 1, whose size evolves according to X .

Interpret each (negative) jump of X as the division of a cell, that is if

$\Delta X(t) = X(t) - X(t-) = -y < 0$, the cell divides at time t into a mother cell (with size $X(t)$) and one daughter cell (of size y).

 After the division, the size of the daughter cell evolves as an independent version of X (started from y), independently of all the other evolutions.

And so one for the daughters, great grand-daughters, and so on...

By [Bertoin '15](#), for every $t \geq 0$, the family of all the cells alive at time t is cube summable

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

– Start at time 0 with one cell of size 1, whose size evolves according to X .

Interpret each (negative) jump of X as the division of a cell, that is if

$\Delta X(t) = X(t) - X(t-) = -y < 0$, the cell divides at time t into a mother cell (with size $X(t)$) and one daughter cell (of size y).

\curvearrowright After the division, the size of the daughter cell evolves as an independent version of X (started from y), independently of all the other evolutions.

And so one for the daughters, great grand-daughters, and so on...

By [Bertoin '15](#), for every $t \geq 0$, the family of all the cells alive at time t is cube summable, and can thus be rearranged in decreasing order.

Growth-fragmentations: genealogical vision

We use X to define a self-similar growth-fragmentation process with binary dislocations. We view $X(t)$ as the size of a typical particle or cell at age t .

– Start at time 0 with one cell of size 1, whose size evolves according to X .

Interpret each (negative) jump of X as the division of a cell, that is if $\Delta X(t) = X(t) - X(t-) = -y < 0$, the cell divides at time t into a mother cell (with size $X(t)$) and one daughter cell (of size y).

 After the division, the size of the daughter cell evolves as an independent version of X (started from y), independently of all the other evolutions.

And so one for the daughters, great grand-daughters, and so on...

By [Bertoin '15](#), for every $t \geq 0$, the family of all the cells alive at time t is cube summable, and can thus be rearranged in decreasing order. This yields a random variable with values in ℓ_3^\downarrow denoted by $\mathbb{X}(t) = (X_1(t), X_2(t), \dots)$.

Growth-fragmentations: temporal vision

One can view \mathbb{X} as the evolution of particle sizes that grow and divide as time passes:

Growth-fragmentations: temporal vision

One can view \mathbb{X} as the evolution of particle sizes that grow and divide as time passes:

 \mathbb{X} satisfies a branching property and is self-similar with index $-1/2$

Growth-fragmentations: temporal vision

One can view \mathbb{X} as the evolution of particle sizes that grow and divide as time passes:

 \mathbb{X} satisfies a branching property and is self-similar with index $-1/2$, that is for every $c > 0$, the process $(c\mathbb{X}(c^{-1/2}t), t \geq 0)$ has the same law as \mathbb{X} starting from $(c, 0, 0, \dots)$.

Growth-fragmentations: temporal vision

One can view \mathbb{X} as the evolution of particle sizes that grow and divide as time passes:

→ \mathbb{X} satisfies a branching property and is self-similar with index $-1/2$, that is for every $c > 0$, the process $(c\mathbb{X}(c^{-1/2}t), t \geq 0)$ has the same law as \mathbb{X} starting from $(c, 0, 0, \dots)$.

→ The divisions of \mathbb{X} are binary, i.e. they amount to dividing m into smaller masses m_1 and m_2 with $m_1 + m_2 = m$.

Growth-fragmentations: temporal vision

One can view \mathbb{X} as the evolution of particle sizes that grow and divide as time passes:

→ \mathbb{X} satisfies a branching property and is self-similar with index $-1/2$, that is for every $c > 0$, the process $(c\mathbb{X}(c^{-1/2}t), t \geq 0)$ has the same law as \mathbb{X} starting from $(c, 0, 0, \dots)$.

→ The divisions of \mathbb{X} are binary, i.e. they amount to dividing m into smaller masses m_1 and m_2 with $m_1 + m_2 = m$. Informally, in \mathbb{X} , each size $m > 0$ divides into smaller masses $(xm, (1-x)m)$ at a rate $m^{-1/2}\nu(dx)$, with

$$\nu(dx) = (x(1-x))^{-5/2}dx, \quad x \in (1/2, 1)$$

Growth-fragmentations: temporal vision

One can view \mathbb{X} as the evolution of particle sizes that grow and divide as time passes:

↗ \mathbb{X} satisfies a branching property and is self-similar with index $-1/2$, that is for every $c > 0$, the process $(c\mathbb{X}(c^{-1/2}t), t \geq 0)$ has the same law as \mathbb{X} starting from $(c, 0, 0, \dots)$.

