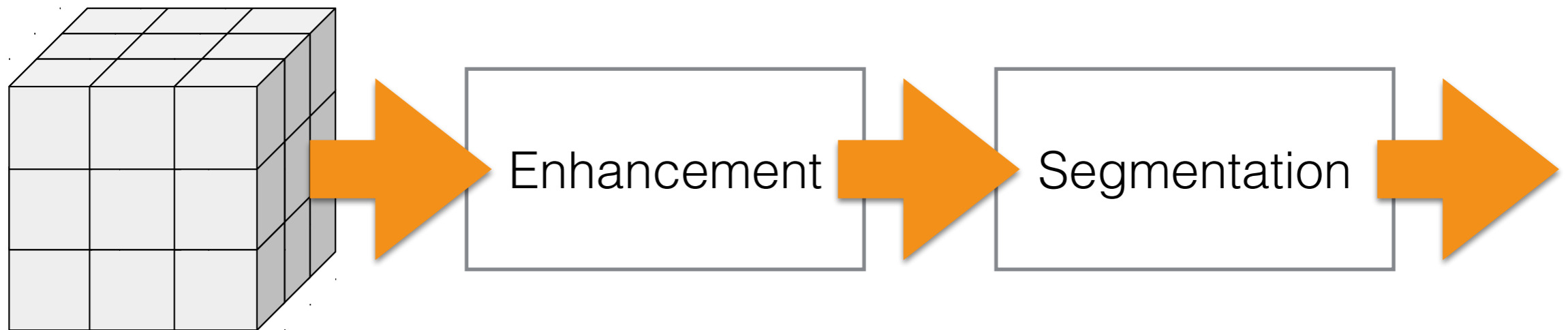


3D from Volume: Part II

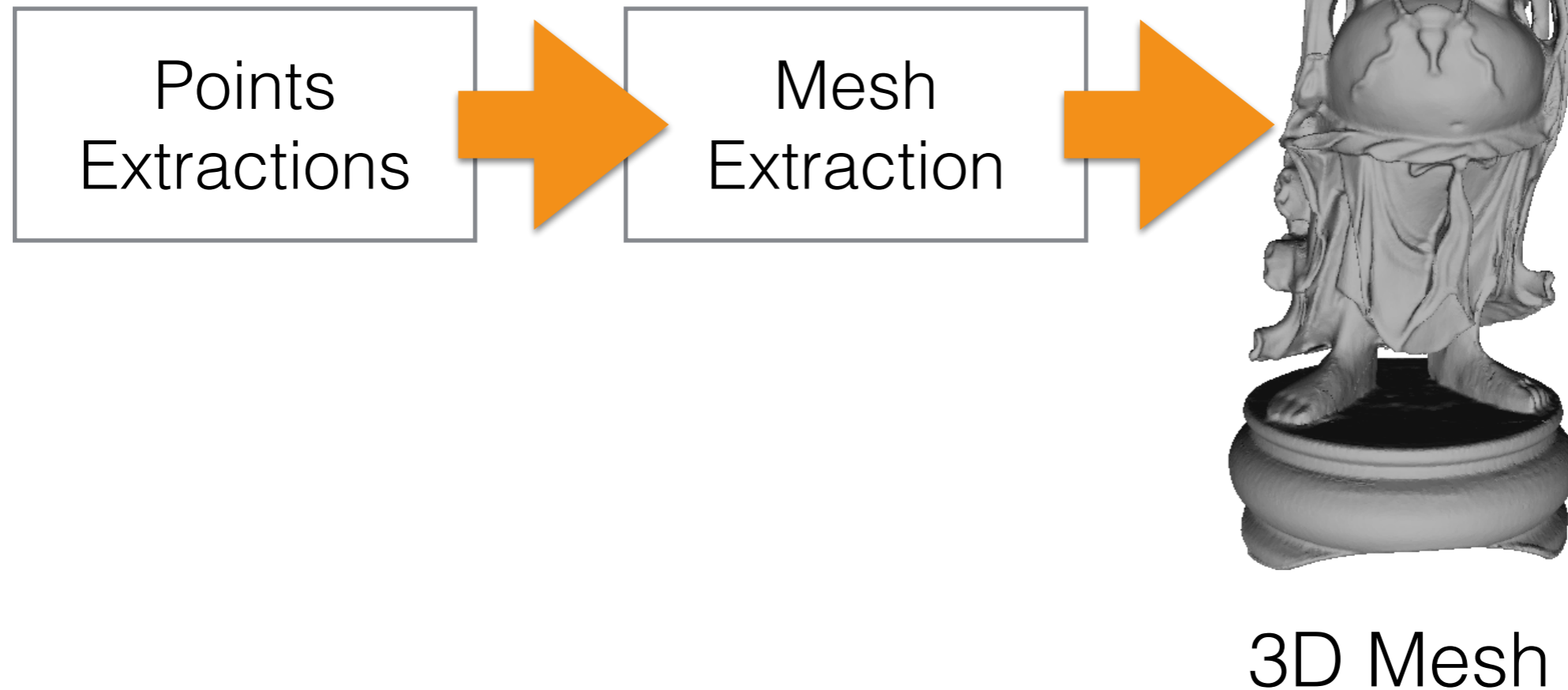
Francesco Banterle, Ph.D.
francesco.banterle@isti.cnr.it

The Processing Pipeline

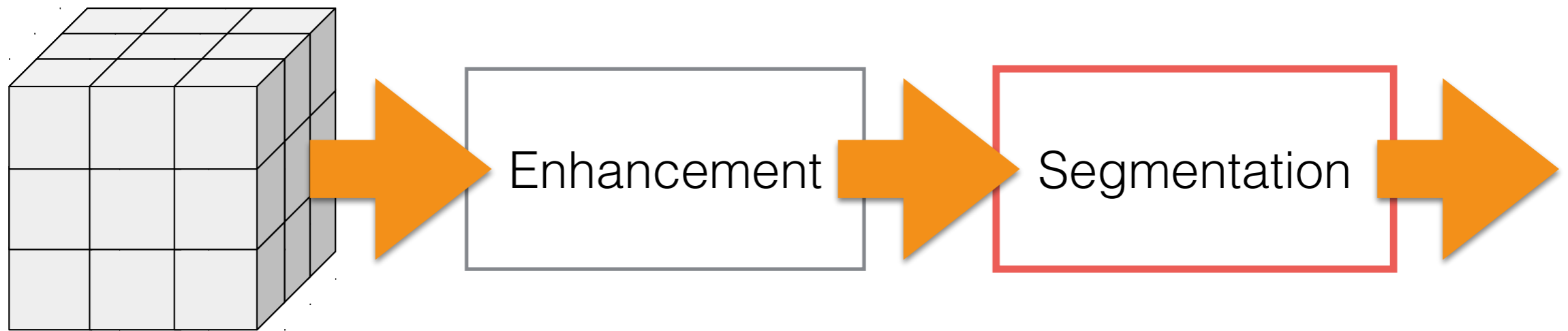


RAW Volume

The Processing Pipeline



The Processing Pipeline



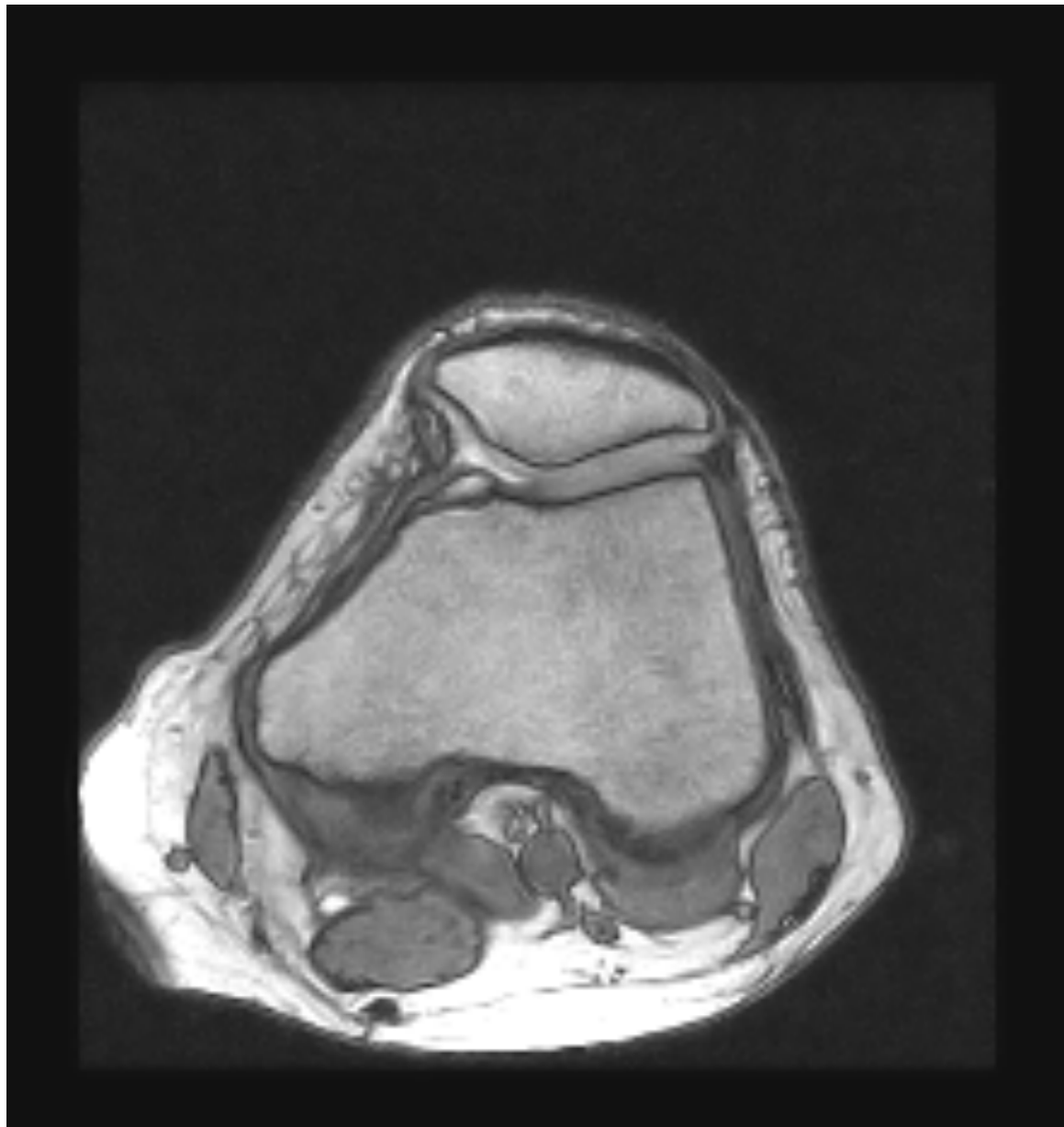
RAW Volume

2D/3D Segmentation

Segmentation

- Segmentation is a process after which we obtain a mask of a structure in an/a image/volume.
- A mask is binary image/volume; i.e., its values can be only either 0 or 1.
- 1 \longrightarrow the pixel/voxel belongs to a structure of our interest
- 0 \longrightarrow the pixel/voxel does not!

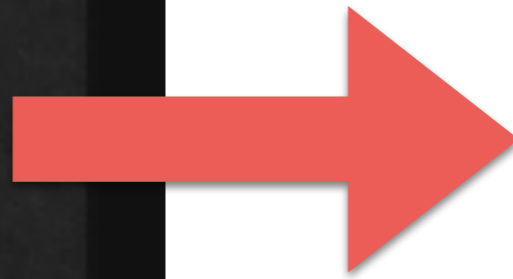
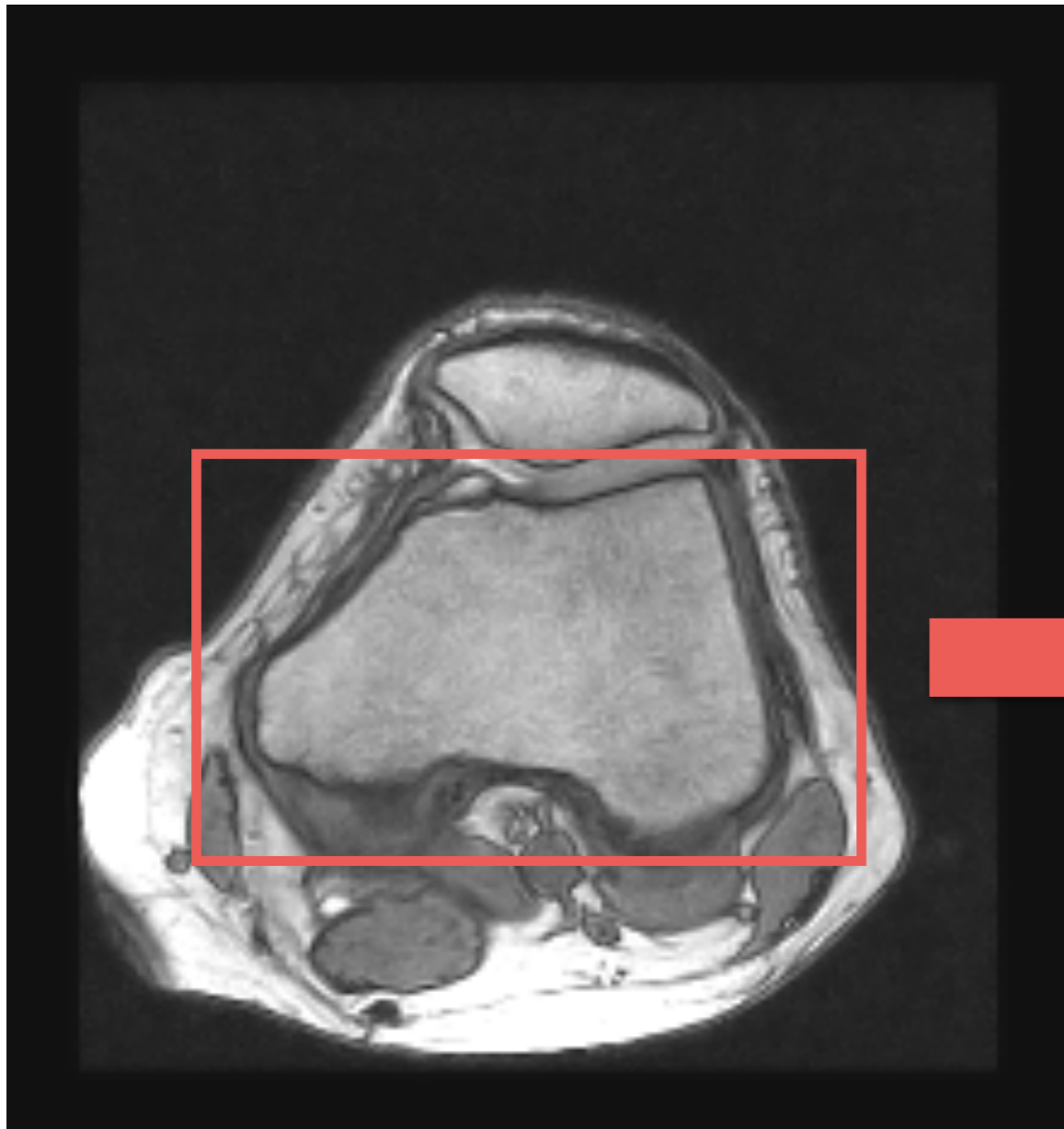
Segmentation Example



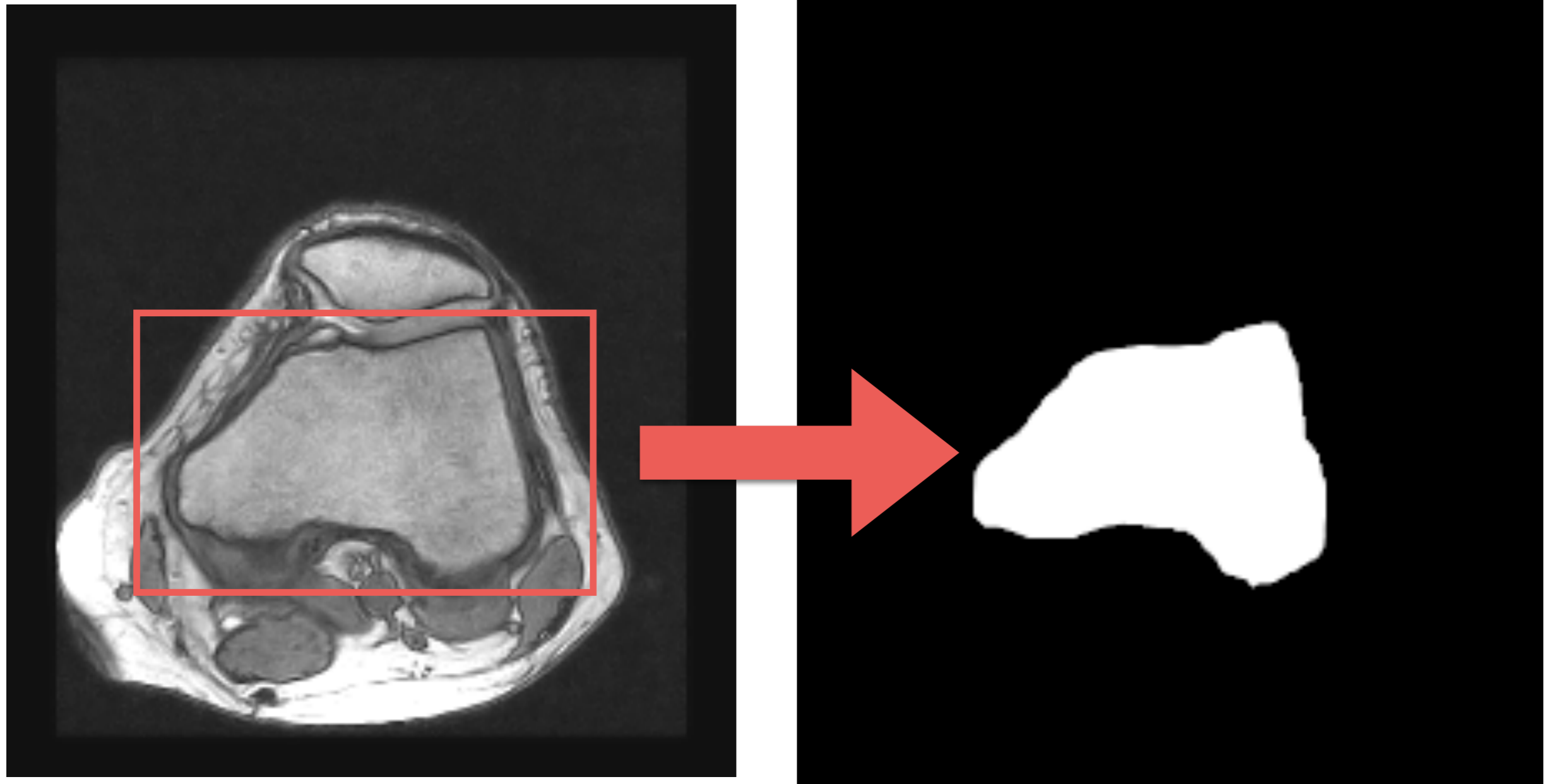
Segmentation Example



Segmentation Example



Segmentation Example



Segmentation

- Obviously, if we need to segment k objects in the image/volume we have two ways to proceed:
 1. We create k -masks, one for each object.
 2. We create an **unsigned integer** mask in which each object is labeled with a number in $[1, k]$. Background is always 0!

3D Segmentation

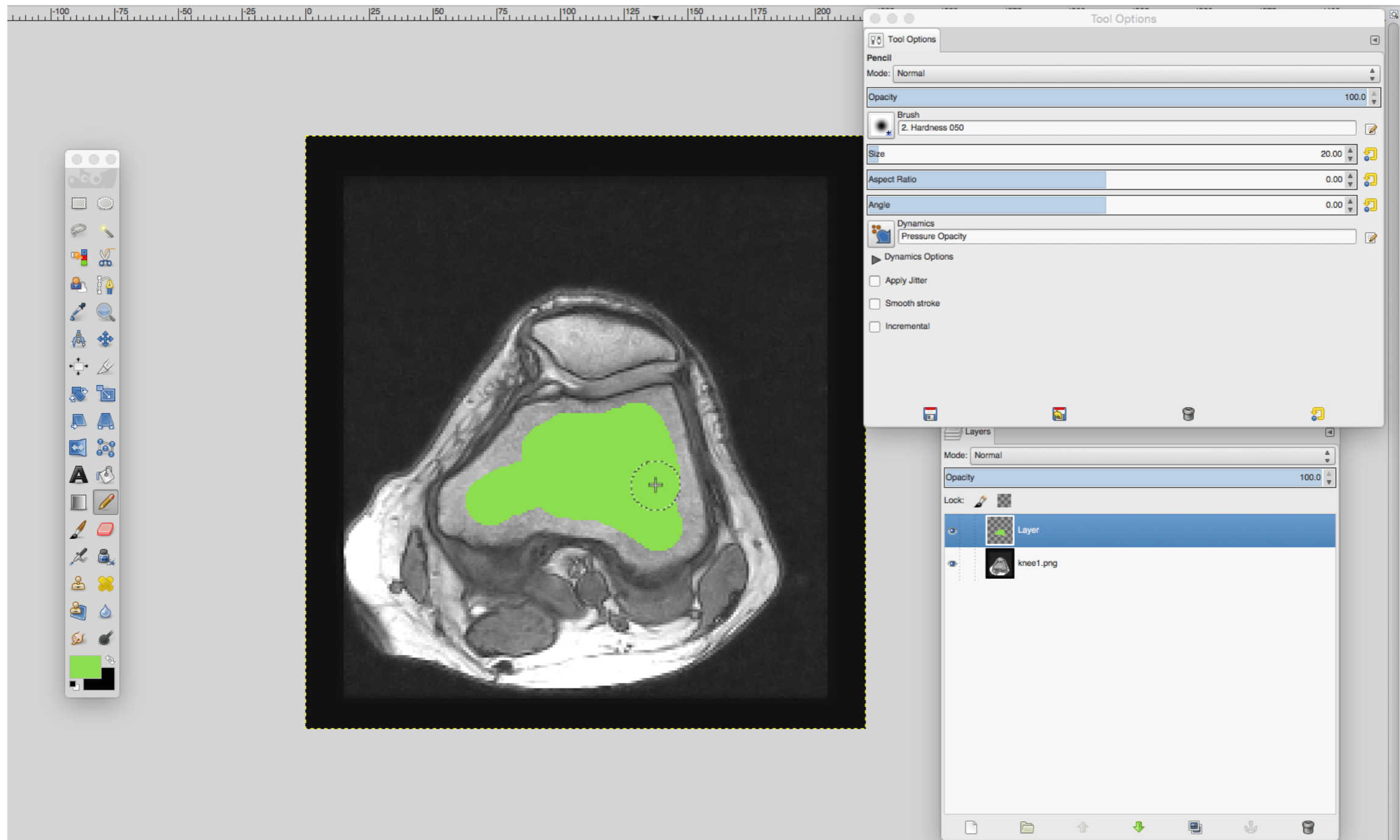
- There are typically two approaches:
 - 2D segmentation for each slice
 - 2D segmentation of a slice and propagation of the segmentation

Manual Segmentation

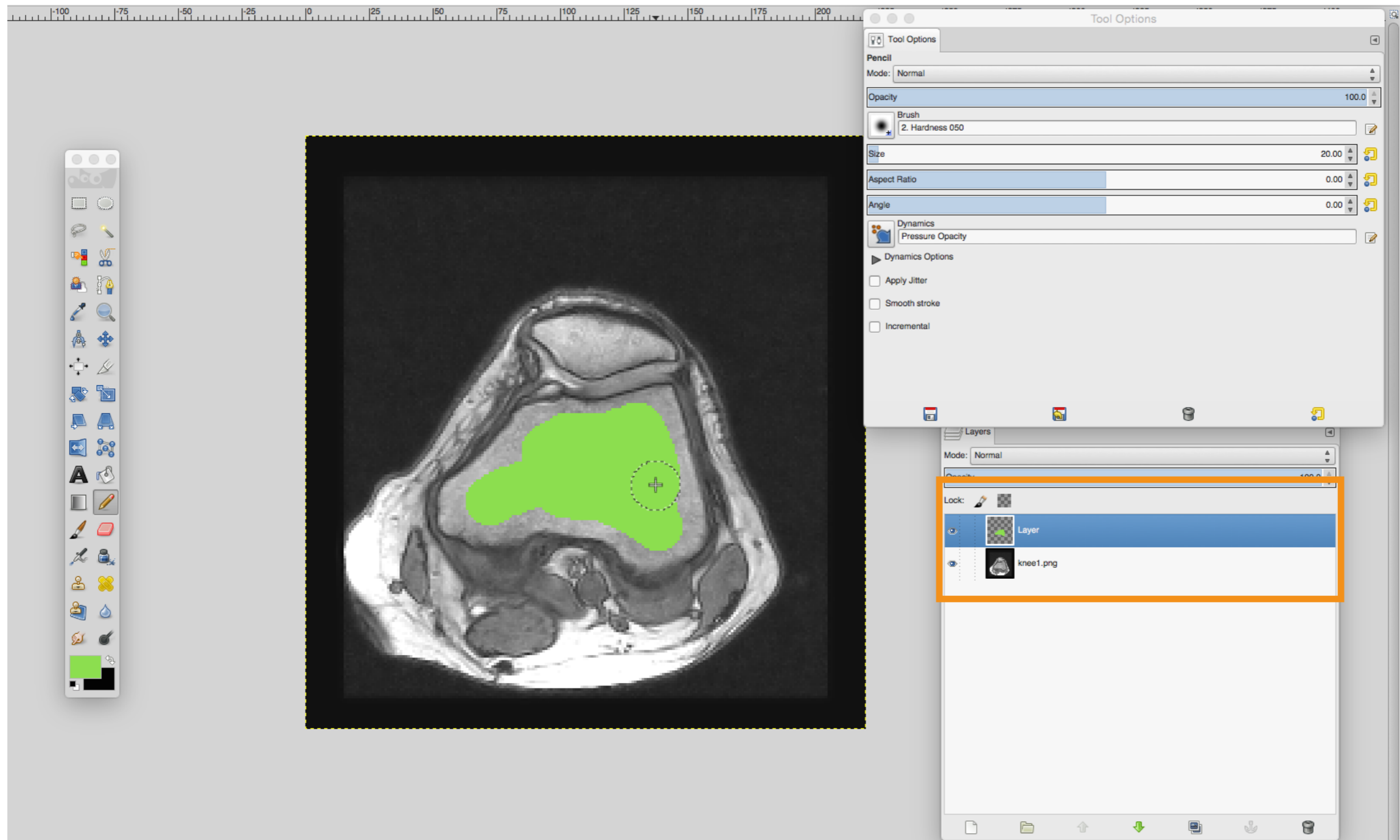
Manual Segmentation: Painting Approach

- We manually paint the mask using a GUI.
- Obviously, the segmentation mask is created in a different layer and not on the input image!

Manual Segmentation: Painting Approach



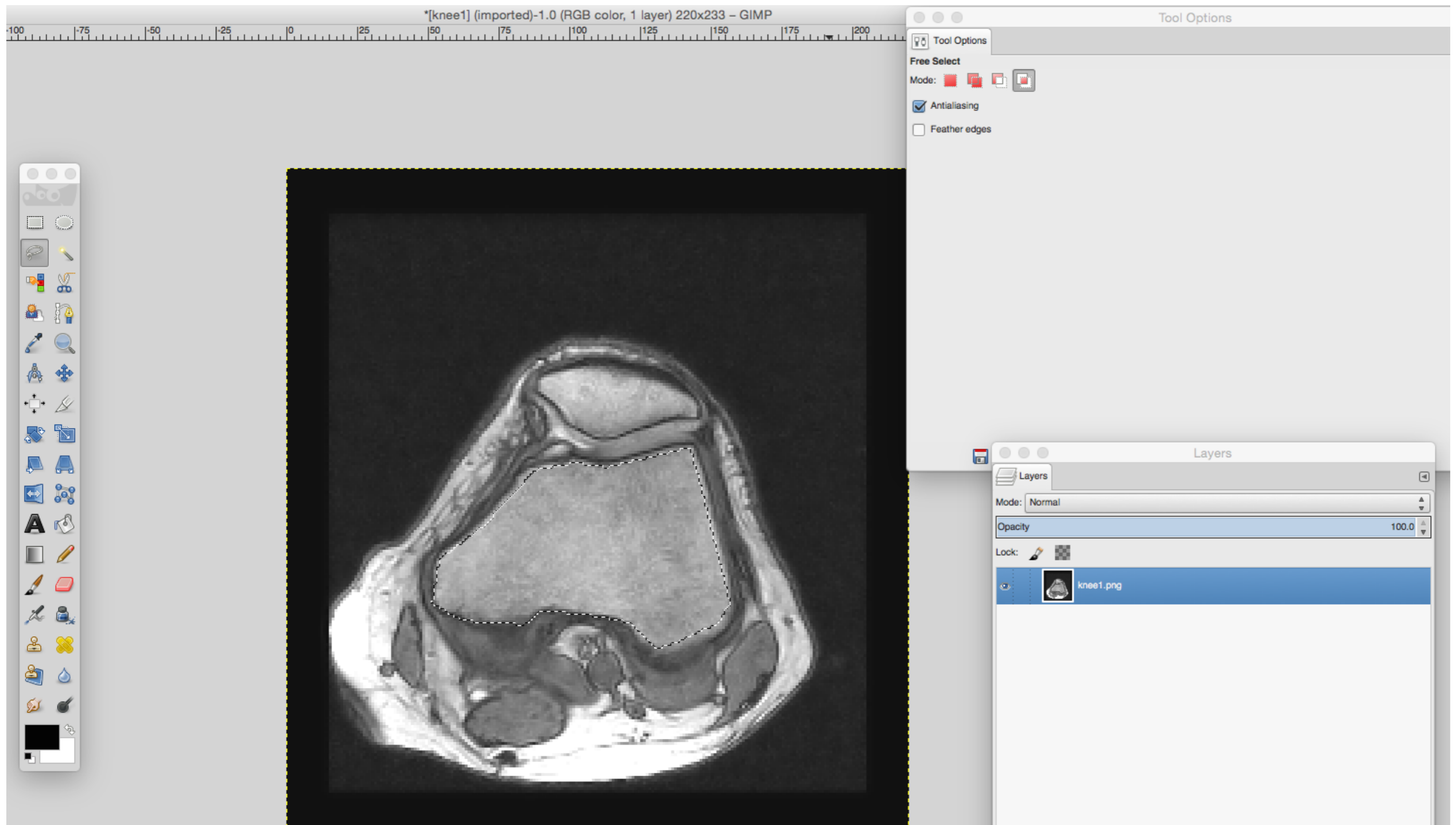
Manual Segmentation: Painting Approach



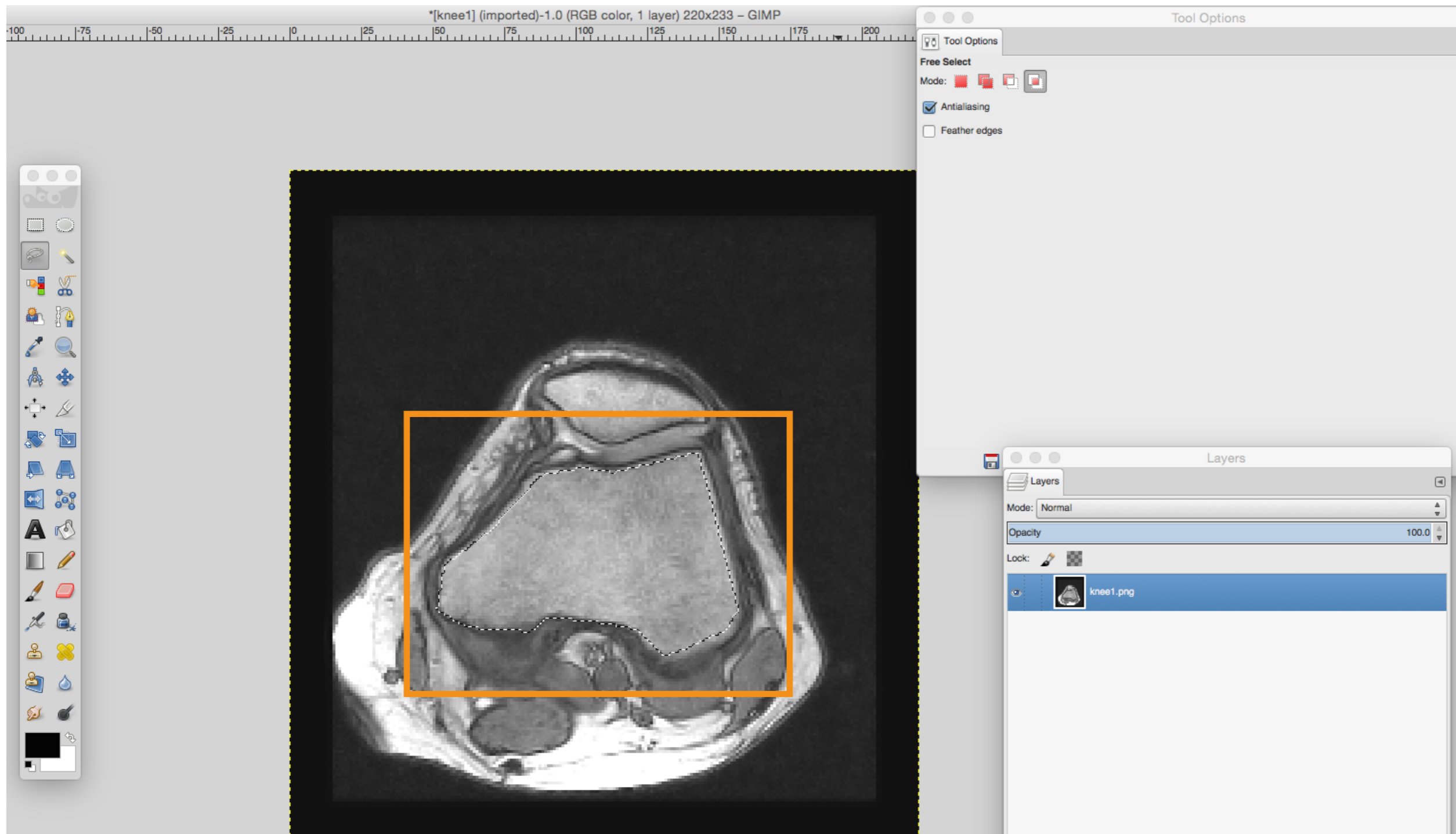
Manual Segmentation: Boundary Definition

- We manually define the mask boundary using a GUI (e.g., GIMP, Adobe PhotoShop, etc.).
- We either define it using polygons or free-hand.
- We can use image gradients and Laplacian to stick polygons to our object of interest.

Manual Segmentation: Boundary Definition



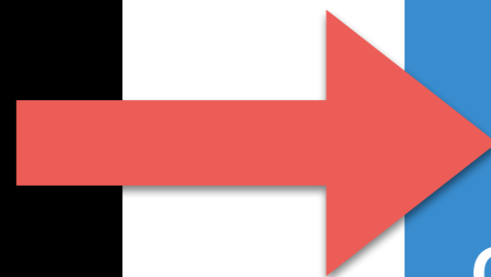
Manual Segmentation: Boundary Definition



Thresholding

Thresholding Example

- We assume that each object in an image/volume has a unique intensity value



Object	Value
Skull	255
Grey Matter	153
Veins	77

Thresholding

- This means:

$$M(i, j) = \begin{cases} 1 & \text{if } d(I(i, j), I_t) < t \\ 0 & \text{otherwise} \end{cases}$$

- We can have different distance functions:

$$d(x, y) = |x - y|$$

$$d(x, y) = (x - y)^2$$

$$d(x, y, \sigma) = \exp\left(-\frac{(x - y)^2}{2\sigma^2}\right)^2$$

Thresholding

- This means:

Reference Value

$$M(i, j) = \begin{cases} 1 & \text{if } d(I(i, j), I_t) < t \\ 0 & \text{otherwise} \end{cases}$$

- We can have different distance functions:

$$d(x, y) = |x - y|$$

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Thresholding

- This means:

$$M(i, j) = \begin{cases} 1 & \text{if } d(I(i, j), I_t) < t \\ 0 & \text{otherwise} \end{cases}$$

Reference Value

Threshold

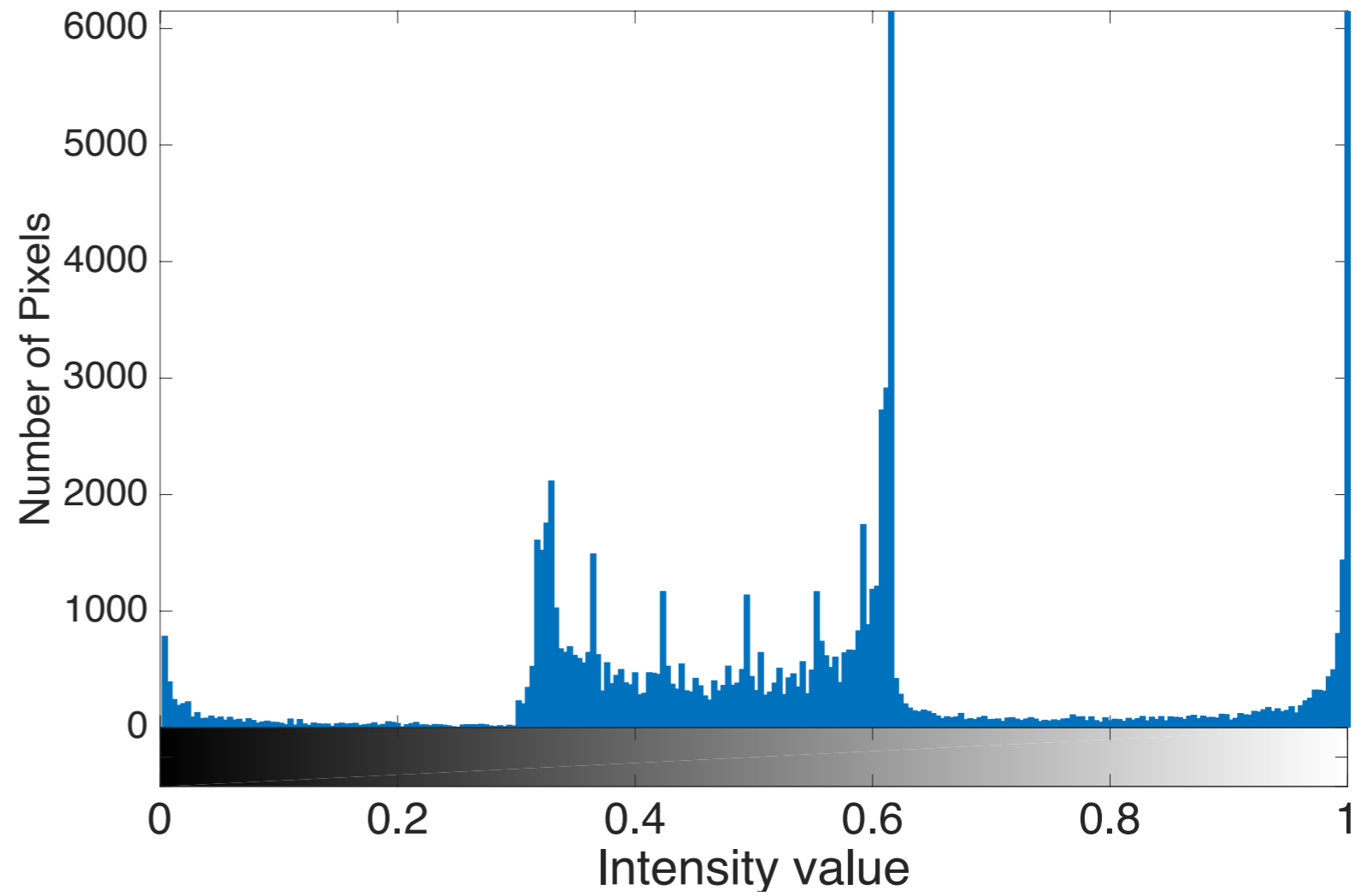
- We can have different distance functions:

$$d(x, y) = |x - y|$$

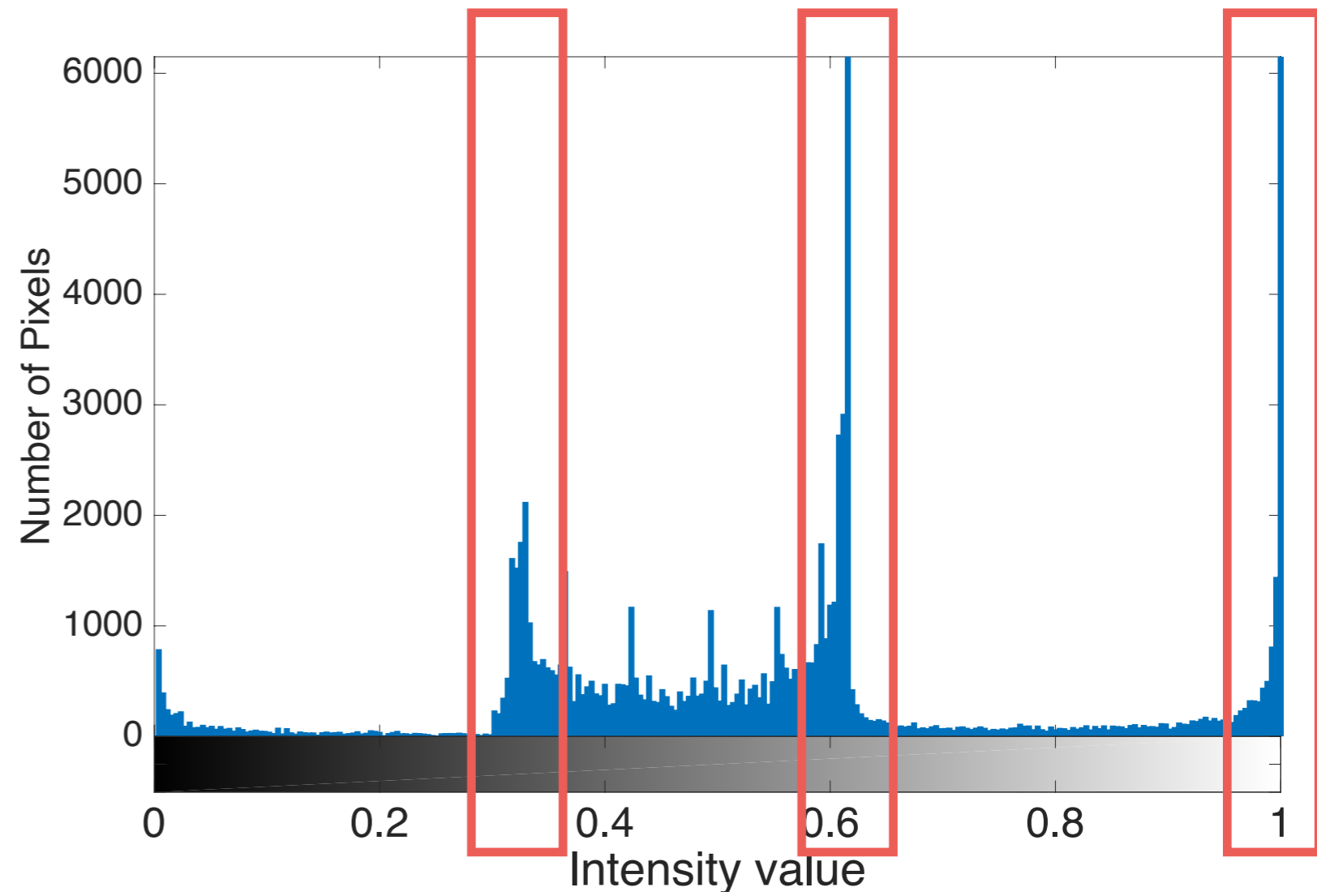
$$d(x, y) = (x - y)^2$$

$$d(x, y, \sigma) = \exp\left(-\frac{(x - y)^2}{2\sigma^2}\right)^2$$

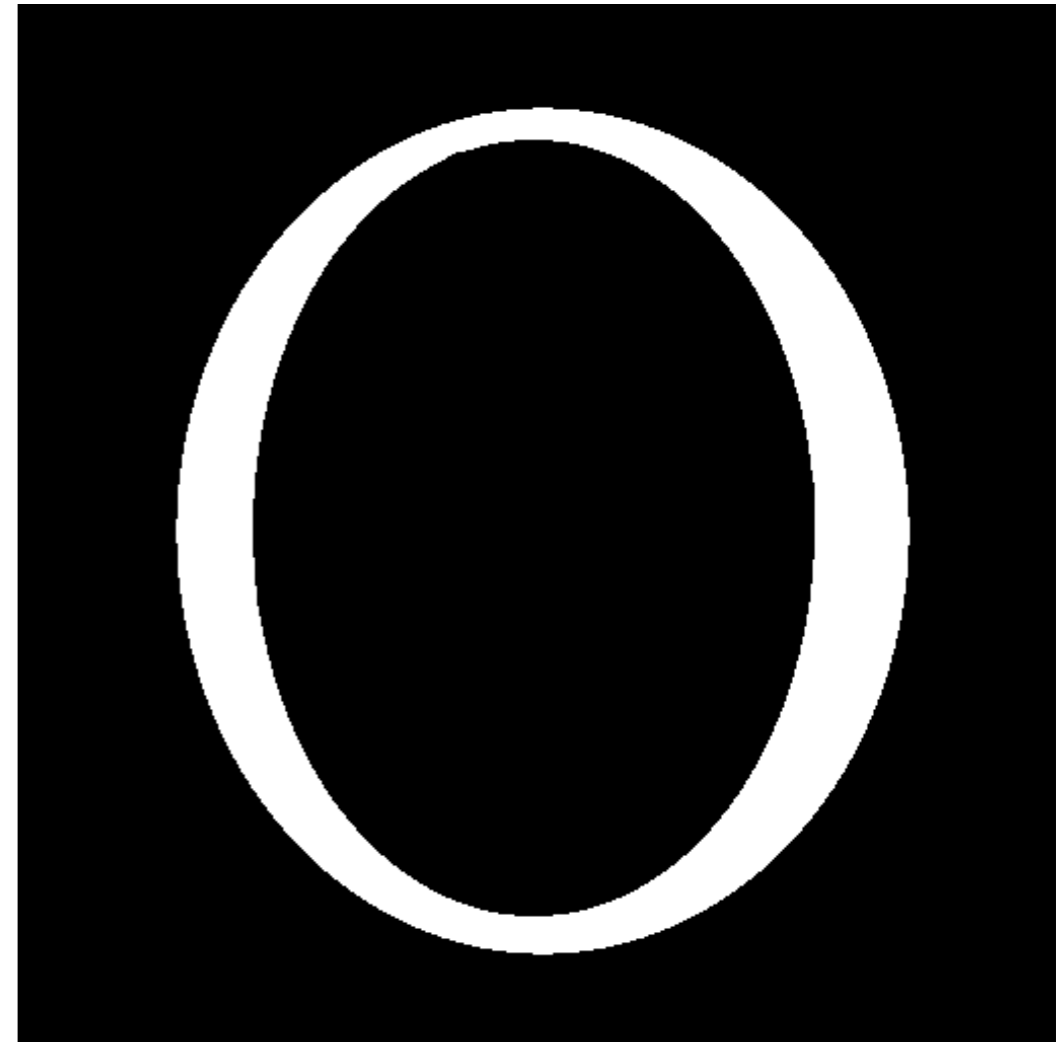
Thresholding



Thresholding

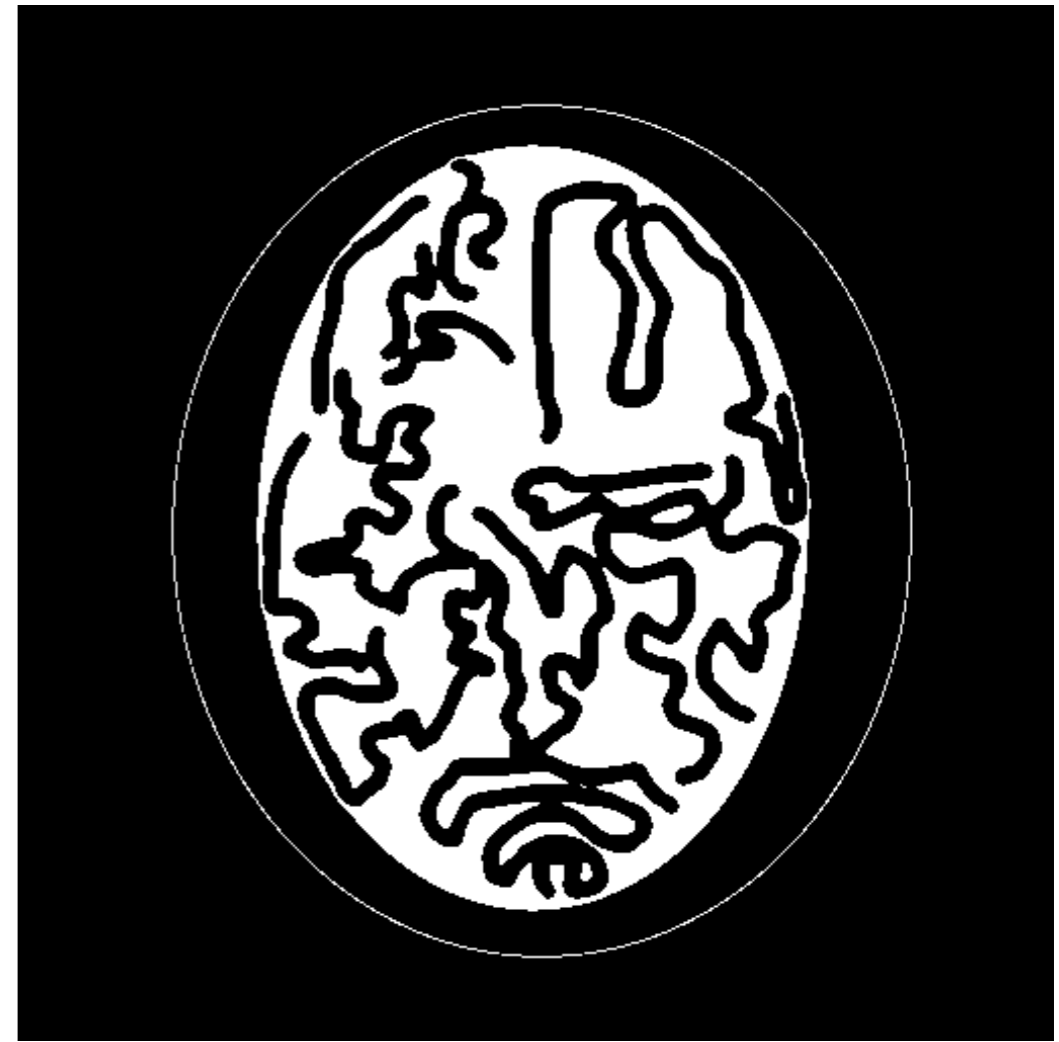


Thresholding Example



$$I_t = 1 \quad t = 0.1$$

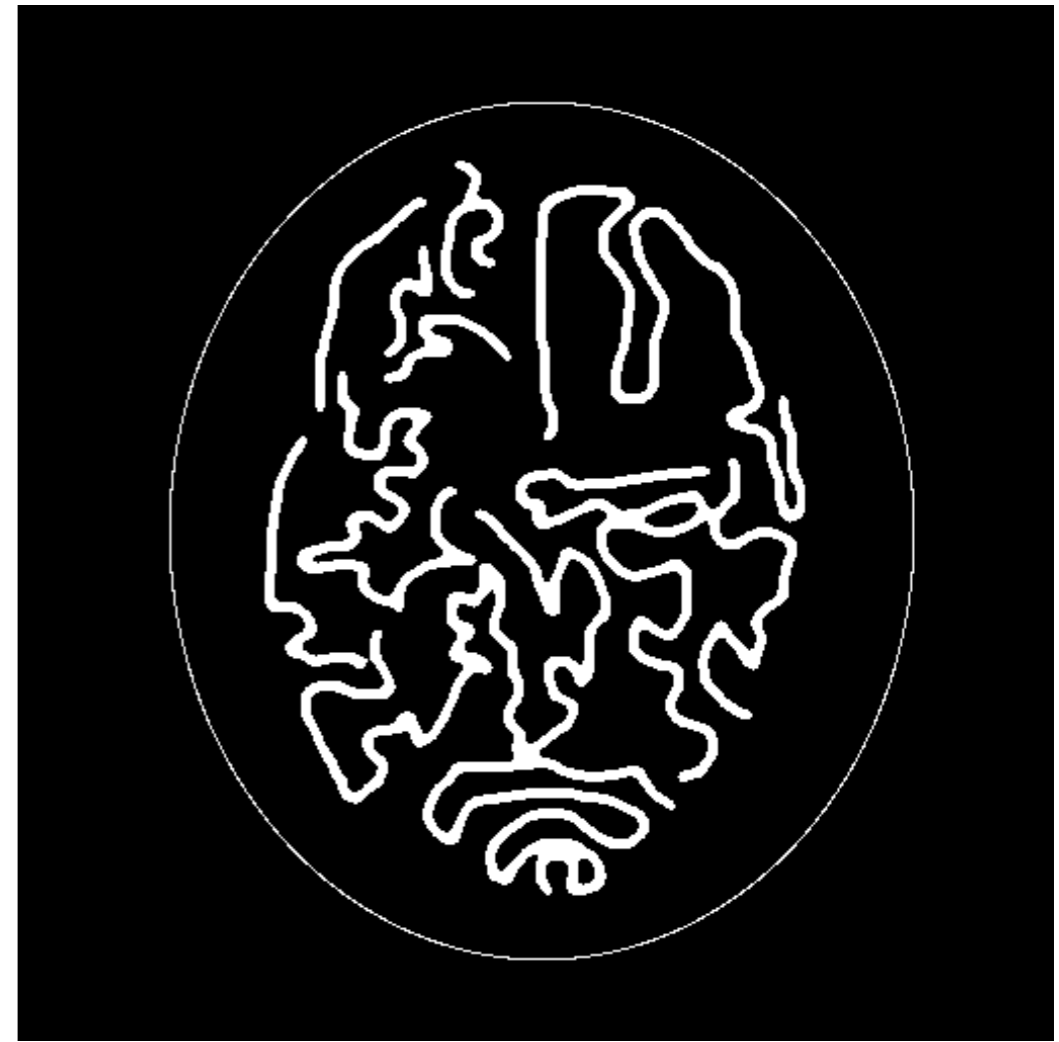
Thresholding Example



$$I_t = 0.6$$

$$t = 0.1$$

Thresholding Example

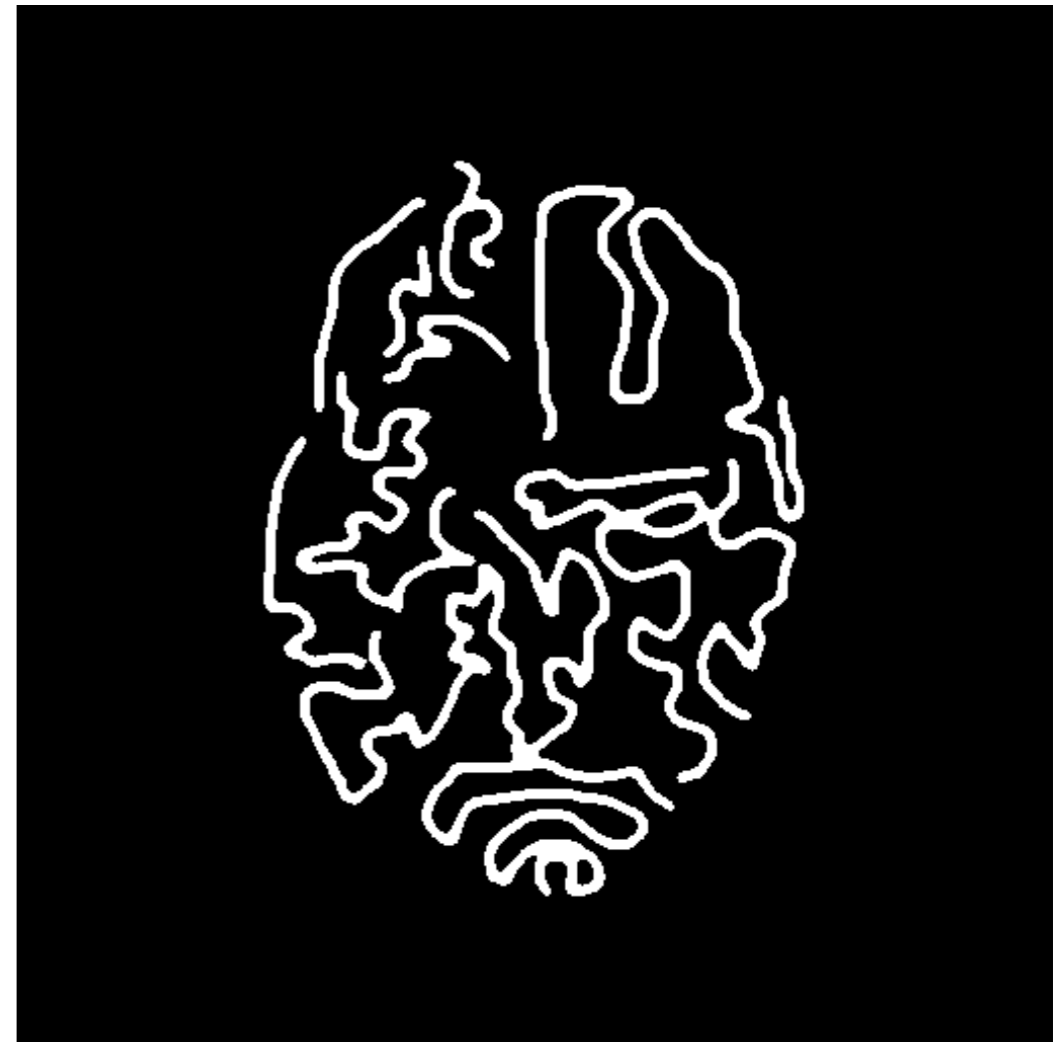


$$I_t = 0.6$$

$$t = 0.1$$

Thresholding: Connected Components

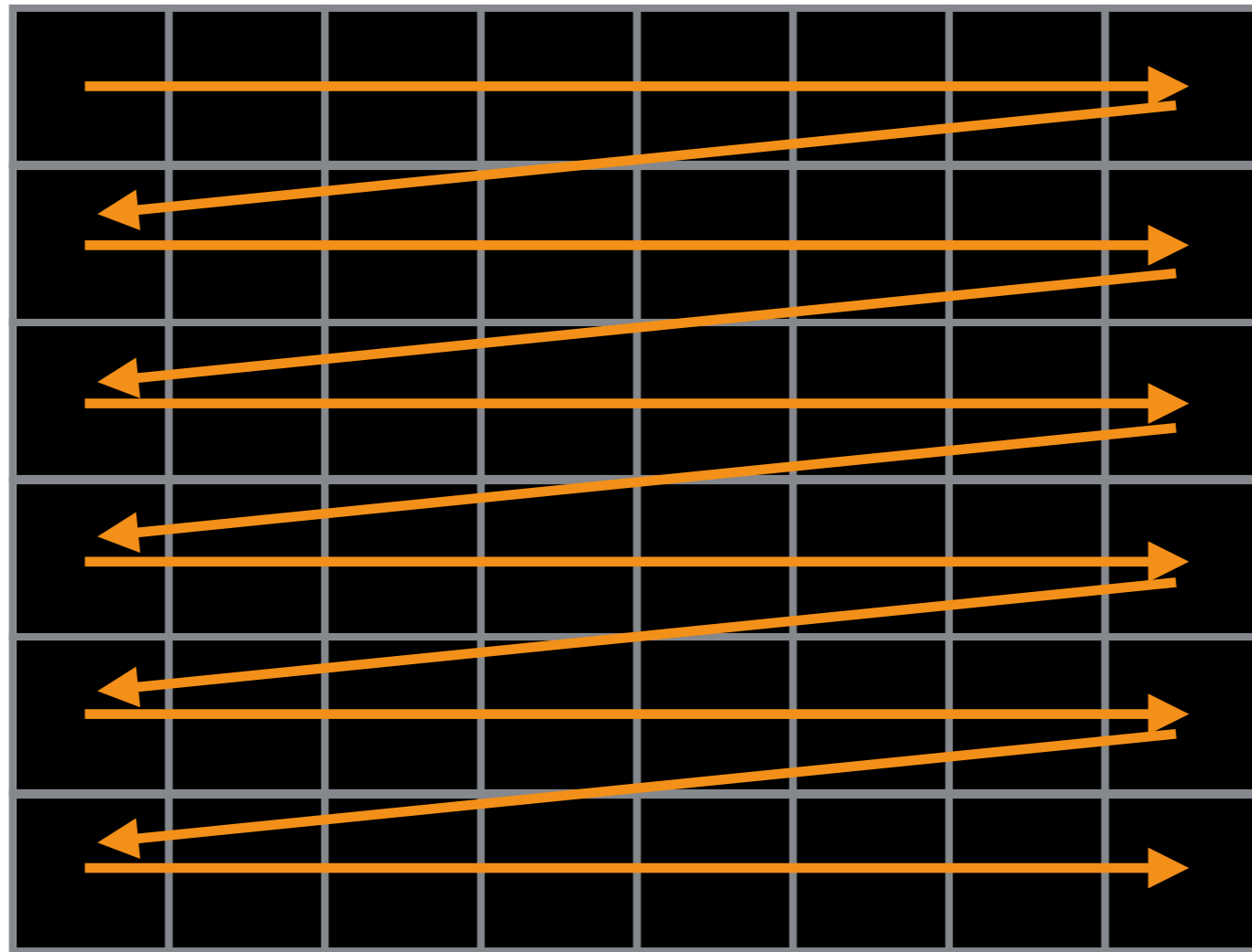
- After segmentation we may end up with different pieces that are not connected.



Thresholding: Connected Components

- A two-pass algorithm that works in scan order (from left to right and from top to bottom).
- 1-Pass: it creates labels to groups of pixel.
- 2-Pass: it merges groups that are connected.

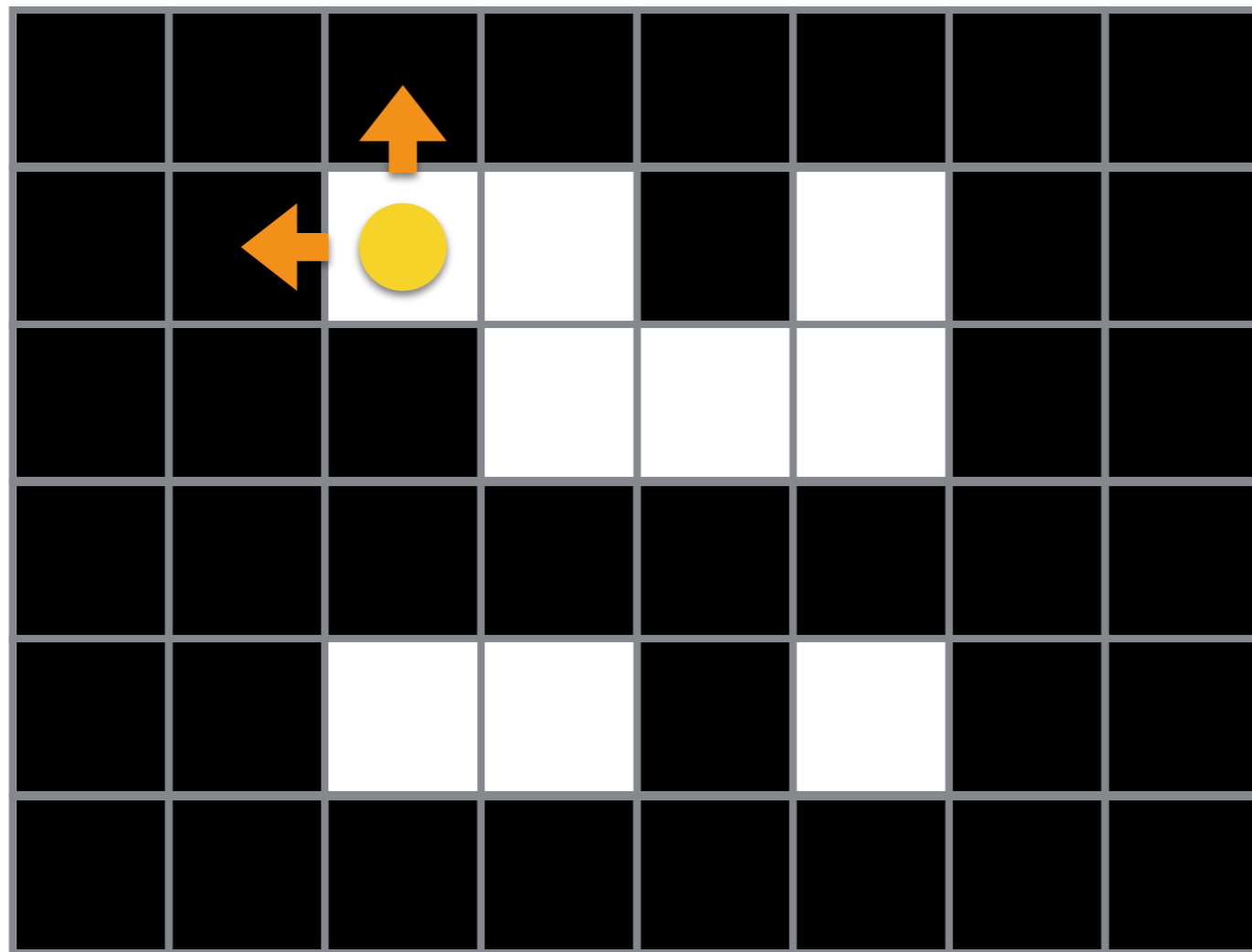
Thresholding: Connected Components



Scan order

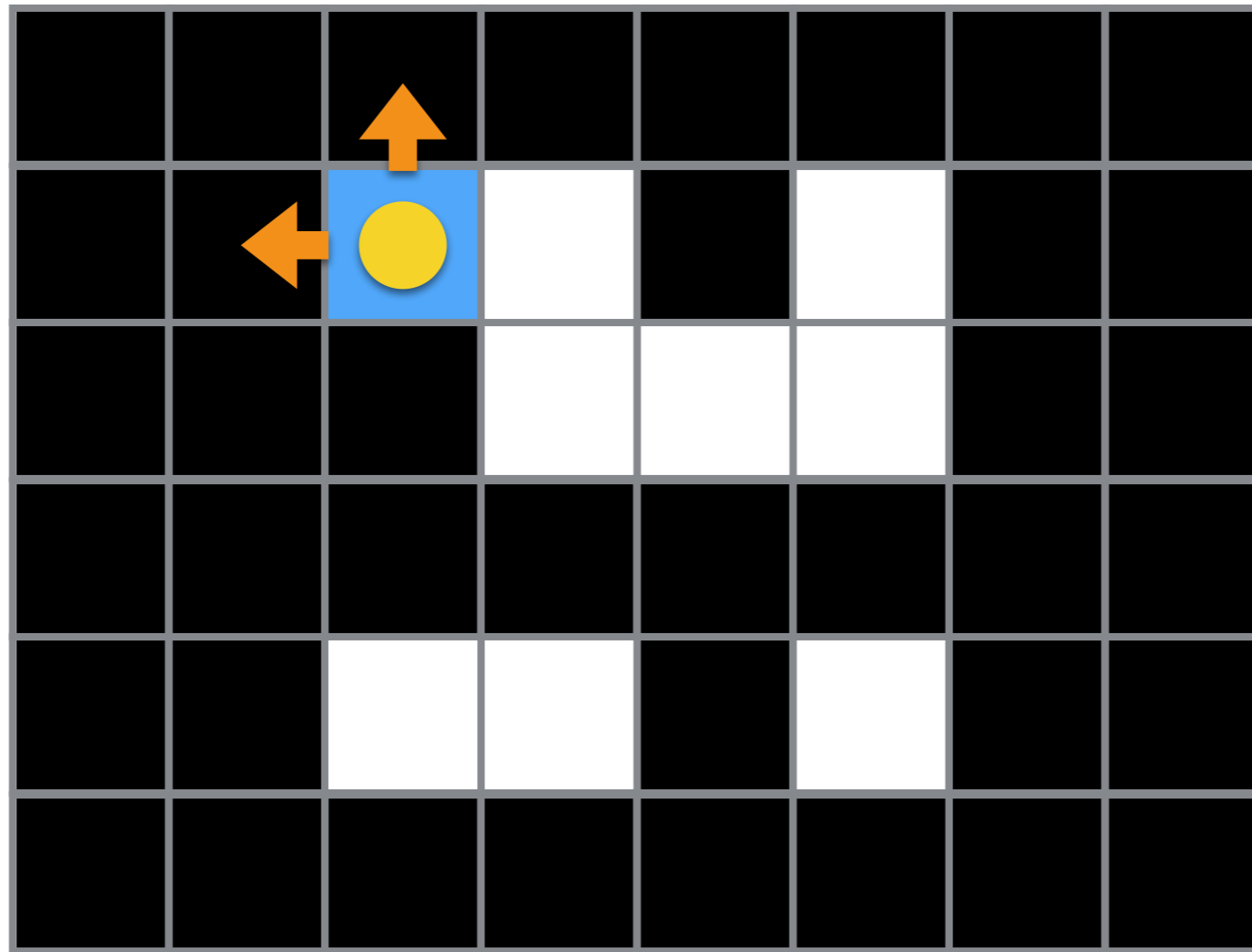
First Pass

Thresholding: Connected Components



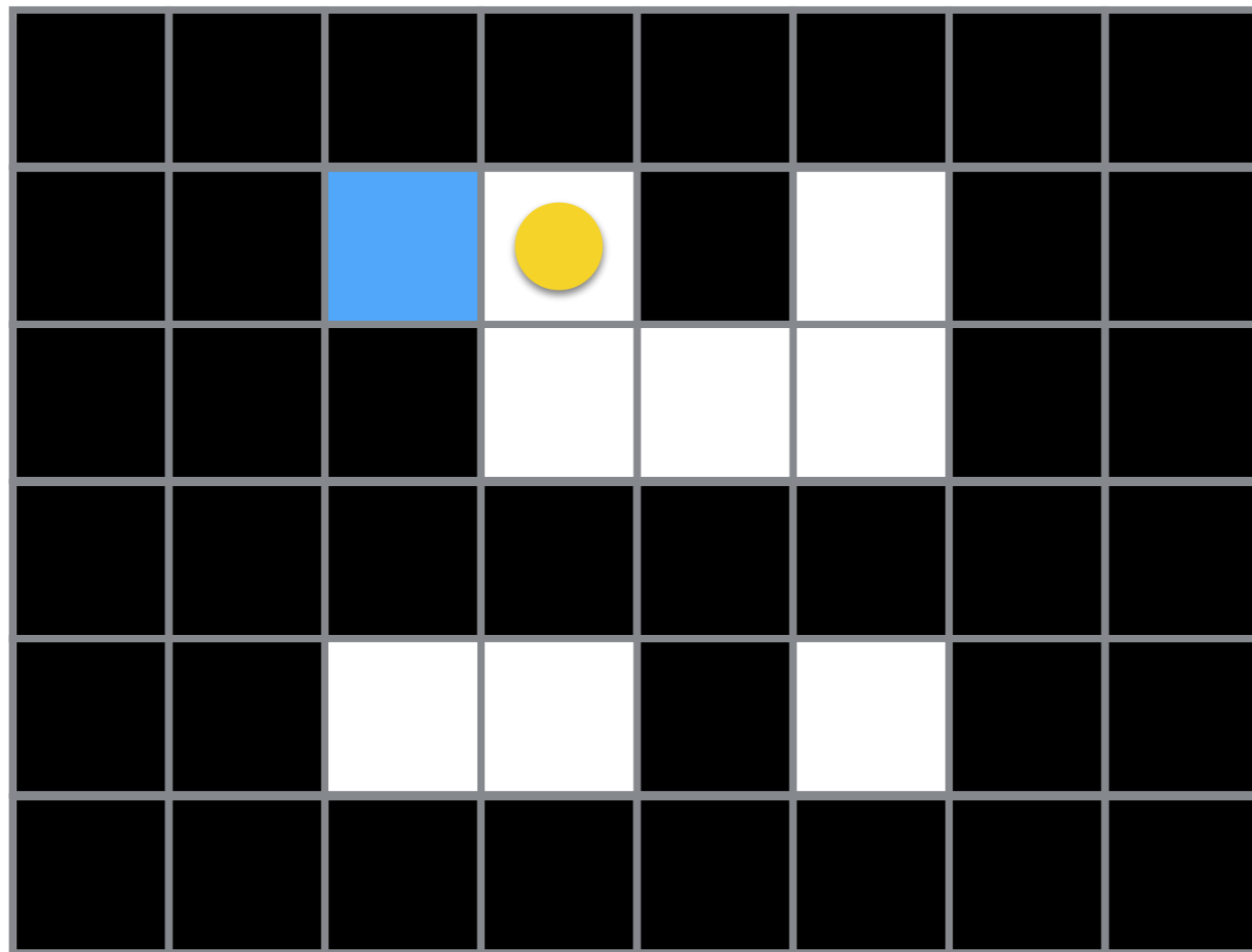
We check up and left neighbors to see if they have a label.