↗ The divisions of \mathbb{X} are binary, i.e. they amount to dividing m into smaller masses m_1 and m_2 with $m_1 + m_2 = m$. Informally, in \mathbb{X} , each size $m > 0$ divides into smaller masses $(xm, (1-x)m)$ at a rate $m^{-1/2}\nu(dx)$, with

$$\nu(dx) = (x(1-x))^{-5/2}dx, \quad x \in (1/2, 1)$$

↗ We have $\int^1 (1-x)^2 \nu(dx) < \infty$, but $\int^1 (1-x) \nu(dx) = \infty$ which underlines the necessity of compensating the dislocations.

An artistic representation of a growth-fragmentation

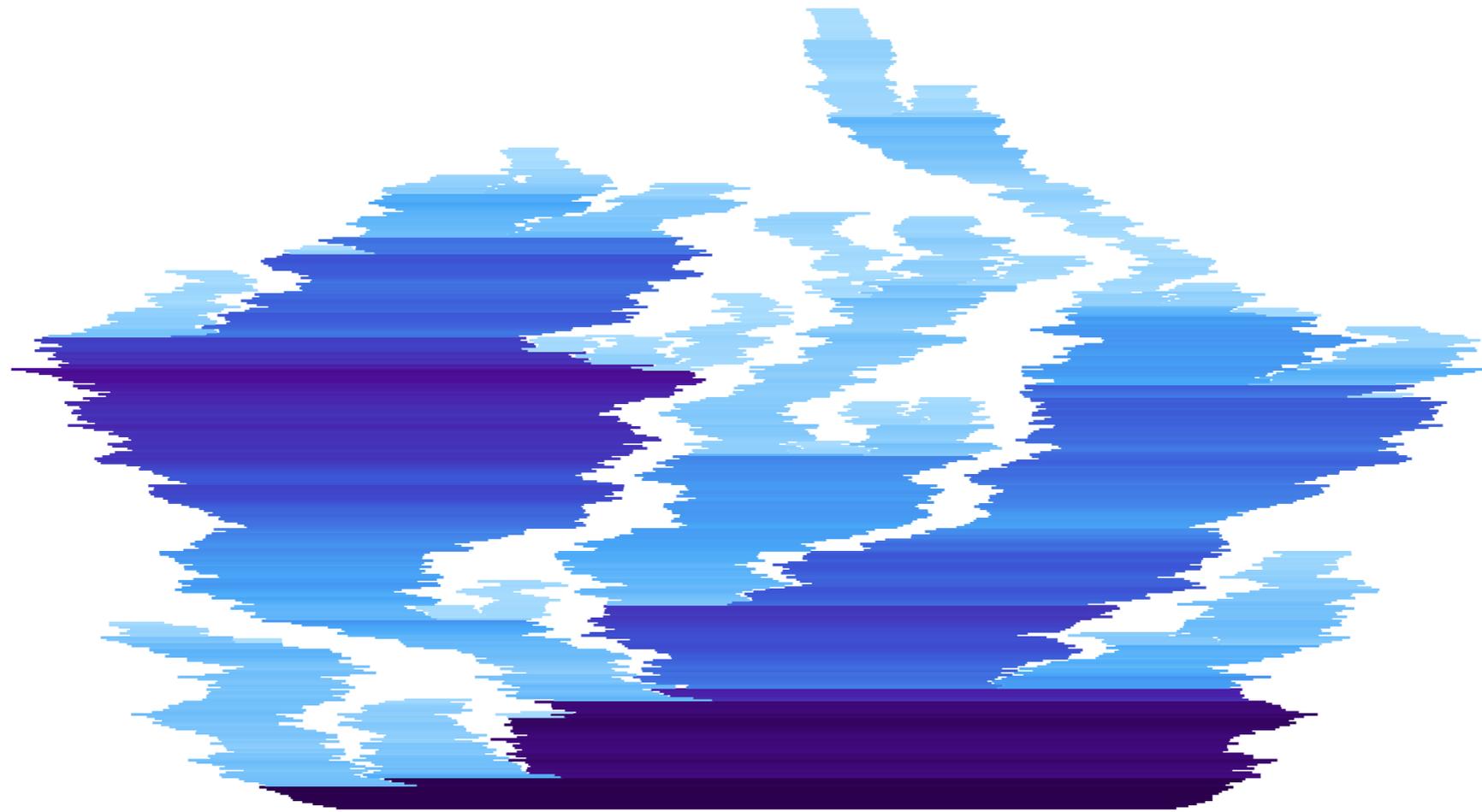


Figure: An artistic representation (by N. Curien) of the cycle lengths at fixed heights of a Boltzmann triangulation with a large boundary: horizontal segments correspond to cycle lengths (the darker the cycle is, the longer it is).