Thresholding: Connected Components



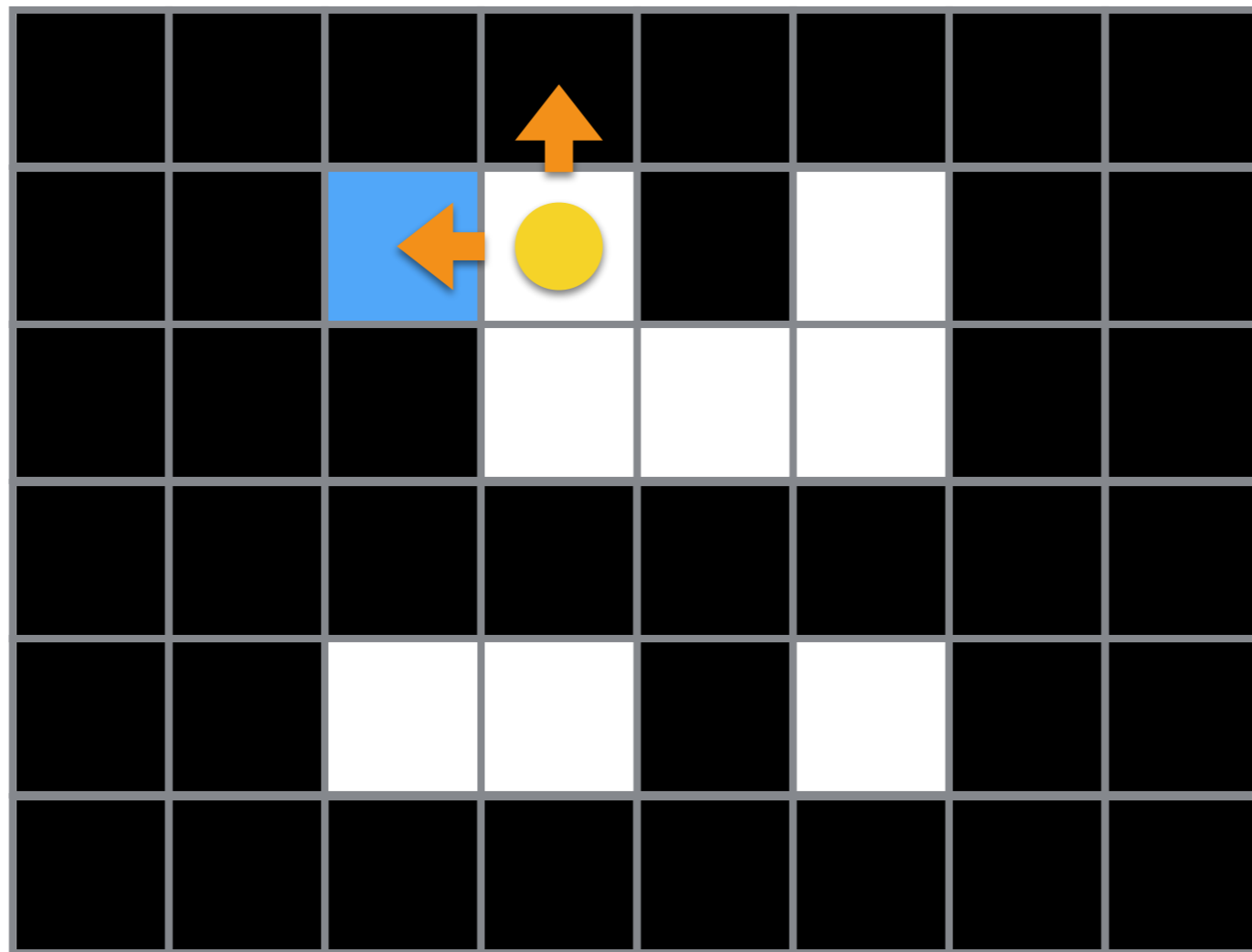
If not we create a
new one.

Thresholding: Connected Components



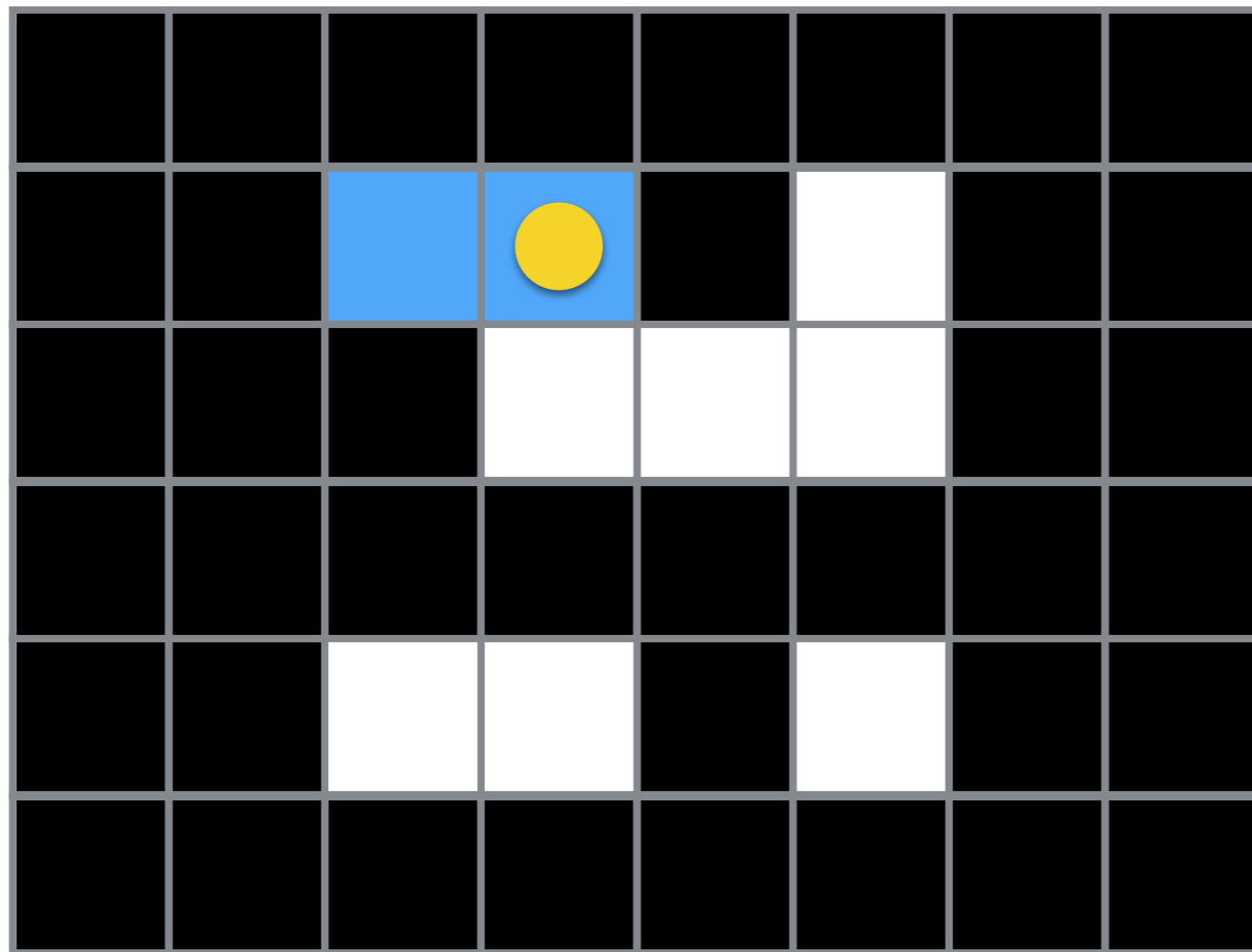
Then, we move
right, and we
repeat the
process.

Thresholding: Connected Components



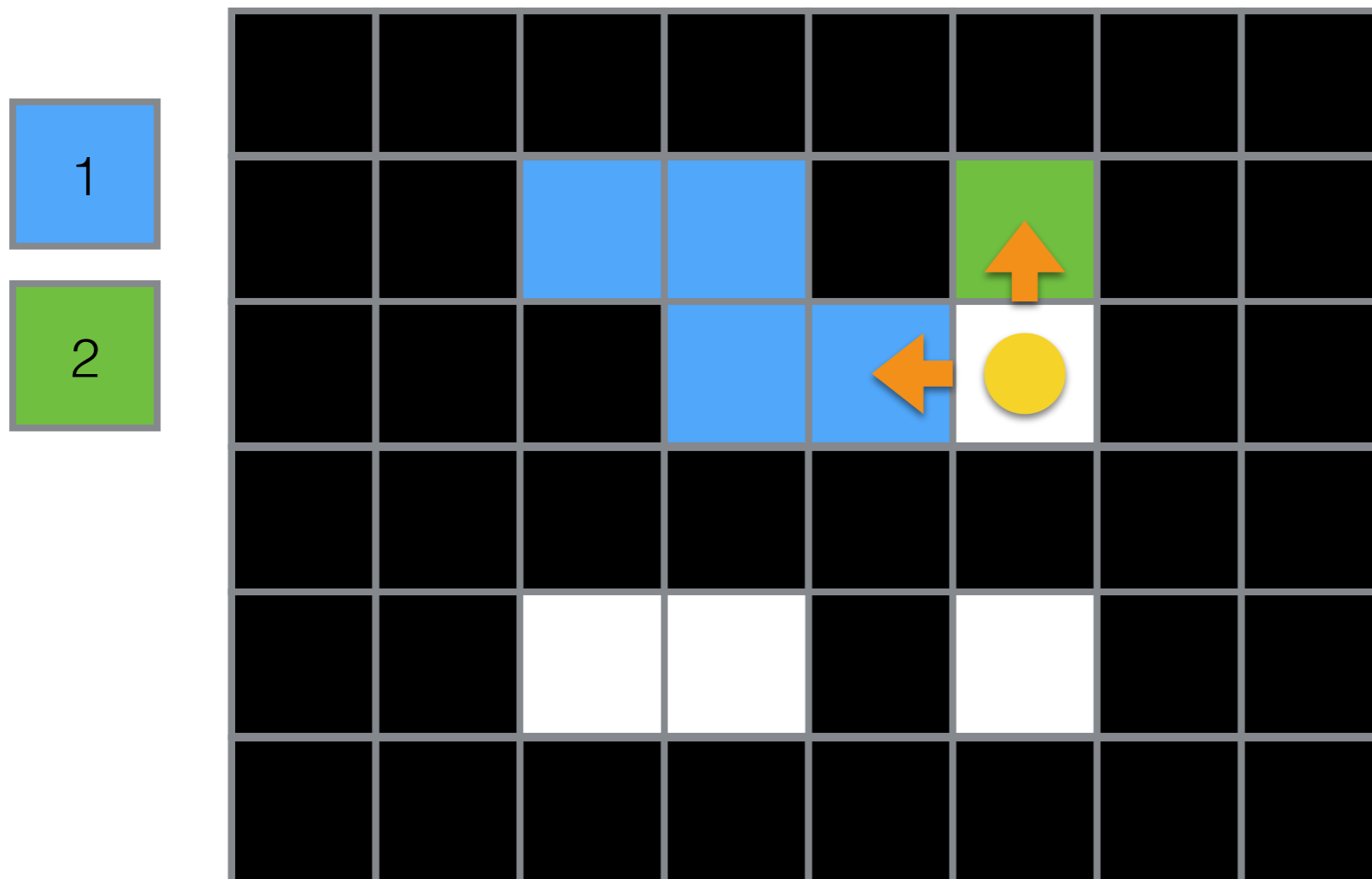
In this case, the left neighbor has a label, so we reuse it.

Thresholding: Connected Components



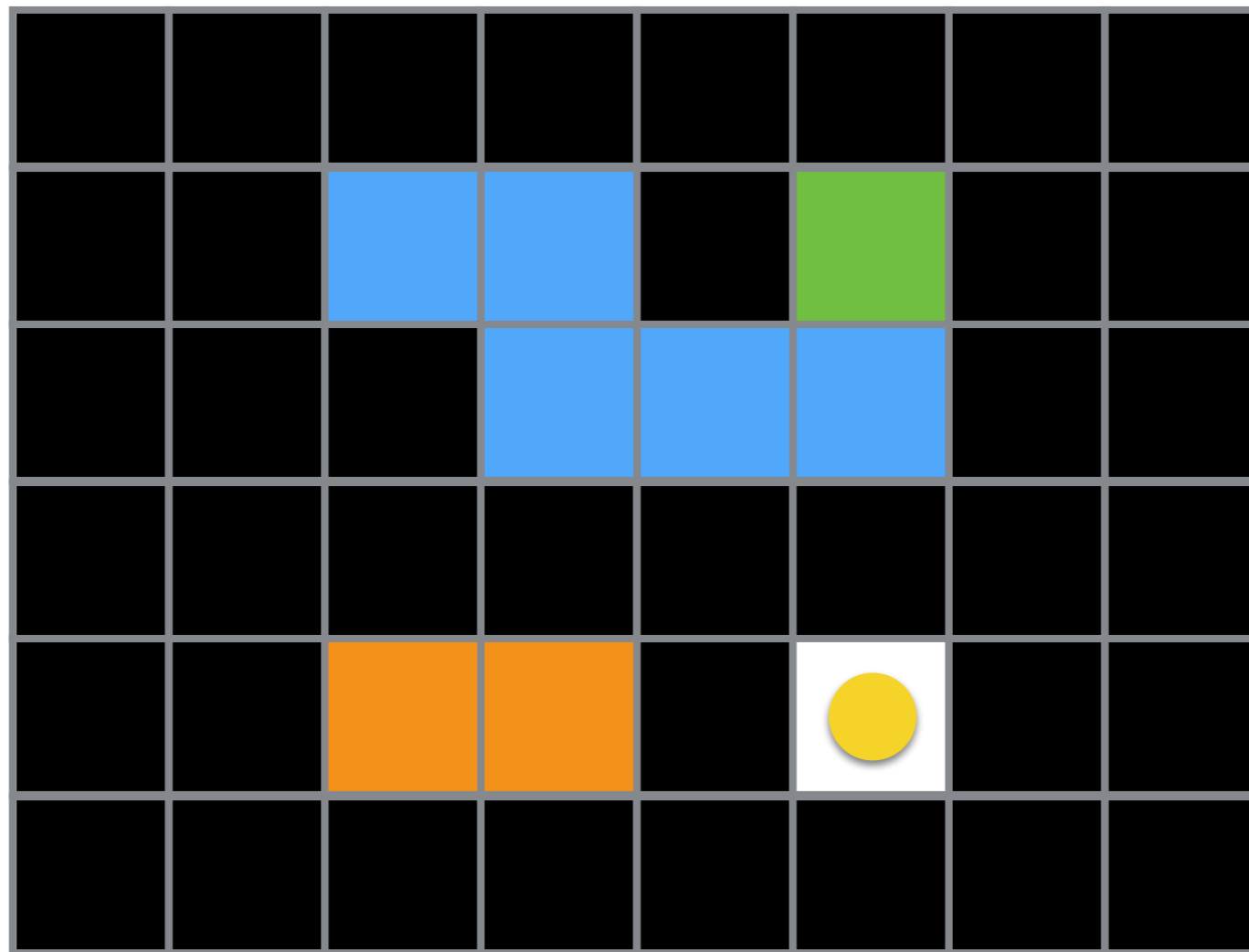
In this case, the left neighbor has a label, so we reuse it.

Thresholding: Connected Components

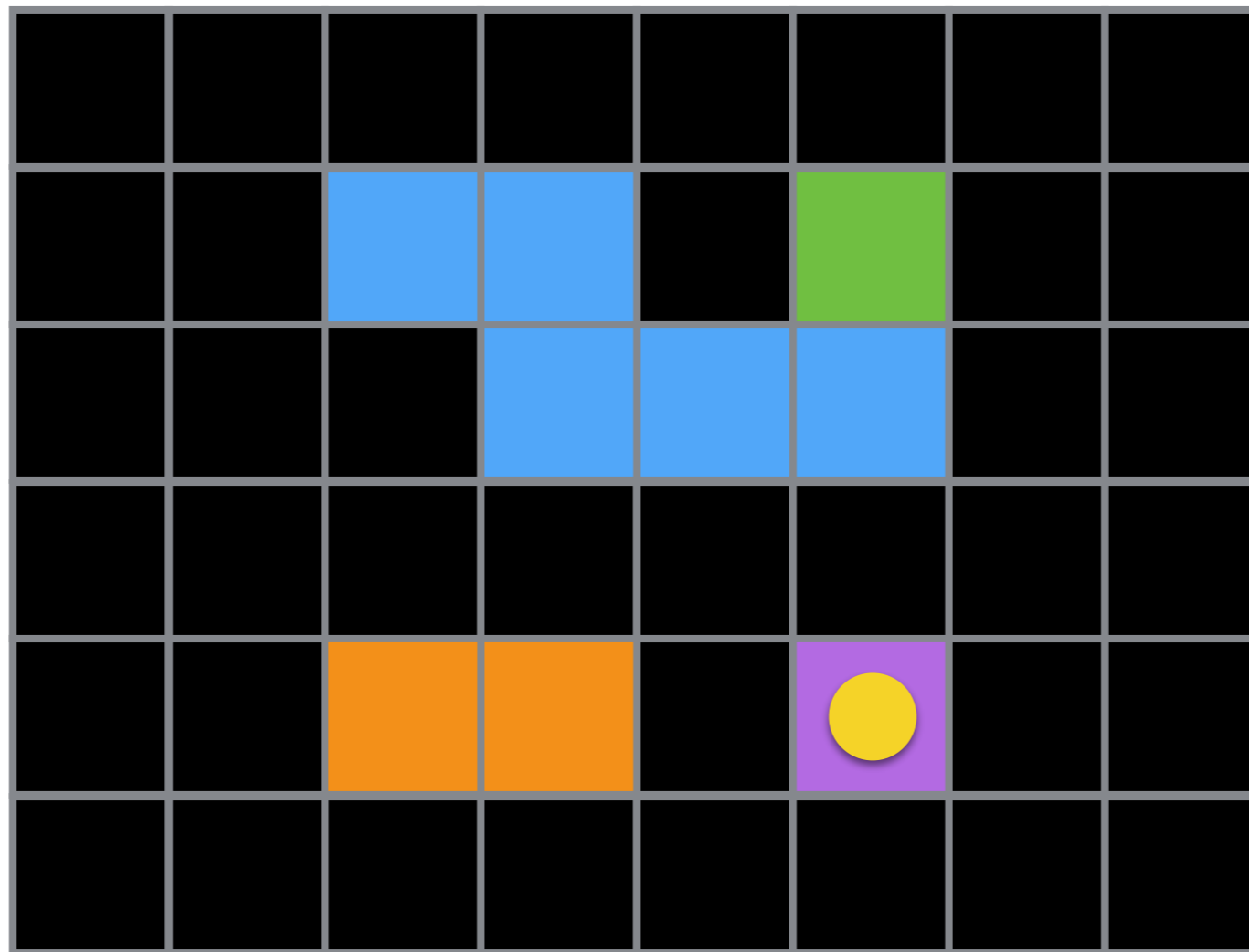


In this case, we choose the lowest label, and we store that 1 is equivalent to 2

Thresholding: Connected Components

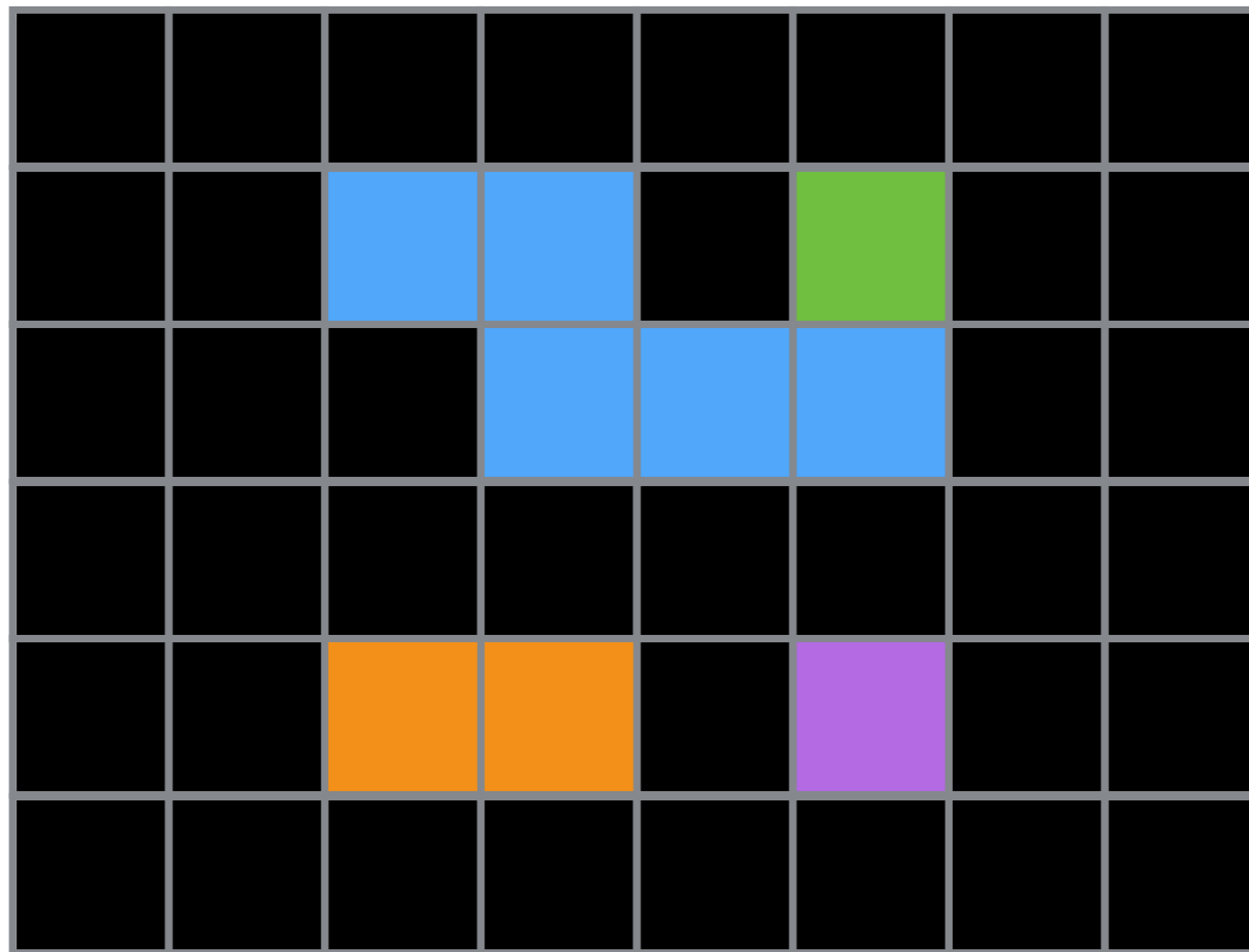
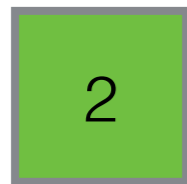
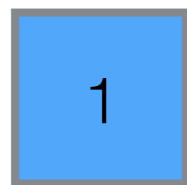


Thresholding: Connected Components



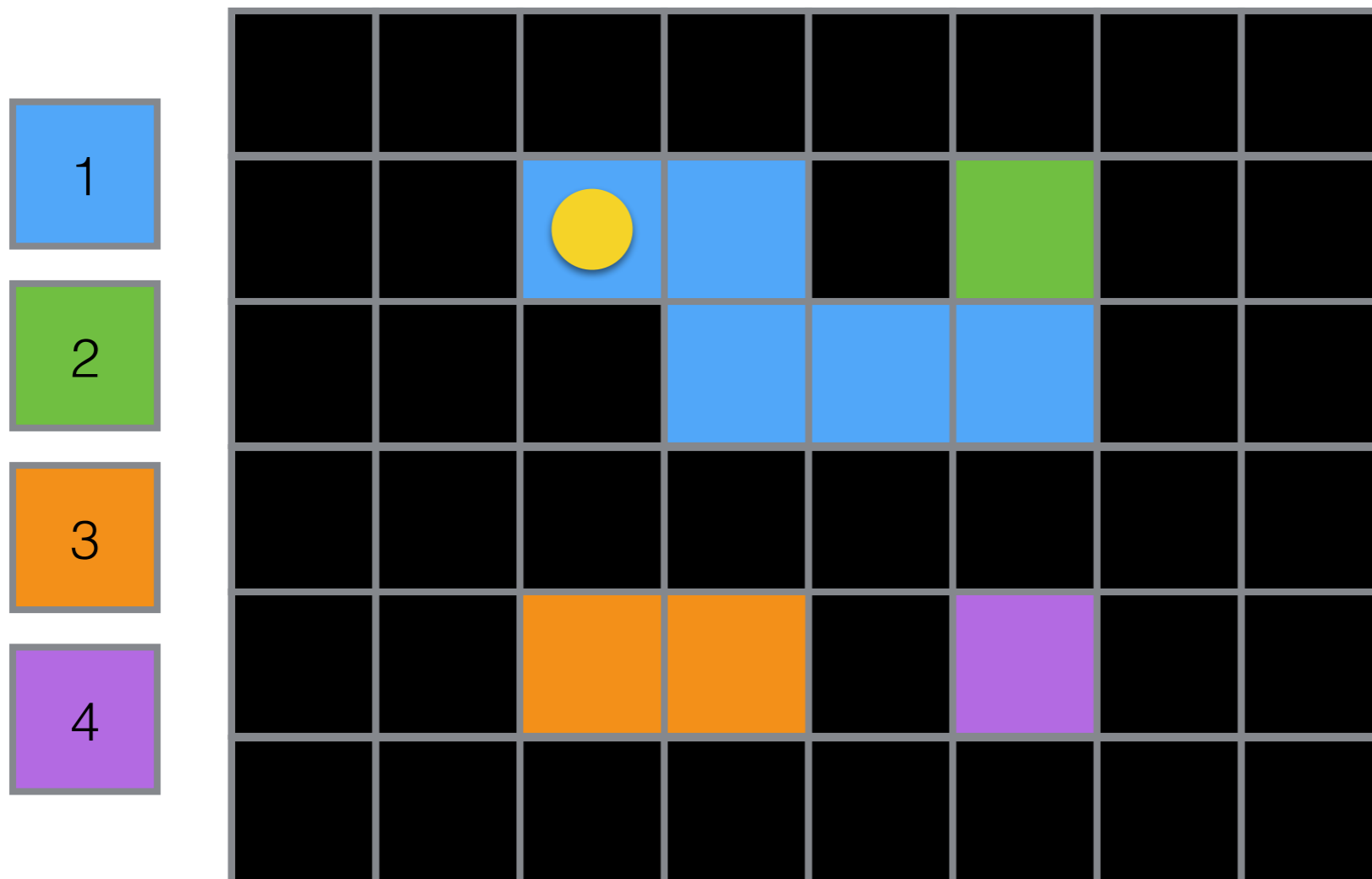
Second Pass

Thresholding: Connected Components

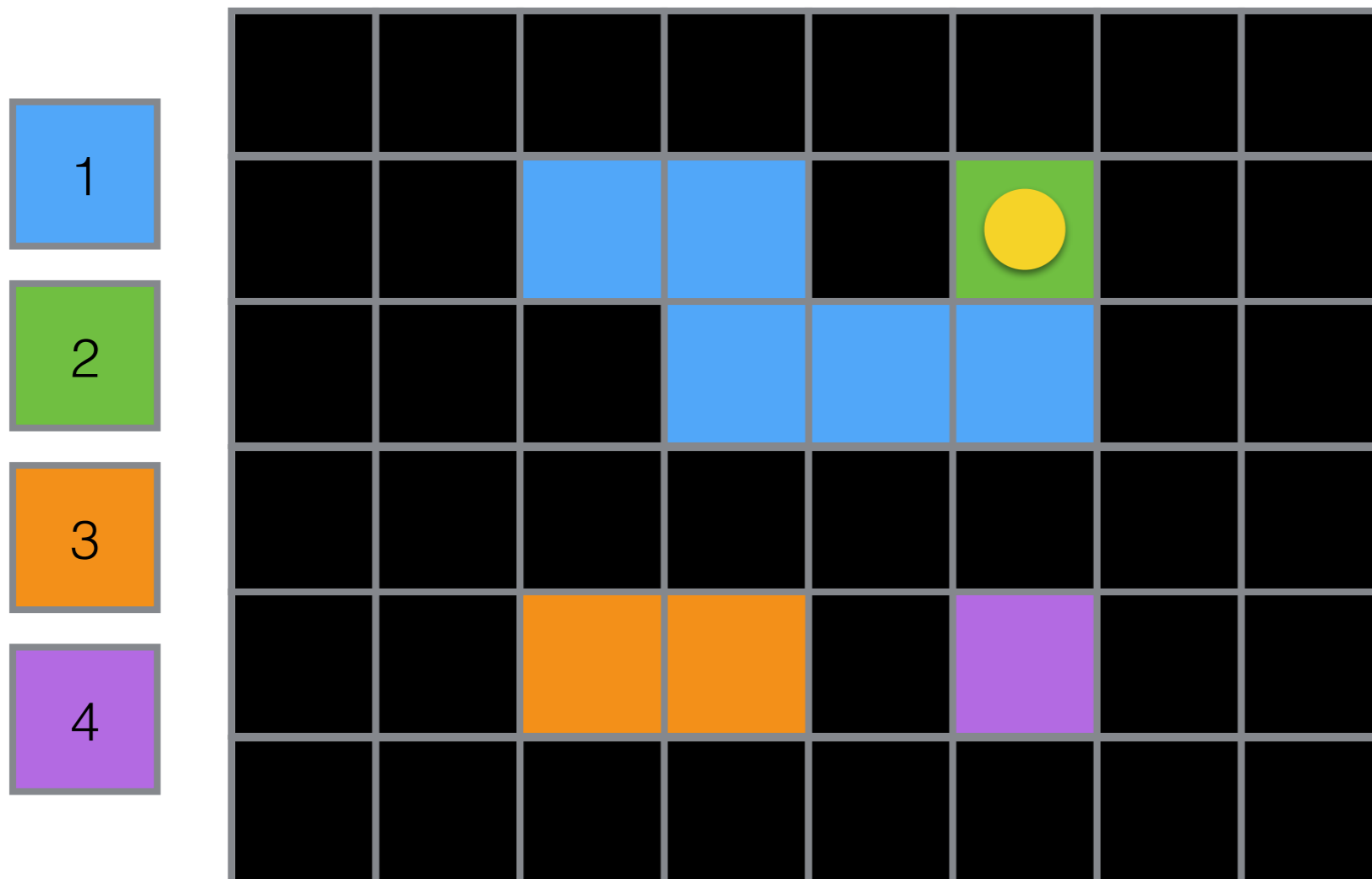


We go through all pixels. For each pixel we set the value of lowest equivalent.

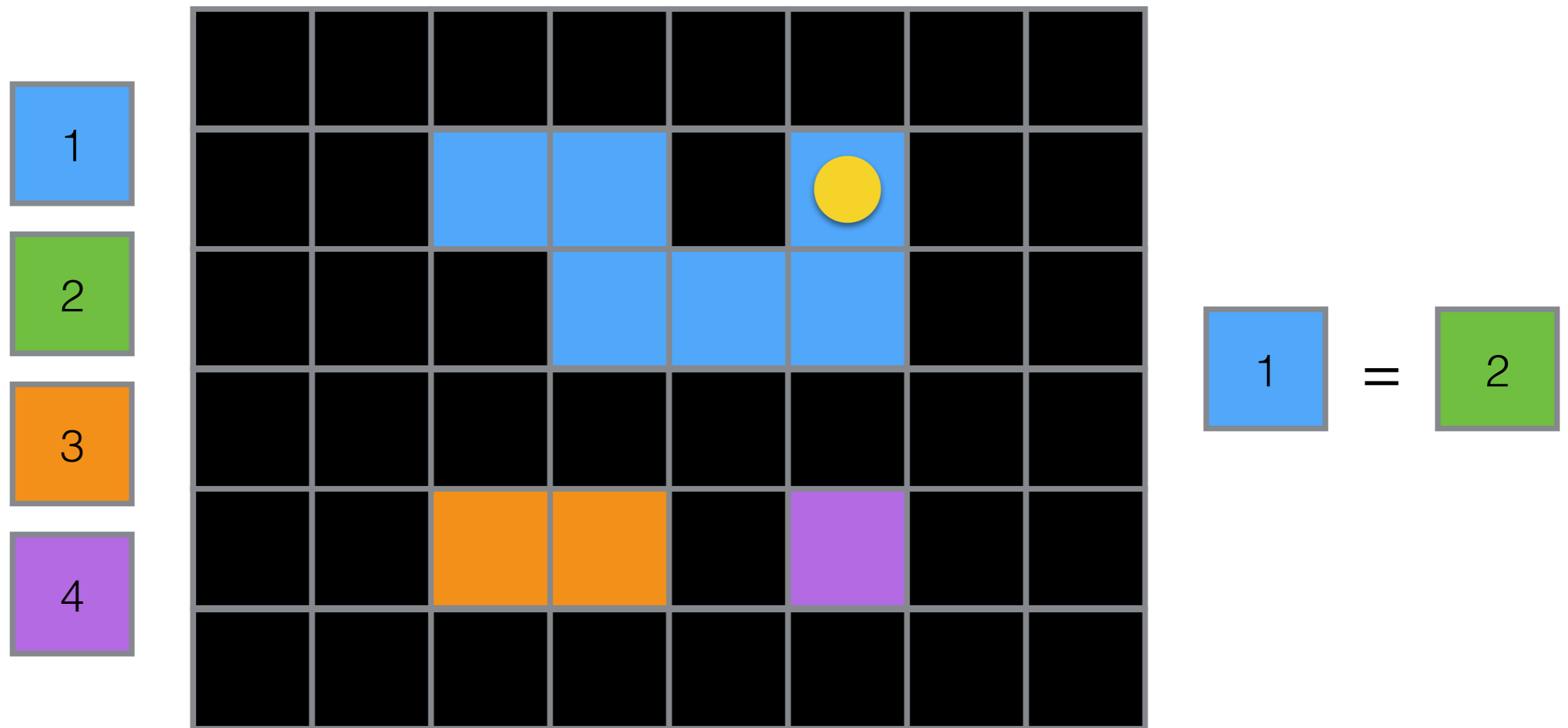
Thresholding: Connected Components



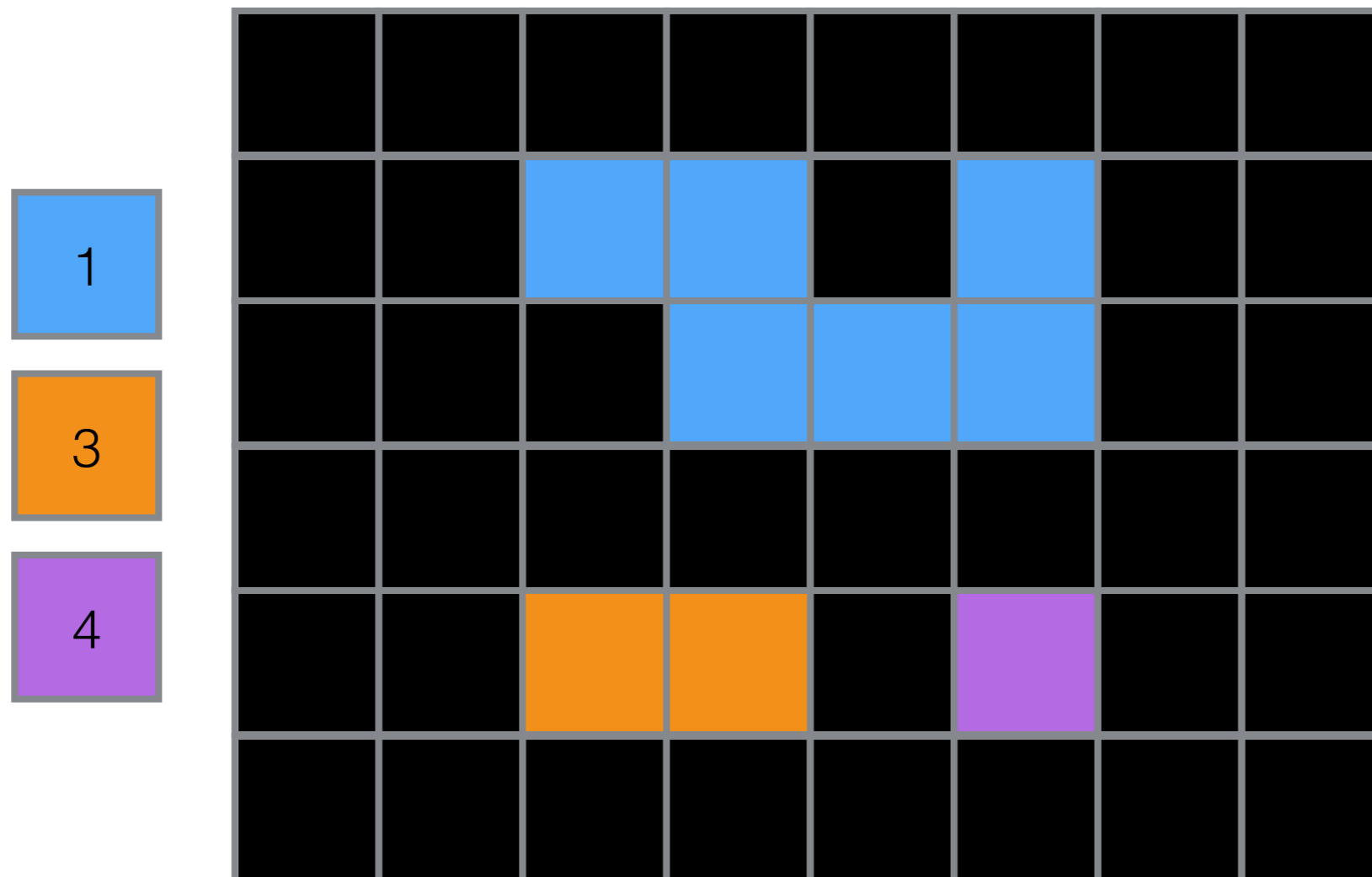
Thresholding: Connected Components



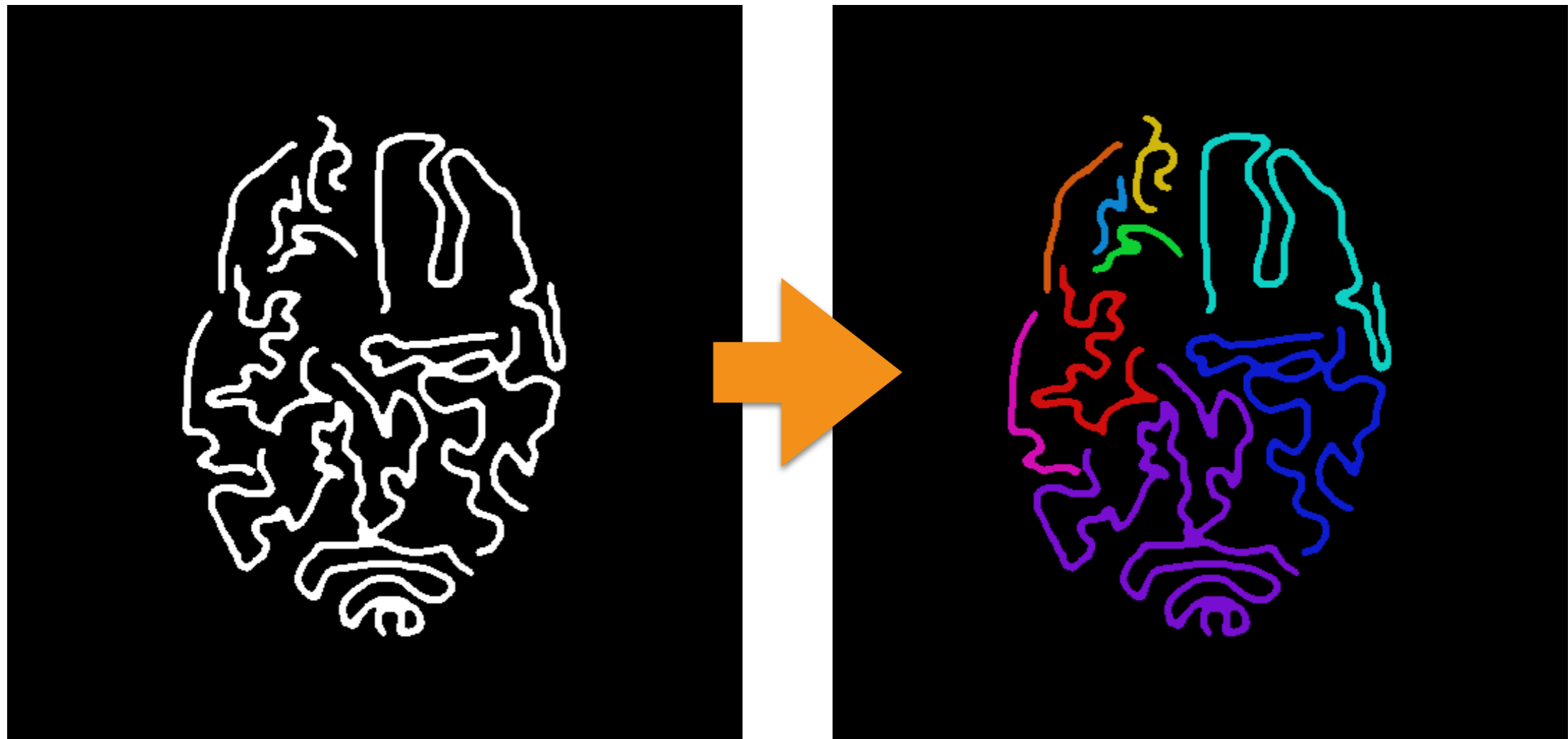
Thresholding: Connected Components



Thresholding: Connected Components



Thresholding: Connected Components Example



Thresholding

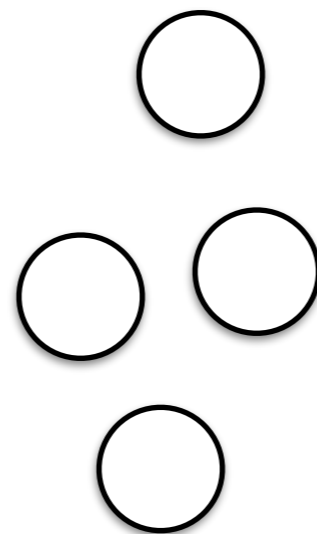
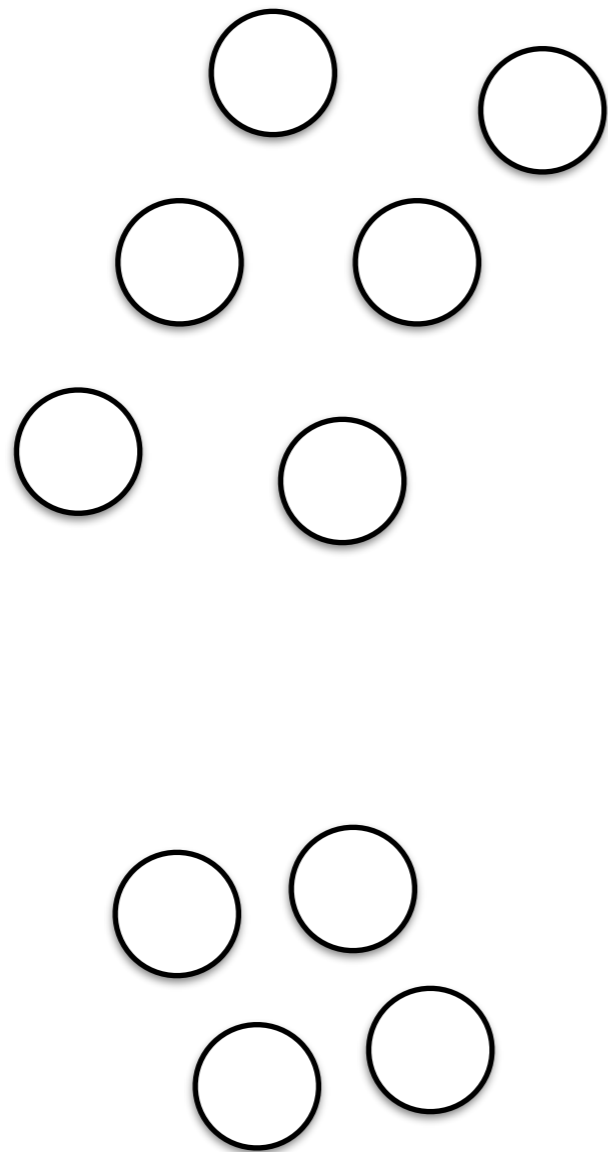
- It works if each object has a unique intensity value/color; this is a very limiting constraint!
 - However, it could be used as a starting point for other algorithms.
- The user needs to set the threshold!
 - The I_t value for each class may be inferred by analyzing the histogram of the input image.
- Its 3D extension is trivial!

k-Means

k -Means

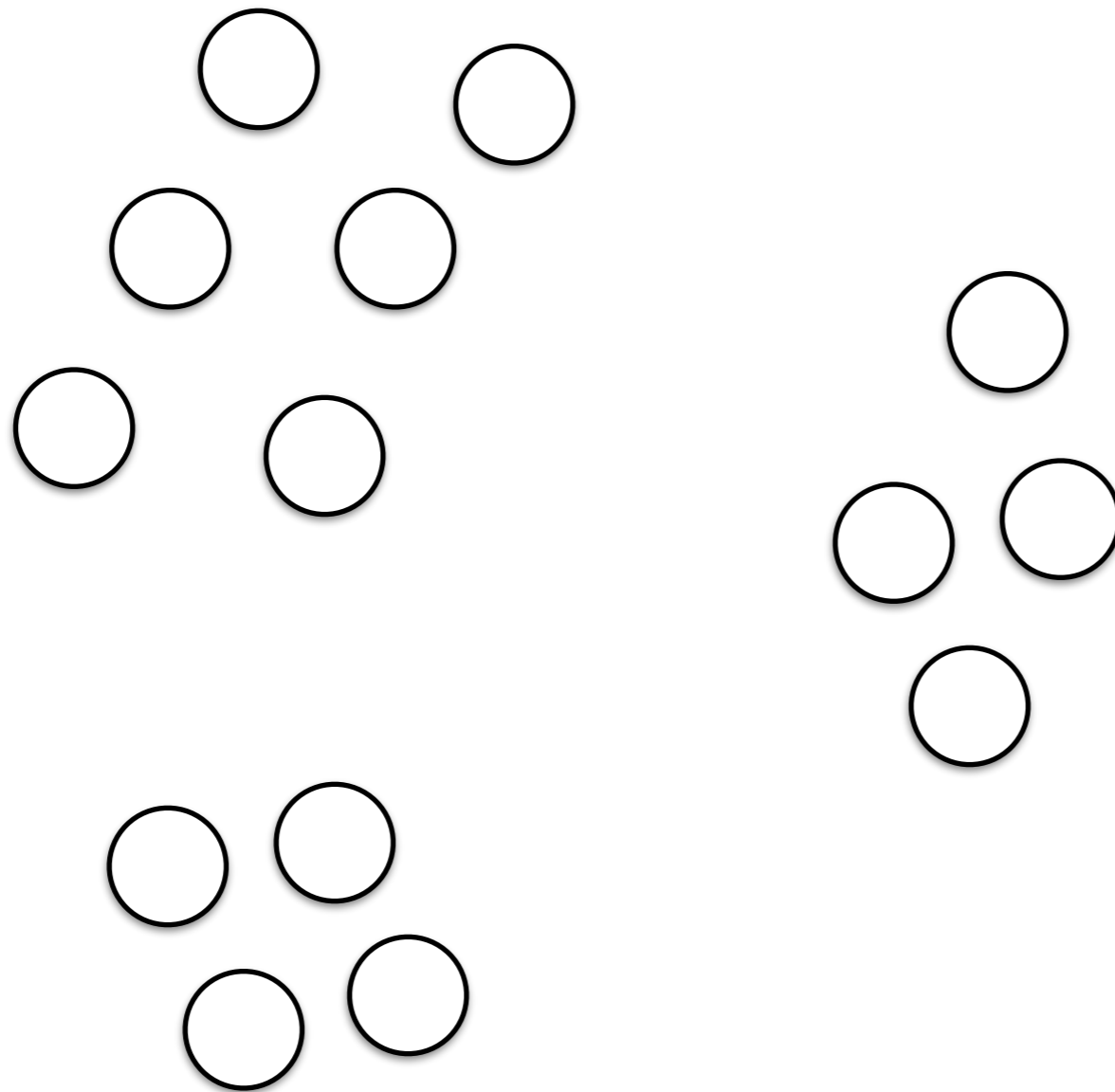
- k -means is a clustering algorithm for clustering n -D vectors/points in an n -D space:
 - A pixel with position (x, y) and intensity l is a 3D vector: $\langle x, y, l \rangle$
 - A voxel with position (x, y, z) and intensity l is a 4D vector: $\langle x, y, z, l \rangle$
- Let's assume we have k objects in the image.
- So we have to determine k -clusters.