The theorem

Recall that $\mathbb{L}^{(p)}(r) = \left(L_1^{(p)}(r), L_2^{(p)}(r), \dots \right)$ are the lengths of the cycles of $B_r(\mathbb{T}^{(p)})$ ranked in decreasing order.

Theorem (Bertoin, Curien, K. '15).

We have

$$\left(\frac{1}{p} \cdot \mathbb{L}^{(p)}(t\sqrt{p}); t \geq 0 \right) \xrightarrow[p \rightarrow \infty]{(d)} \left(\mathbb{X} \left(\frac{3}{2\sqrt{\pi}} \cdot t \right); t \geq 0 \right),$$

in distribution in ℓ_3^\downarrow , where $\mathbb{X} = (\mathbb{X}(t); t \geq 0)$ is a càdlàg process with values in ℓ_3^\downarrow , which is a *self-similar growth-fragmentation process* (Bertoin '15).

Characterization of the growth-fragmentation

↗ The law of the cell process does not characterize the law of the **growth-fragmentation**.

Characterization of the growth-fragmentation

↗ The law of the cell process does not characterize the law of the **growth-fragmentation**.

However by Shi '15, the law of the **growth-fragmentation** is characterized by the so called **cumulant function** κ defined by

$$\kappa(q) = \Psi(q) + \int_{(-\infty, 0)} (1 - e^y)^q \Lambda(dy),$$

where Ψ is the Laplace exponent of the Lévy process associated to the self-similar cell process and Λ is its Lévy measure.

Characterization of the growth-fragmentation

↗ The law of the cell process does not characterize the law of the **growth-fragmentation**.

However by Shi '15, the law of the **growth-fragmentation** is characterized by the so called **cumulant function** κ defined by

$$\kappa(q) = \Psi(q) + \int_{(-\infty, 0)} (1 - e^y)^q \Lambda(dy),$$

where Ψ is the Laplace exponent of the Lévy process associated to the self-similar cell process and Λ is its Lévy measure.

In our case,

$$\kappa(q) = \frac{4\sqrt{\pi}}{3} \frac{\Gamma(q - \frac{3}{2})}{\Gamma(q - 3)}, \quad q > 3/2.$$

Characterization of the growth-fragmentation

↗ The law of the cell process does not characterize the law of the **growth-fragmentation**.

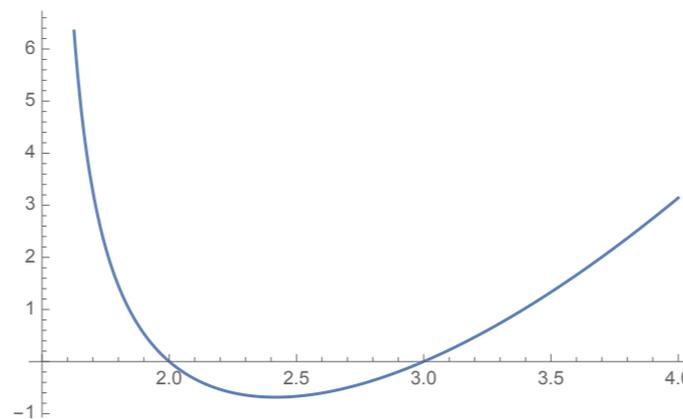
However by Shi '15, the law of the **growth-fragmentation** is characterized by the so called **cumulant function** κ defined by

$$\kappa(q) = \Psi(q) + \int_{(-\infty, 0)} (1 - e^y)^q \Lambda(dy),$$

where Ψ is the Laplace exponent of the Lévy process associated to the self-similar cell process and Λ is its Lévy measure.

In our case,

$$\kappa(q) = \frac{4\sqrt{\pi}}{3} \frac{\Gamma(q - \frac{3}{2})}{\Gamma(q - 3)}, \quad q > 3/2.$$



Useful tools: martingales

Bertoin, Budd, Curien, K:

 Zeros of the **cumulant function** allow to define **martingales**. In our case, two martingales: one for $\omega_- = 2$ and one for $\omega_+ = 3$.

Useful tools: martingales

Bertoin, Budd, Curien, K:

→ Zeros of the **cumulant function** allow to define **martingales**. In our case, two martingales: one for $\omega_- = 2$ and one for $\omega_+ = 3$.