k -Means: How it Works



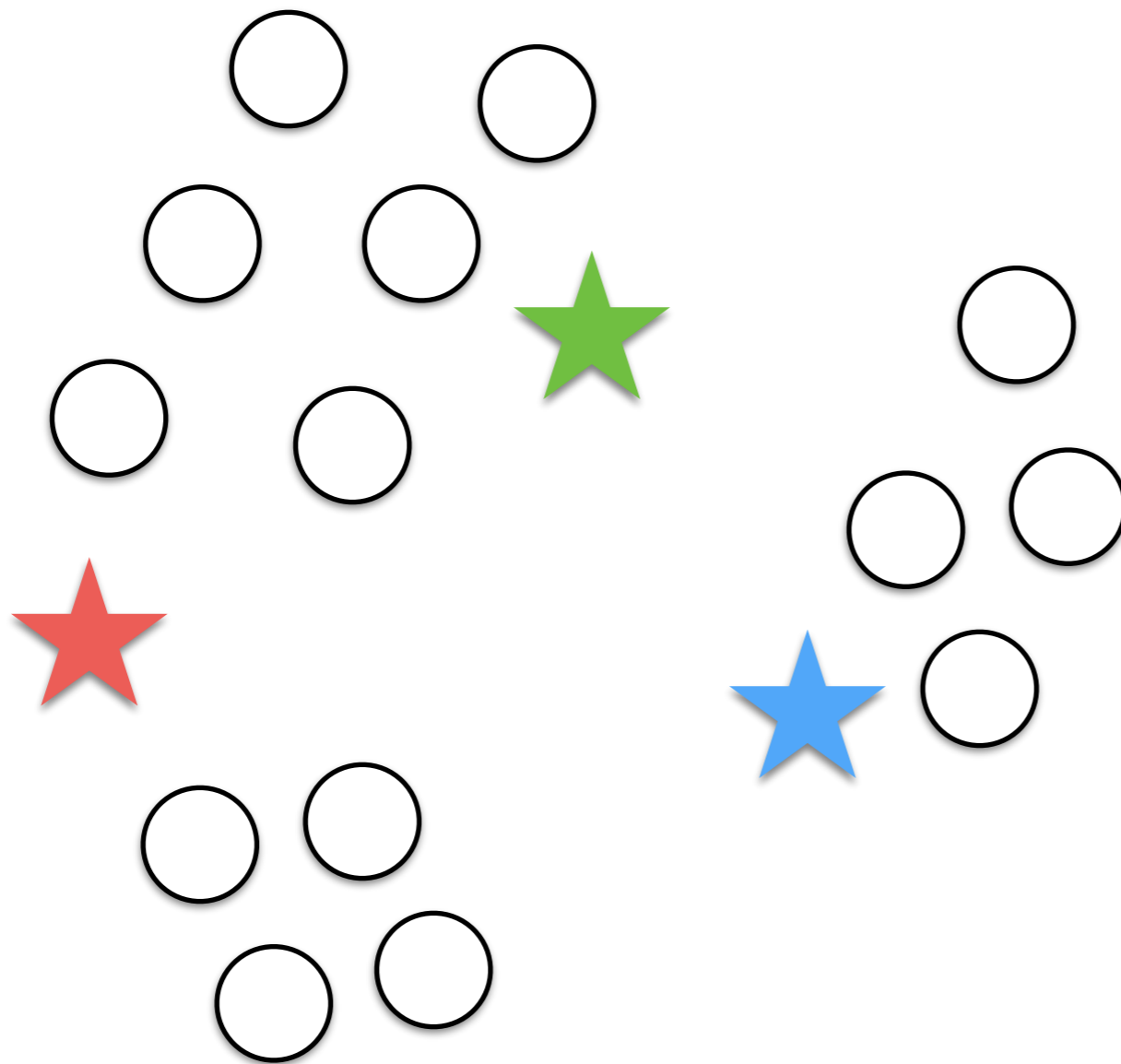
- A 2D example

k -Means: Initialization



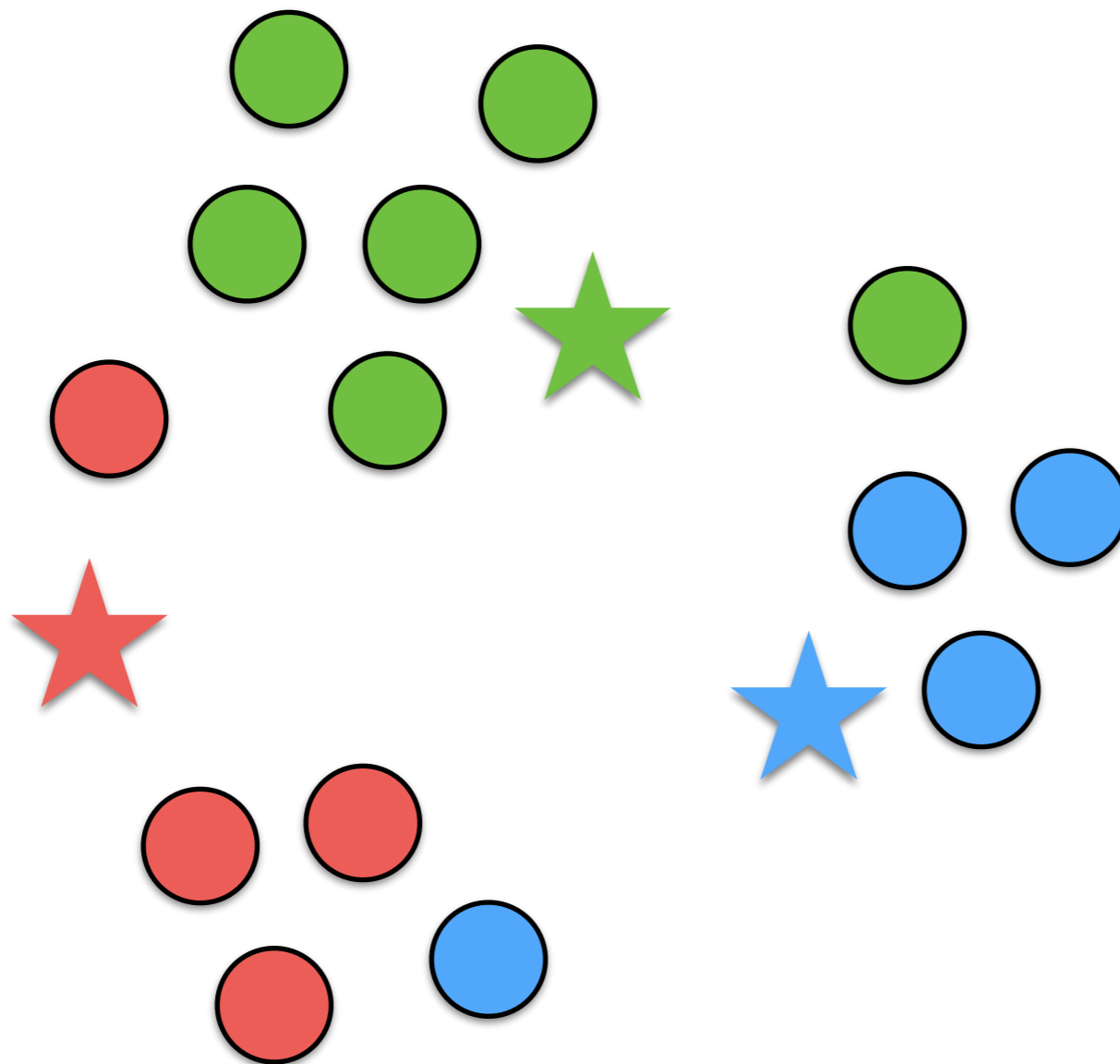
- Let's assume $k = 3$
- We make a random guess on the k -centroids; i.e., the stars.

k -Means: Initialization



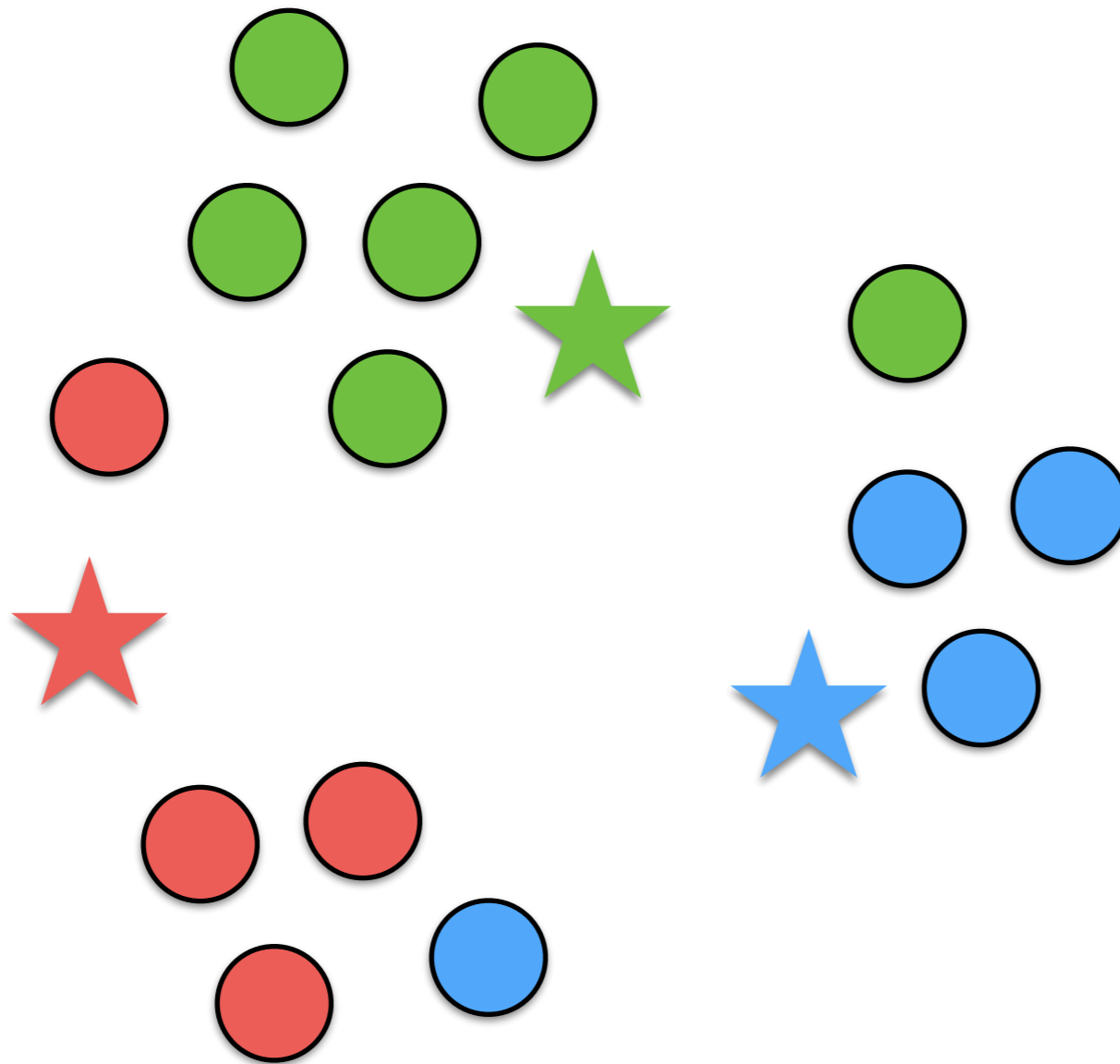
- Let's assume $k = 3$
- We make a random guess on the k -centroids; i.e., the stars.

k -Means: Iteration



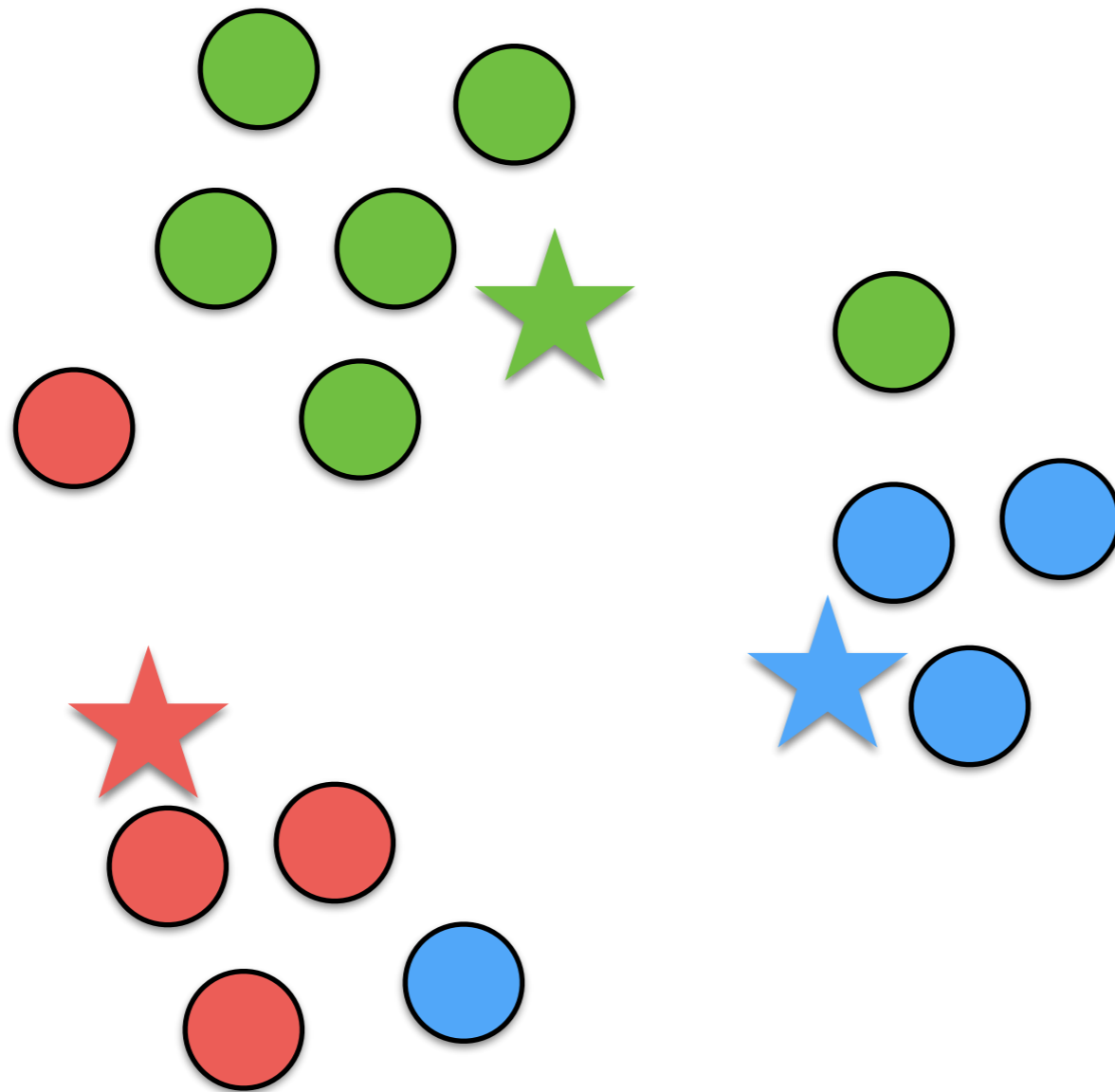
- We now assign a sample to a cluster if the distance (L1, L2, etc.), between a centroid is the minimum.

k -Means: Iteration



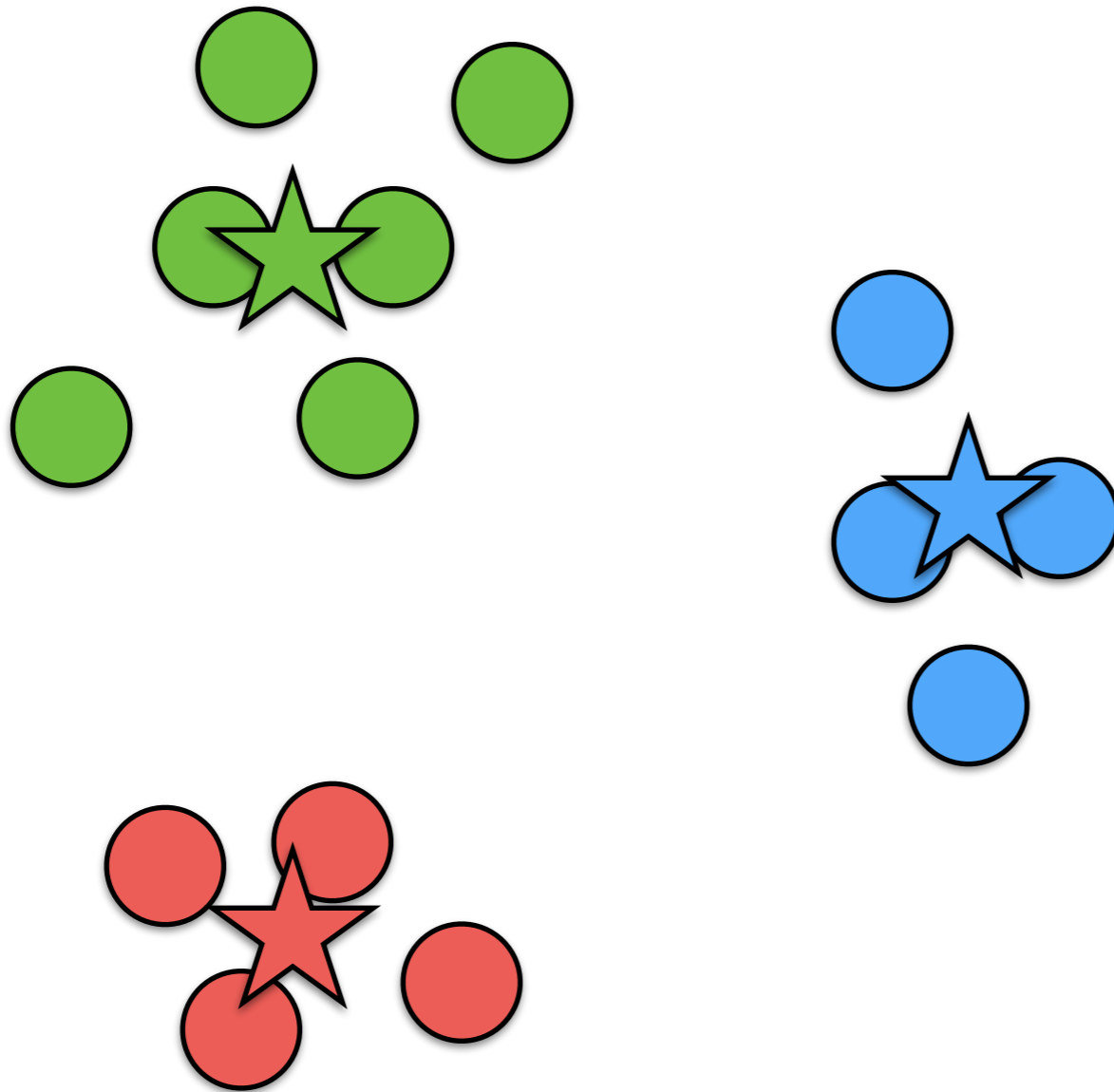
- We re-compute the centroid as the mean of samples of a cluster.

k -Means: Iteration



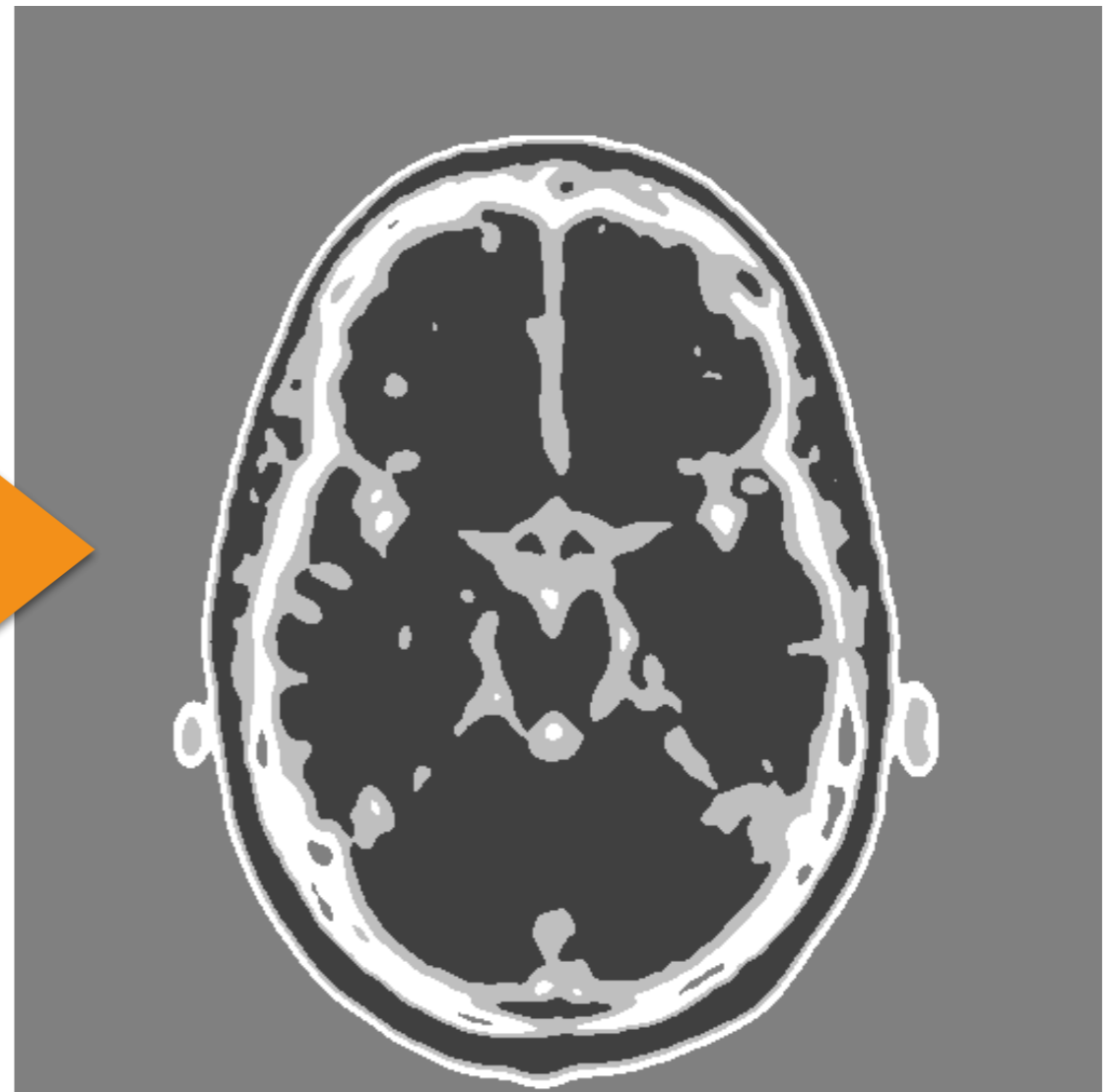
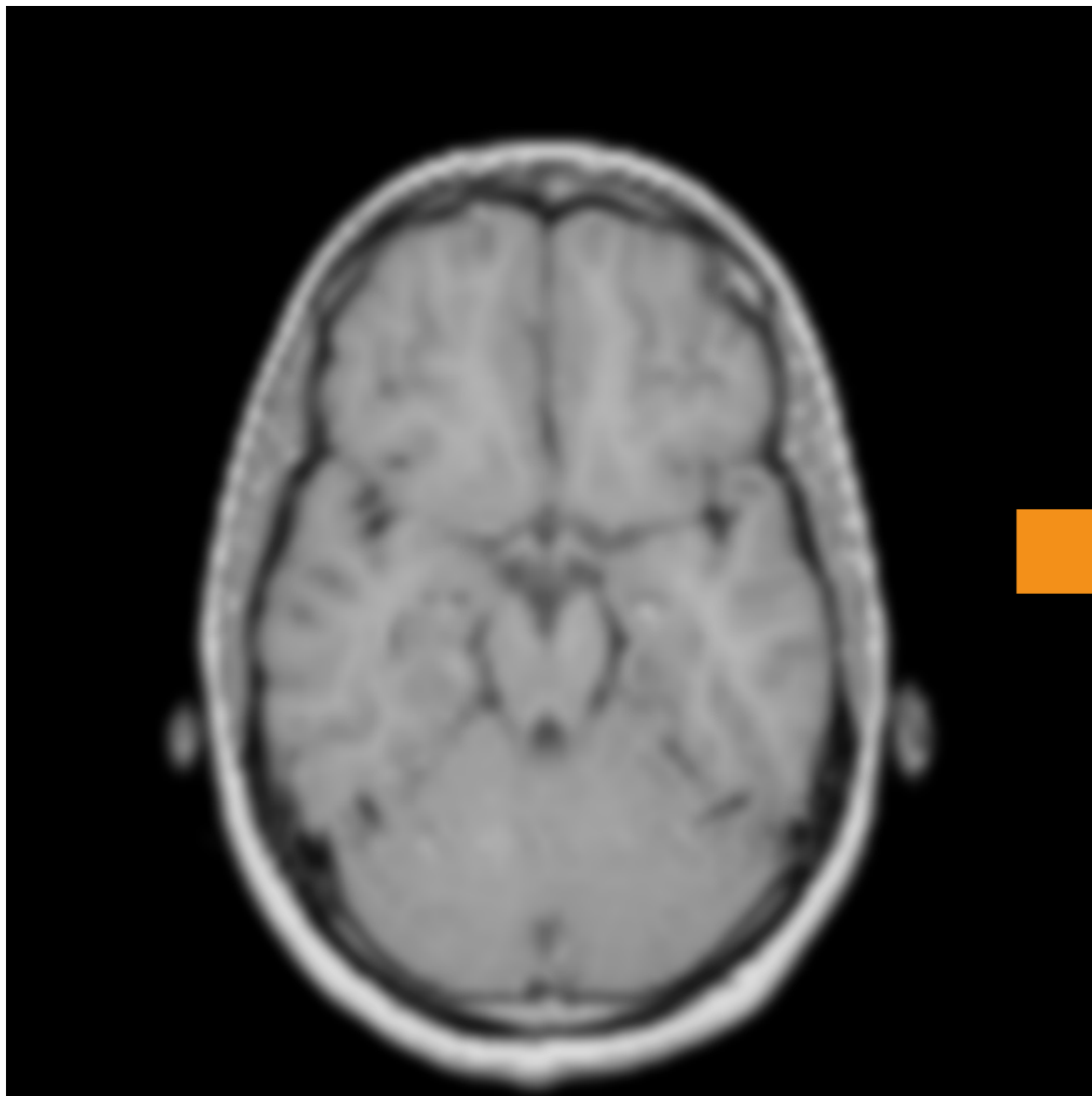
- We repeat the process until convergence (no more changes) or after m iterations.

k -Means: Iteration

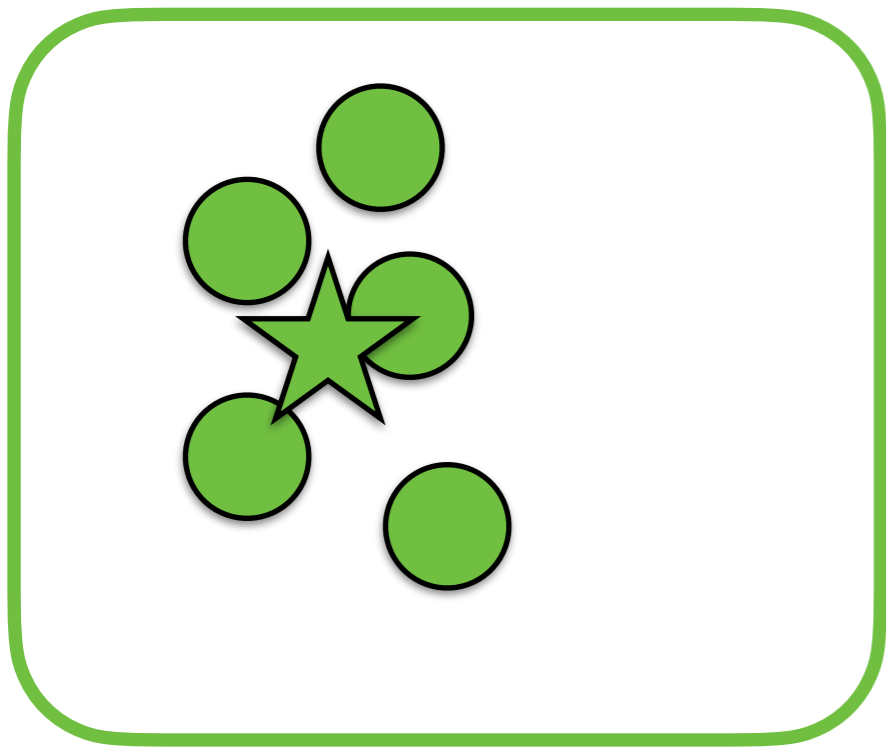


- We repeat the process until convergence (no more changes) or after m iterations.

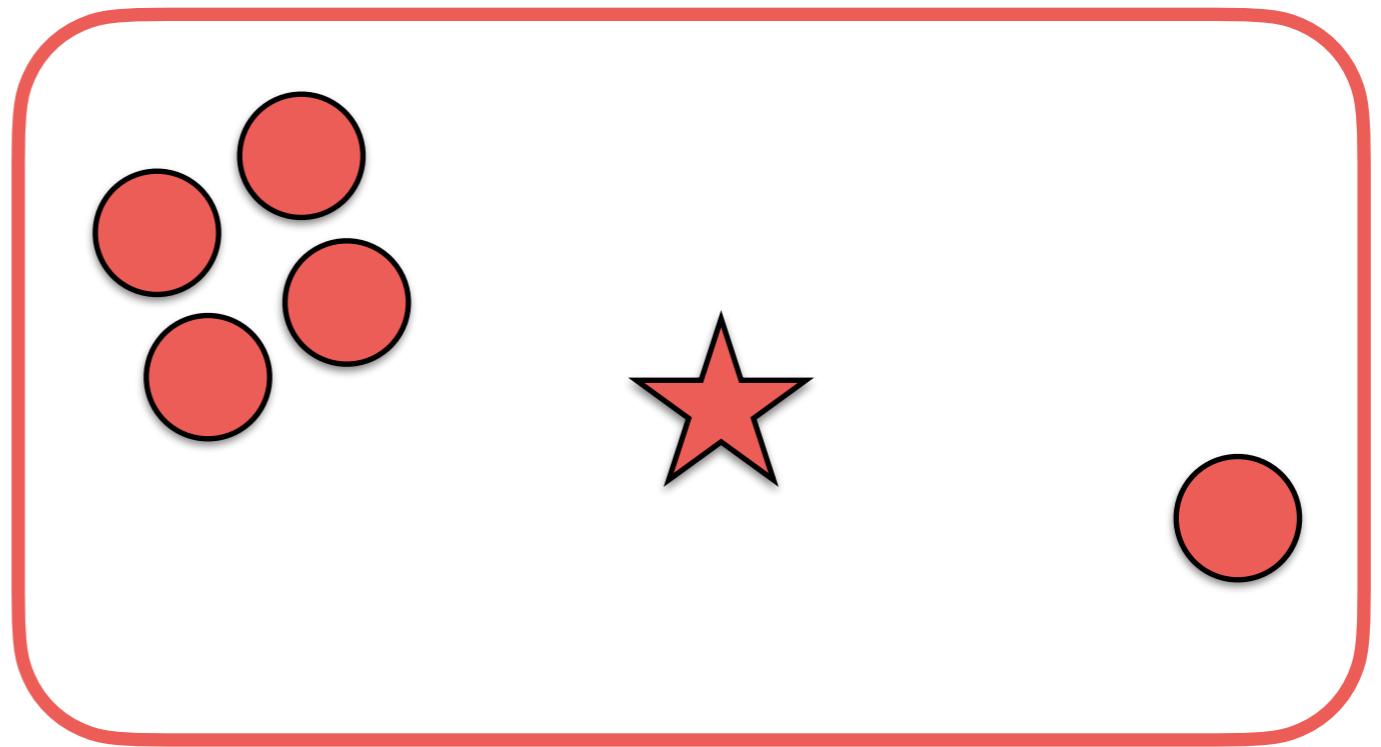
k -Means Example



k -Means: Outliers

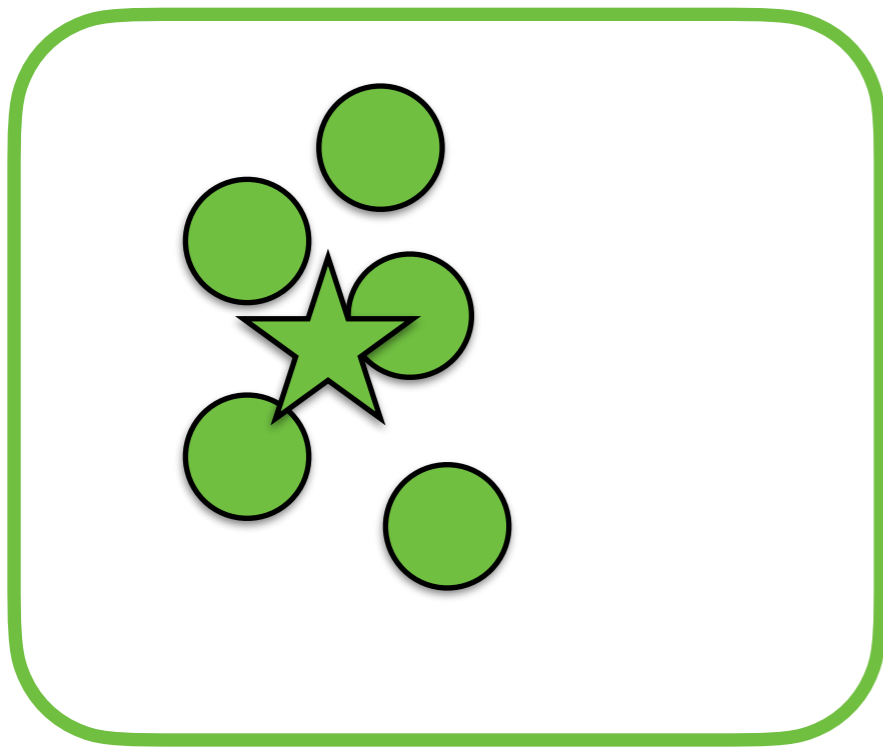


Cluster A

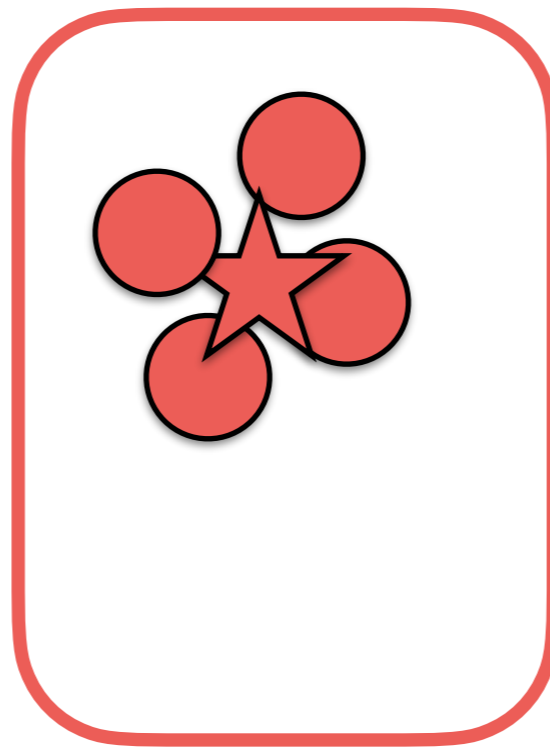


Cluster B

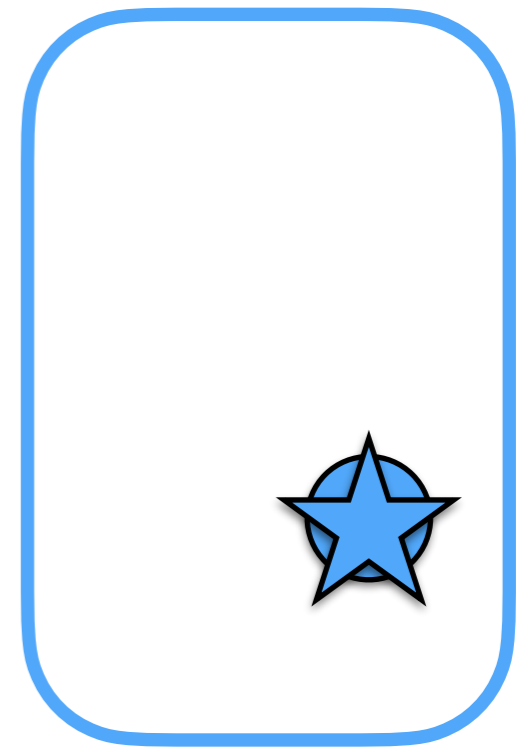
k -Means: Outliers Solution 1



Cluster A

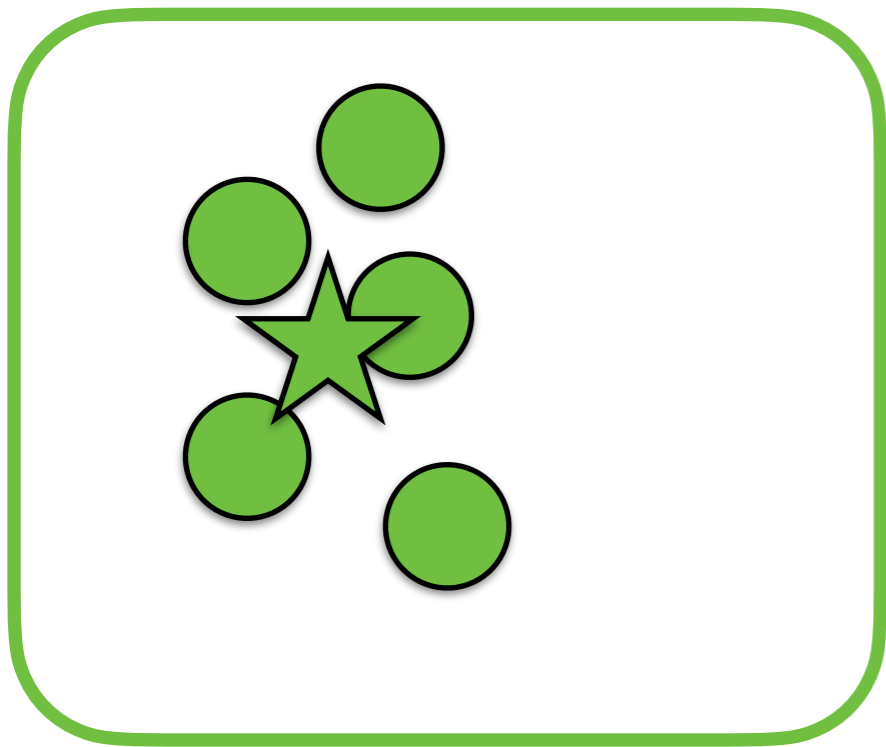


Cluster B

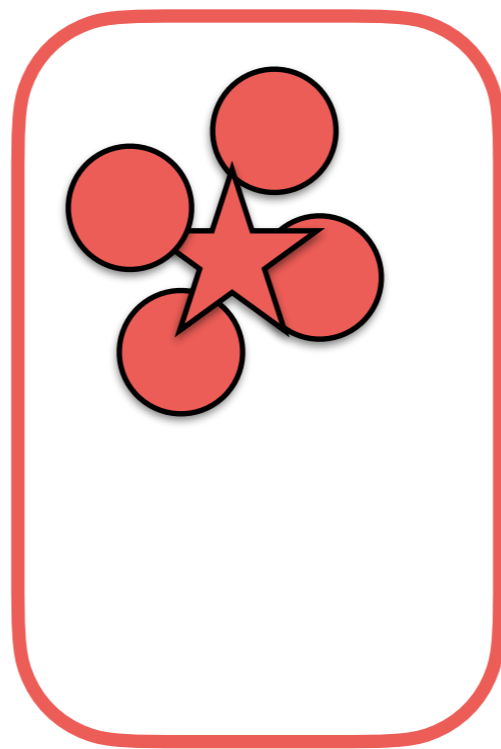


Cluster C

k -Means: Outliers Solution 2

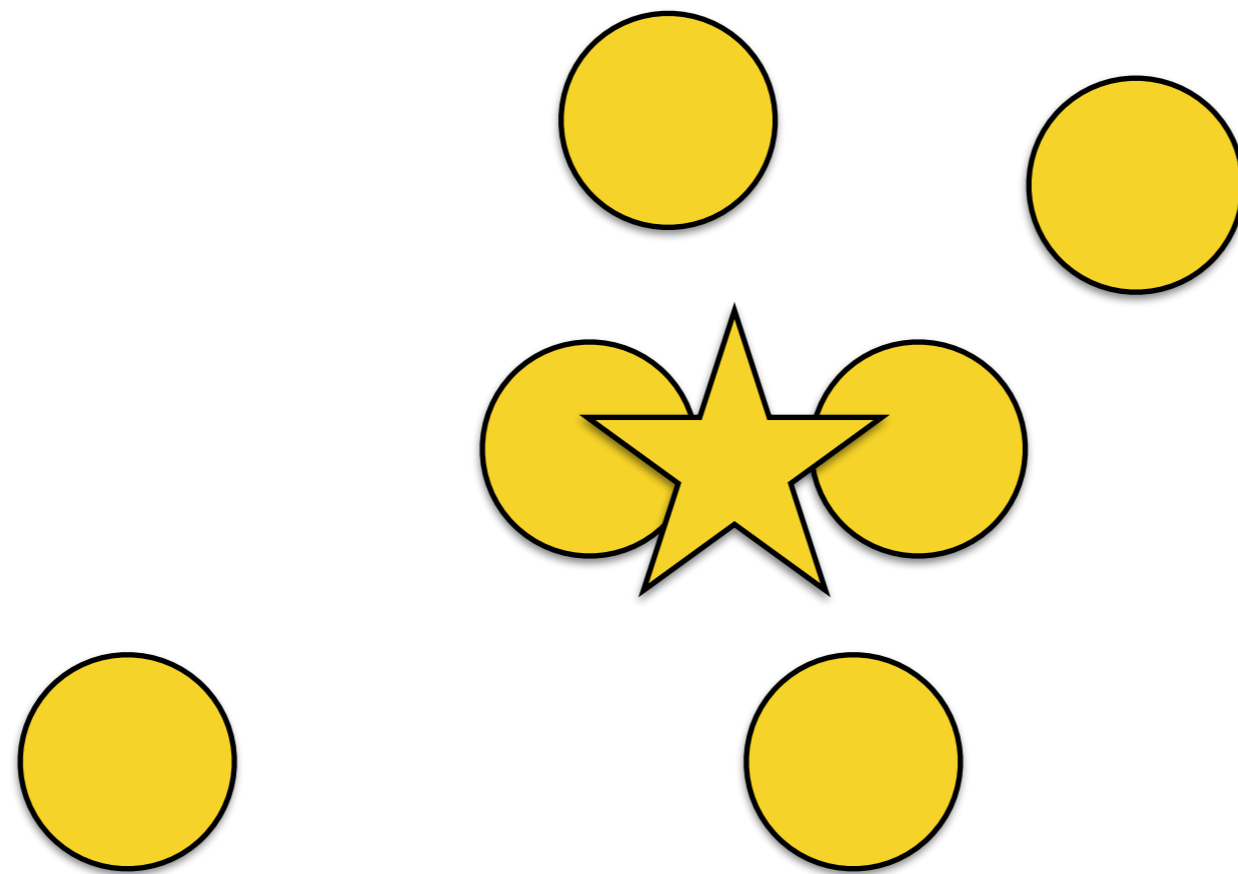


Cluster A



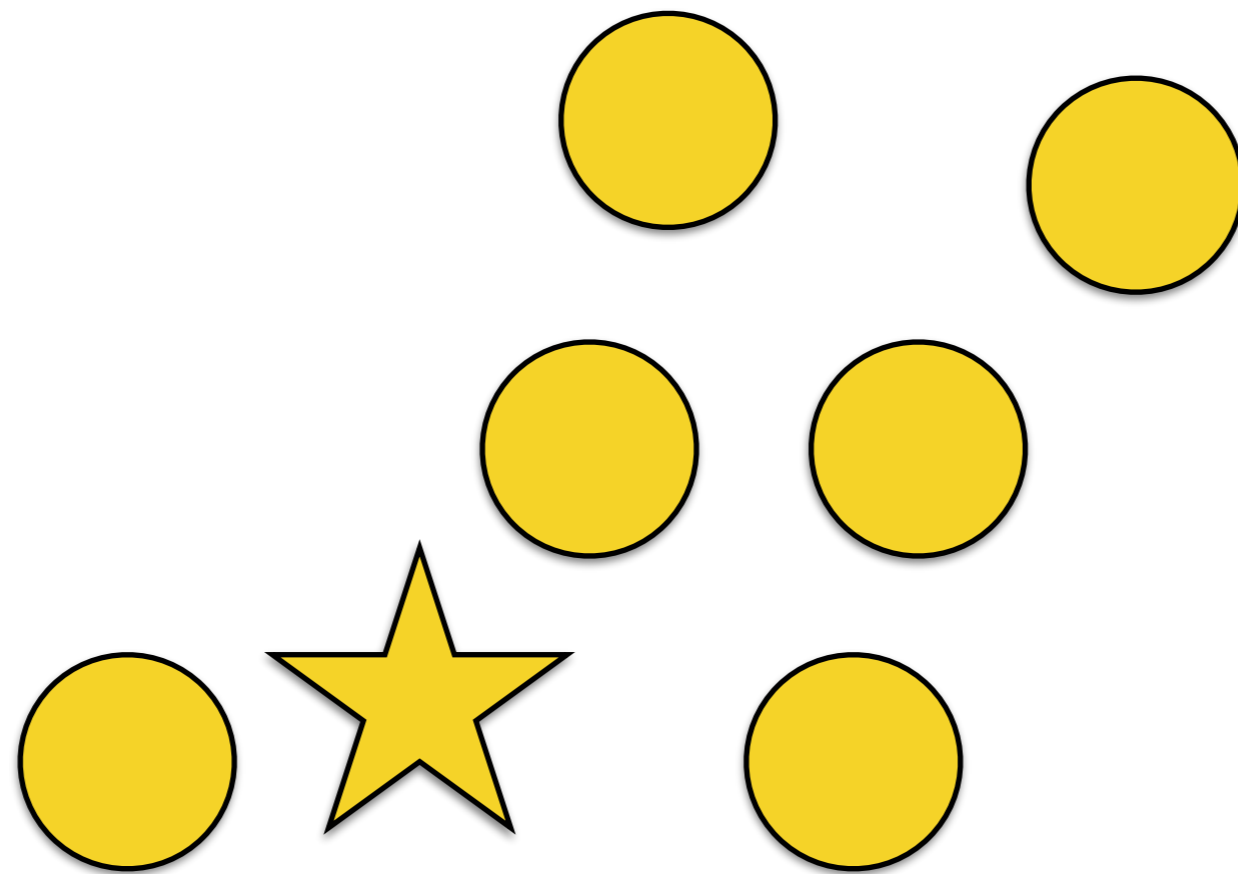
Cluster B

k -Means: Oscillation



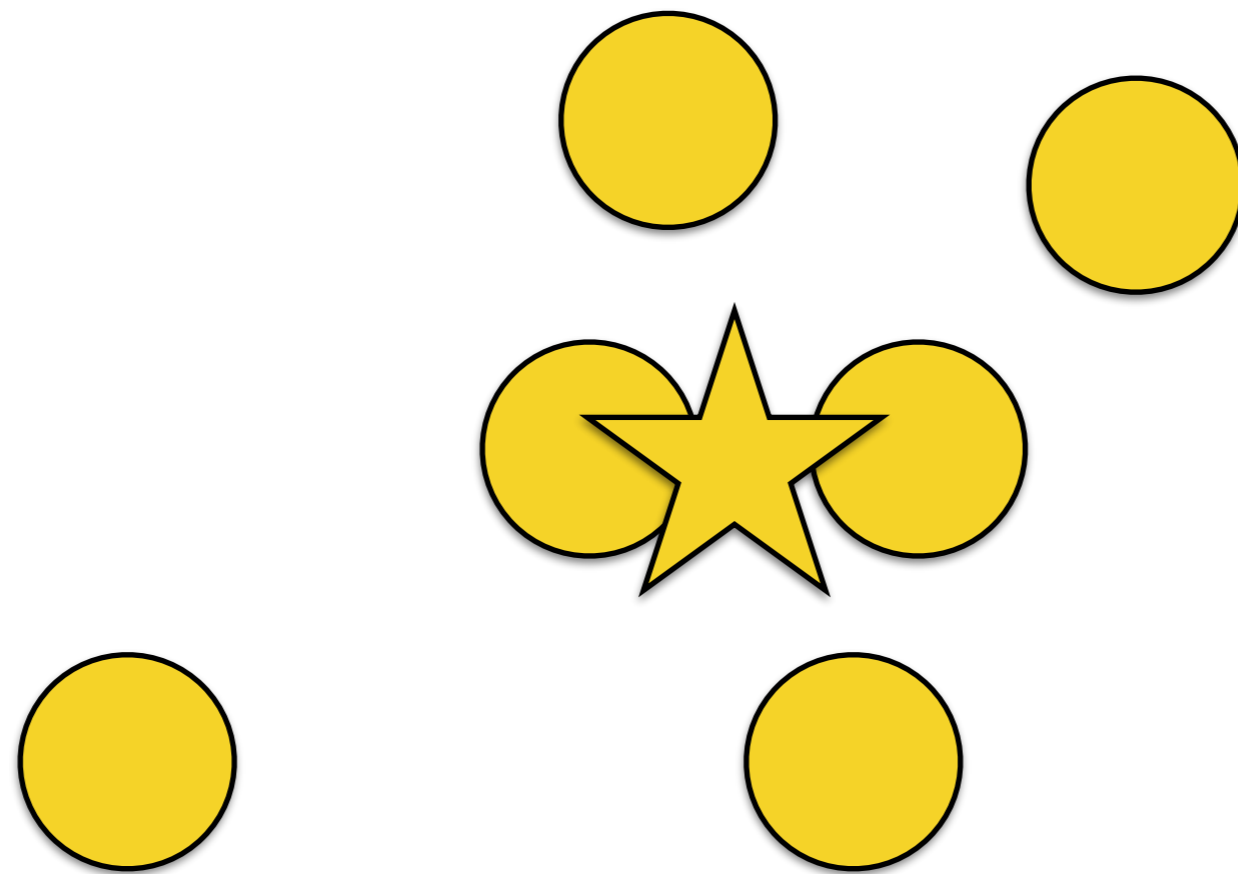
Even Iteration

k -Means: Oscillation



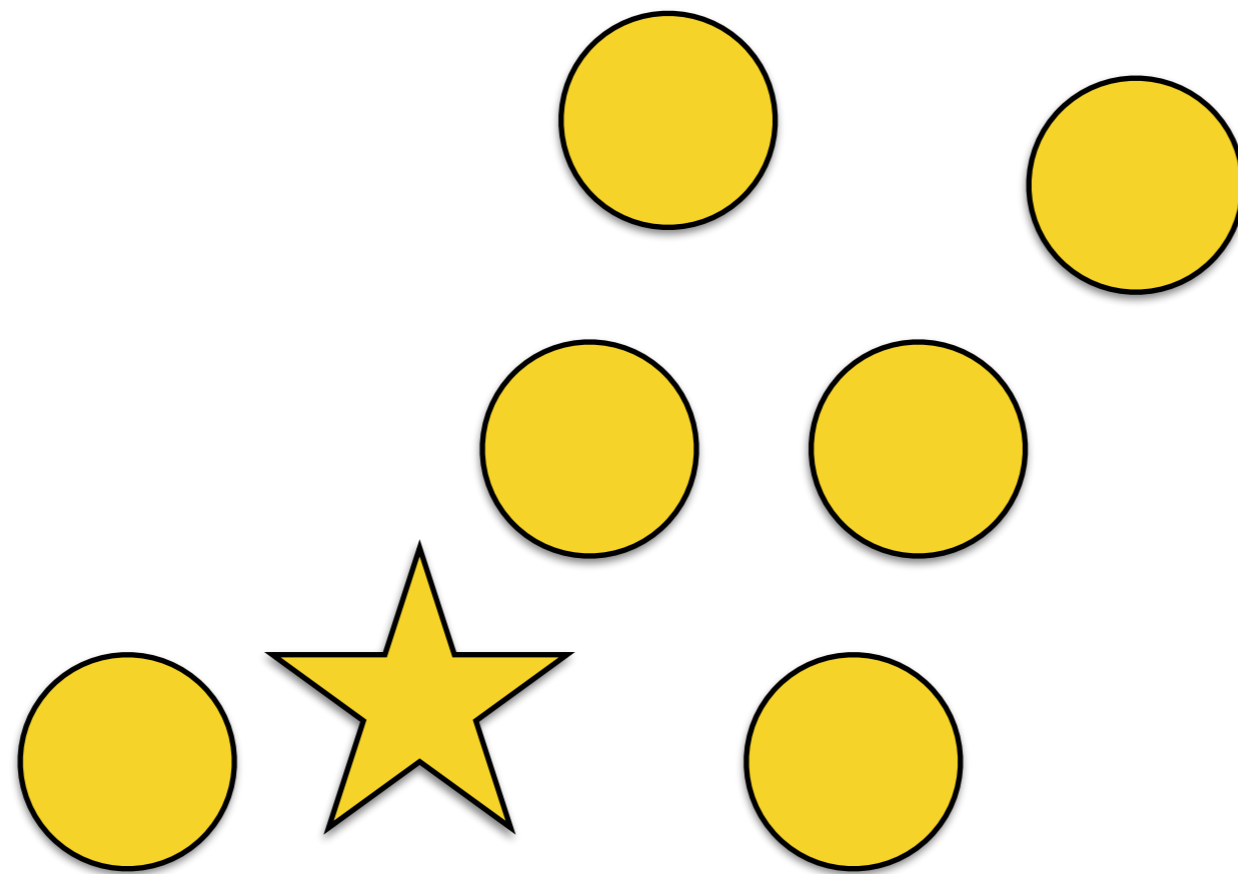
Odd Iteration

k -Means: Oscillation



Even Iteration

k -Means: Oscillation



Odd Iteration

k -Means: Advantages

- The method is fully automatic.
- This works for 2D and 3D volumes.
- This can “understand” neighbors in an implicit way.