→ These **martingales** can be used to bias the genealogical structure *à la* Lyons–Pemantle–Peres.

Useful tools: martingales

Bertoin, Budd, Curien, K:

-  Zeros of the **cumulant function** allow to define **martingales**. In our case, two martingales: one for $\omega_- = 2$ and one for $\omega_+ = 3$.
-  These **martingales** can be used to bias the genealogical structure *à la* Lyons–Pemantle–Peres.
-  The evolution of the size of the tagged cell when biasing with the **martingale** associated with $\omega_- = 2$ is a **spectrally negative 3/2-stable process conditioned to die at 0 continuously** (Caballero & Chaumont), which can be interpreted as the evolution of the cycle targeting a random leaf.

Useful tools: martingales

Bertoin, Budd, Curien, K:

↗ Zeros of the **cumulant function** allow to define **martingales**. In our case, two martingales: one for $\omega_- = 2$ and one for $\omega_+ = 3$.

↗ These **martingales** can be used to bias the genealogical structure *à la* Lyons–Pemantle–Peres.

↗ The evolution of the size of the tagged cell when biasing with the **martingale** associated with $\omega_- = 2$ is a **spectrally negative 3/2-stable process conditioned to die at 0 continuously** (Caballero & Chaumont), which can be interpreted as the evolution of the cycle targeting a random leaf.

↗ Conversely, if one assumes that the evolution of the tagged cell when biasing with the **martingale** associated with ω_- is a **spectrally negative α -stable process conditioned to die at 0 continuously**, then $\alpha = 3/2$ and

$$\kappa(q) = \frac{4\sqrt{\pi}}{3} \frac{\Gamma(q - \frac{3}{2})}{\Gamma(q - 3)}, \quad q > 3/2 \text{ (use Kuznetsov \& Pardo).}$$

EXTENSION TO OTHER MODELS OF PLANAR MAPS



Extension to other models

In [Bertoin, Budd, Curien, K](#), we consider a different family of random planar maps which have large degrees

Extension to other models

In Bertoin, Budd, Curien, K, we consider a different family of random planar maps which have large degrees, for which the level set process scales to a one parameter family of **self-similar growth-fragmentations** with **cumulant functions** $(\kappa_\theta)_{1/2 < \theta \leq 3/2}$ given by

$$\kappa_\theta(q) = \frac{\cos(\pi(q - \theta))}{\sin(\pi(q - 2\theta))} \cdot \frac{\Gamma(q - \theta)}{\Gamma(q - 2\theta)}, \quad \theta < q < 2\theta + 1.$$

Extension to other models

In Bertoin, Budd, Curien, K, we consider a different family of random planar maps which have large degrees, for which the level set process scales to a one parameter family of **self-similar growth-fragmentations** with **cumulant functions** $(\kappa_\theta)_{1/2 < \theta \leq 3/2}$ given by

$$\kappa_\theta(q) = \frac{\cos(\pi(q - \theta))}{\sin(\pi(q - 2\theta))} \cdot \frac{\Gamma(q - \theta)}{\Gamma(q - 2\theta)}, \quad \theta < q < 2\theta + 1.$$

In this case $\omega_- = \theta + 1/2$, $\omega_+ = \theta + 3/2$, and the evolution of the size of the tagged cell when biasing with the **martingale** associated to ω_- is a **θ -stable process**, with **positivity parameter** ρ such that $\theta(1 - \rho) = 1/2$, conditioned die at 0 continuously.

Extension to other models

In Bertoin, Budd, Curien, K, we consider a different family of random planar maps which have large degrees, for which the level set process scales to a one parameter family of **self-similar growth-fragmentations** with **cumulant functions** $(\kappa_\theta)_{1/2 < \theta \leq 3/2}$ given by

$$\kappa_\theta(q) = \frac{\cos(\pi(q - \theta))}{\sin(\pi(q - 2\theta))} \cdot \frac{\Gamma(q - \theta)}{\Gamma(q - 2\theta)}, \quad \theta < q < 2\theta + 1.$$

In this case $\omega_- = \theta + 1/2$, $\omega_+ = \theta + 3/2$, and the evolution of the size of the tagged cell when biasing with the **martingale** associated to ω_- is a **θ -stable process**, with **positivity parameter** ρ such that $\theta(1 - \rho) = 1/2$, conditioned die at 0 continuously.

Question. Find the asymptotic behavior of the tail of the extinction time of these growth-fragmentations.