k -Means: Disadvantages

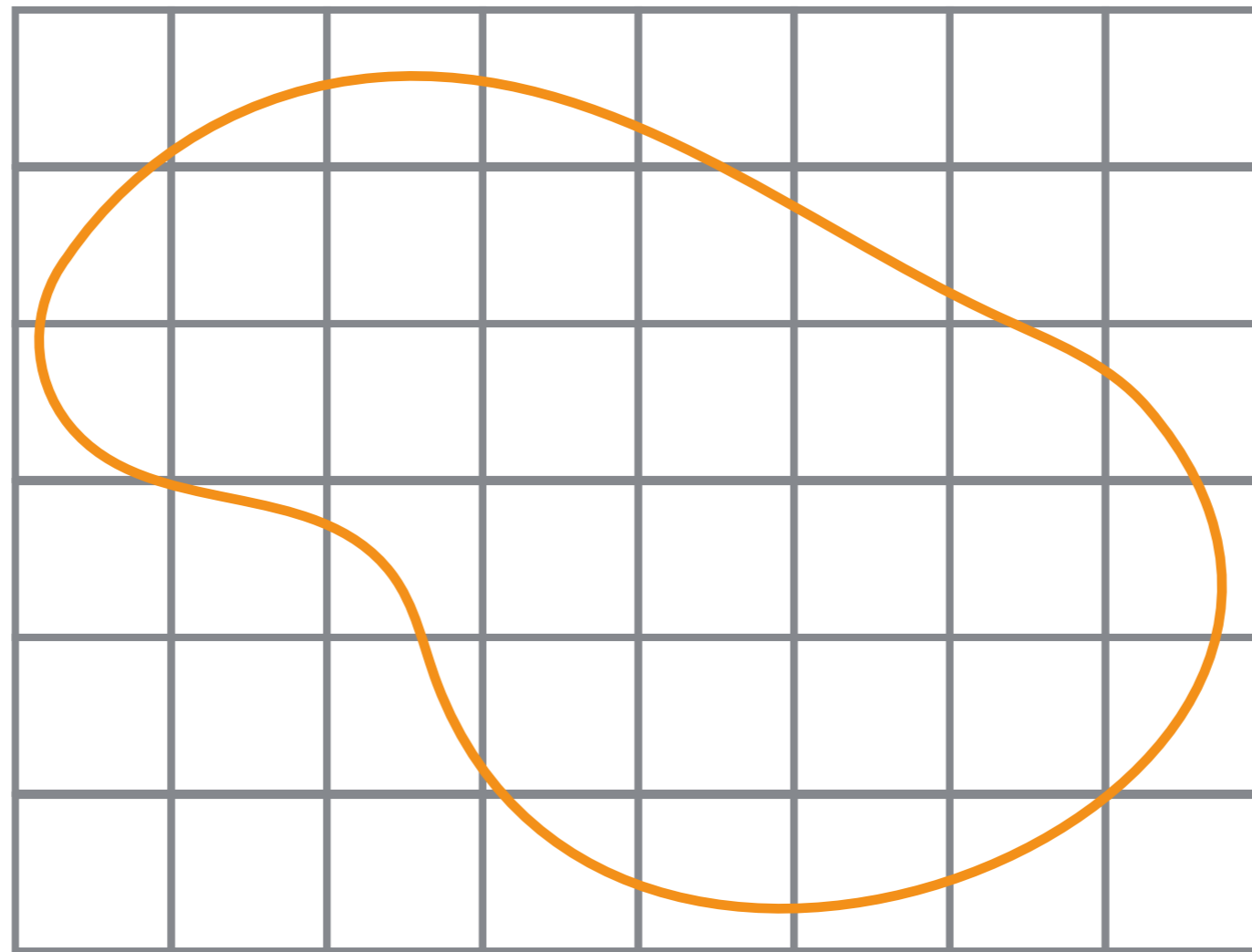
- We need to know how many objects (including the background) are in the image:
 - We may run k-means multiple times until a certain criterion is met; e.g., reaching the 80% of percentage of explained variance.
- Outliers:
 - better initialization (sampling).
- The method **may not** converge.
- We need to set a **maximum** number of iterations.

Region Growing

Region Growing

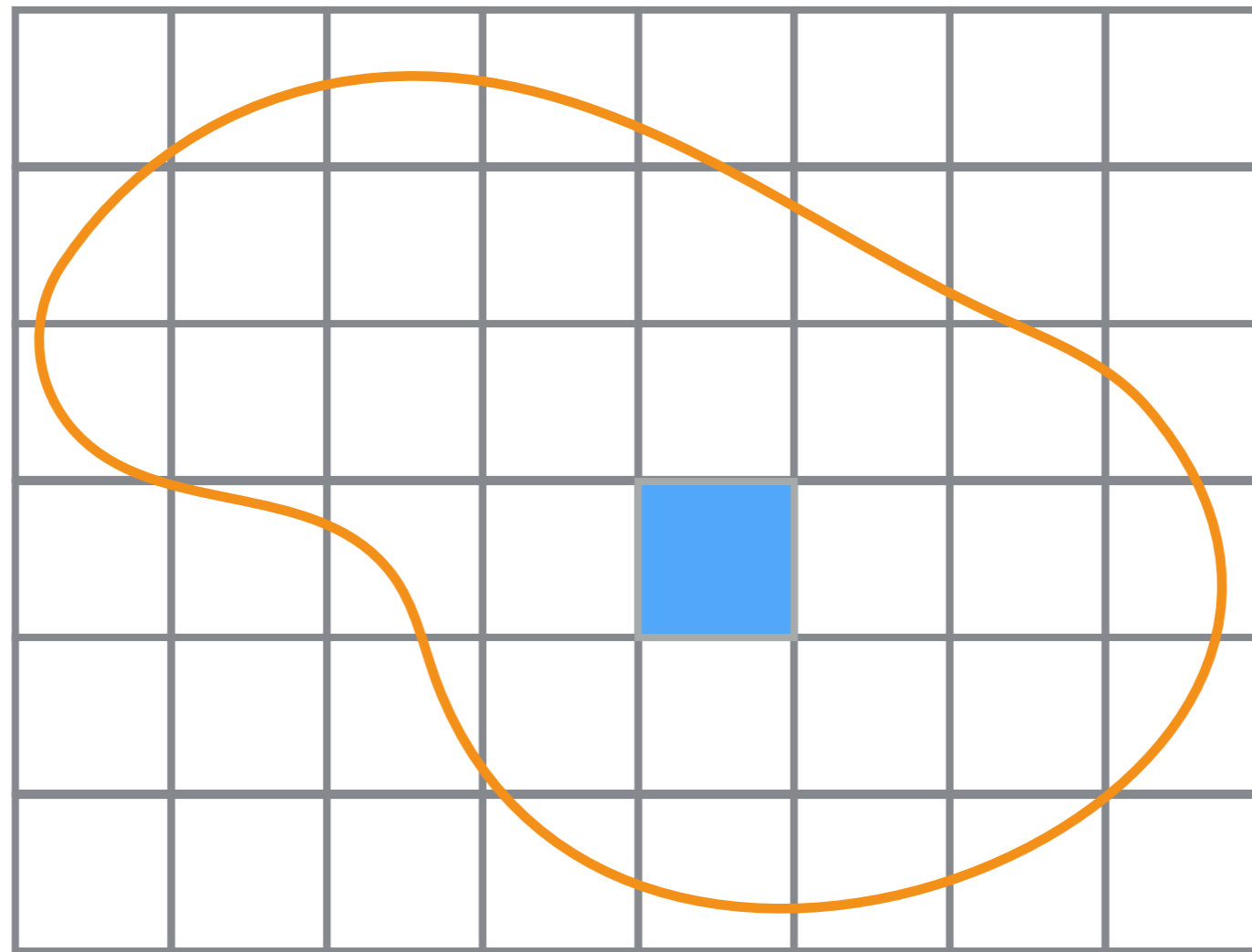
- This algorithm expands a painted initial mask until it reaches strong edges
- Therefore, we need to compute edges first!

Region Growing



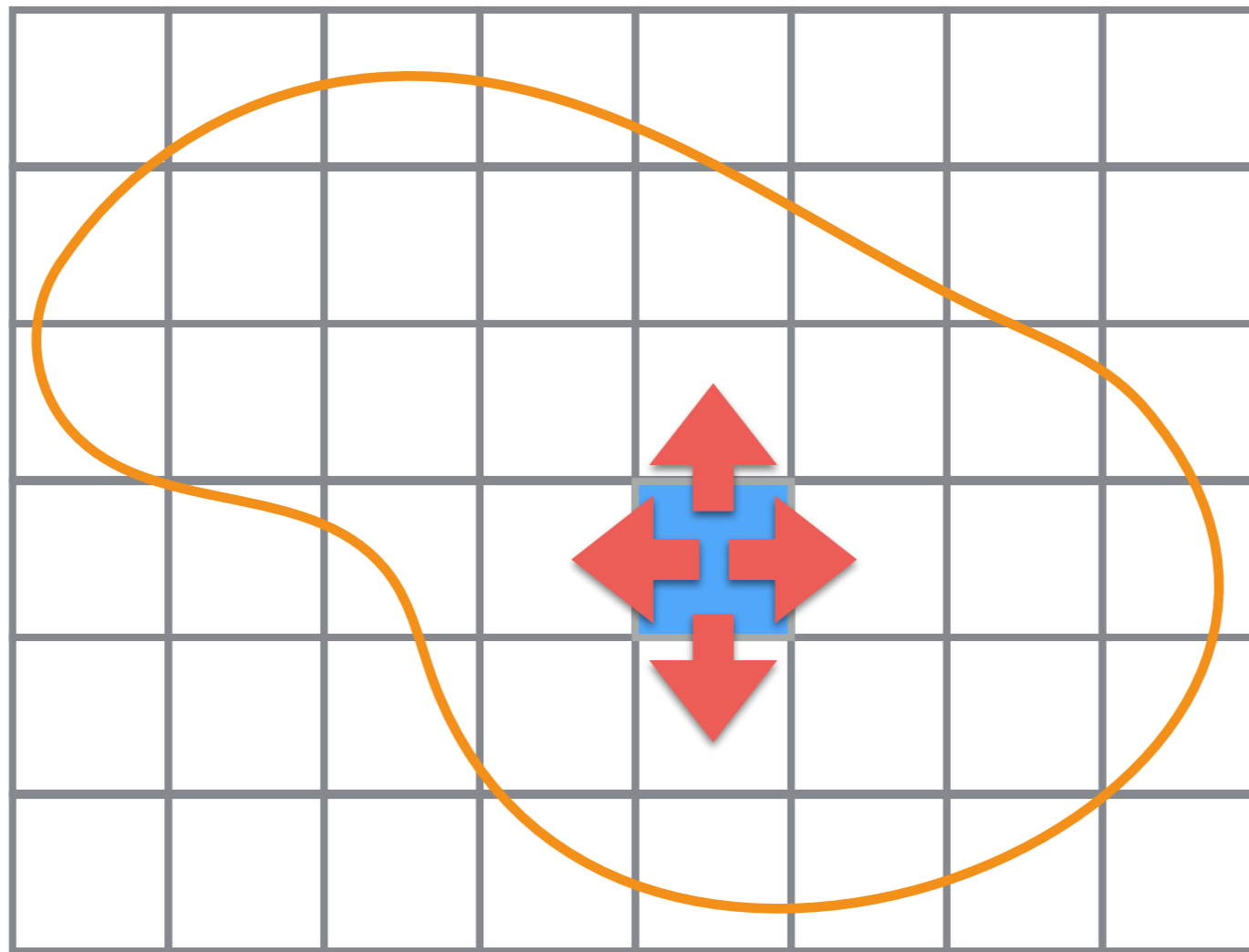
— Structure
Edge

Region Growing

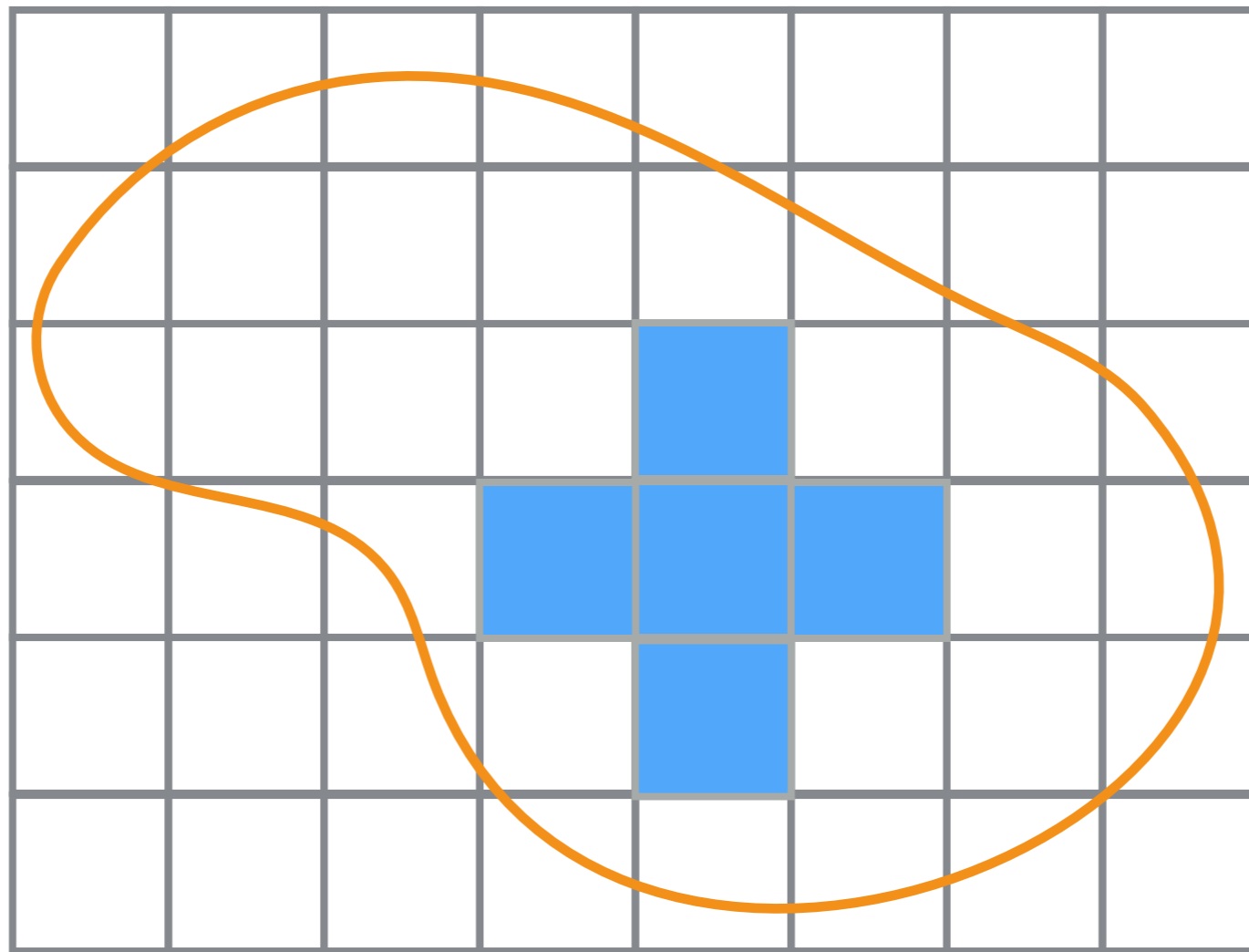


Seed

Region Growing

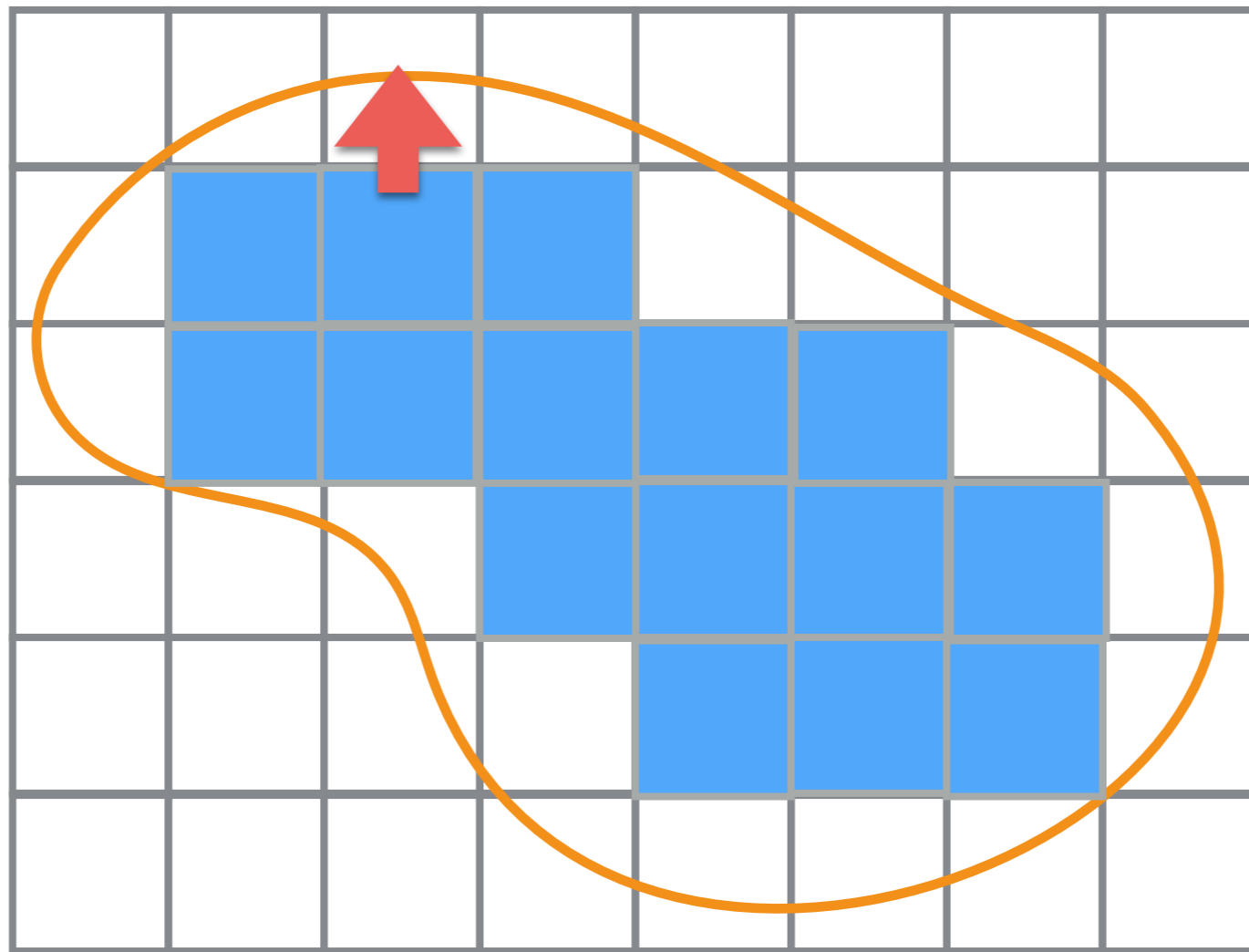


Region Growing

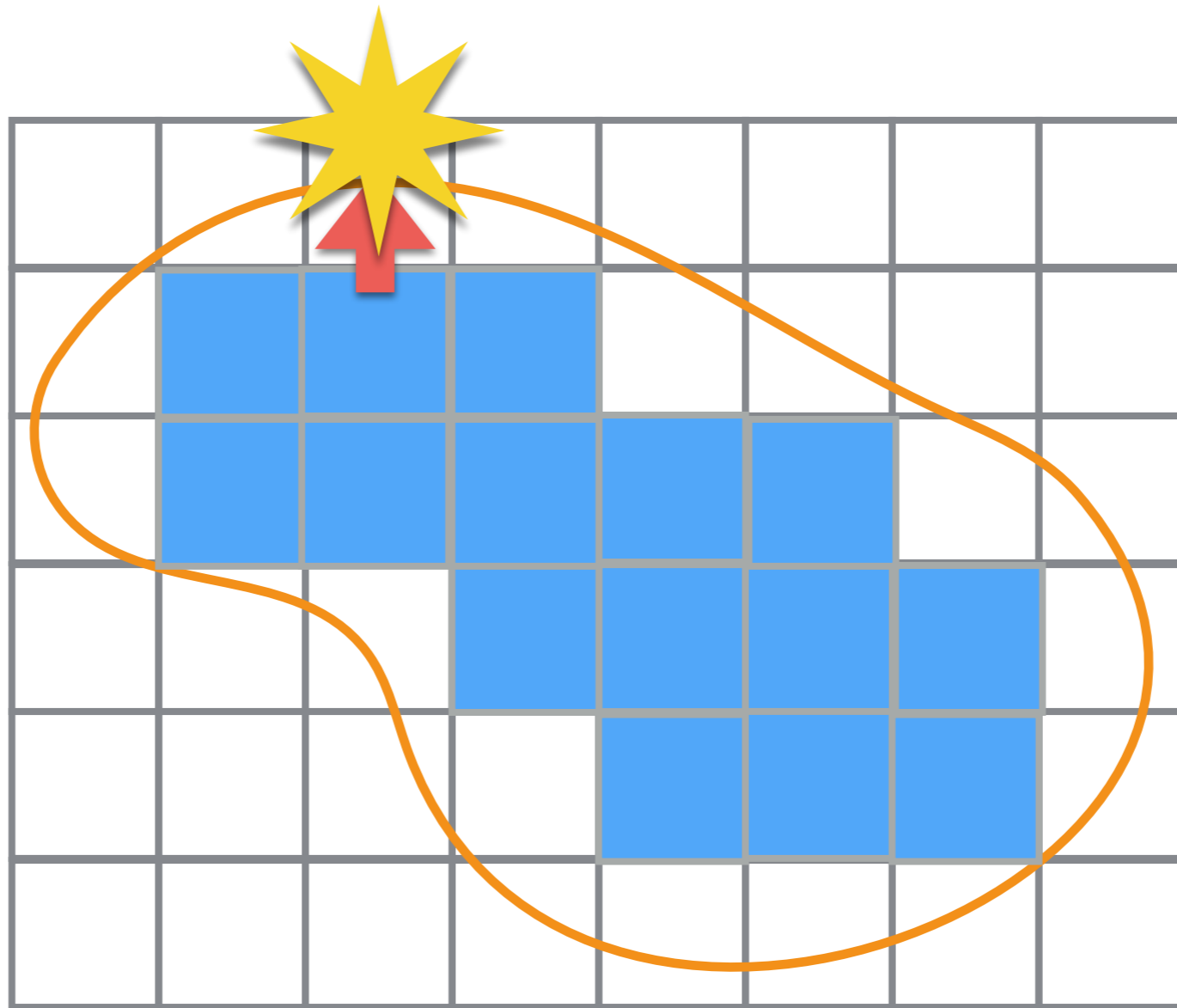


after a while...

Region Growing



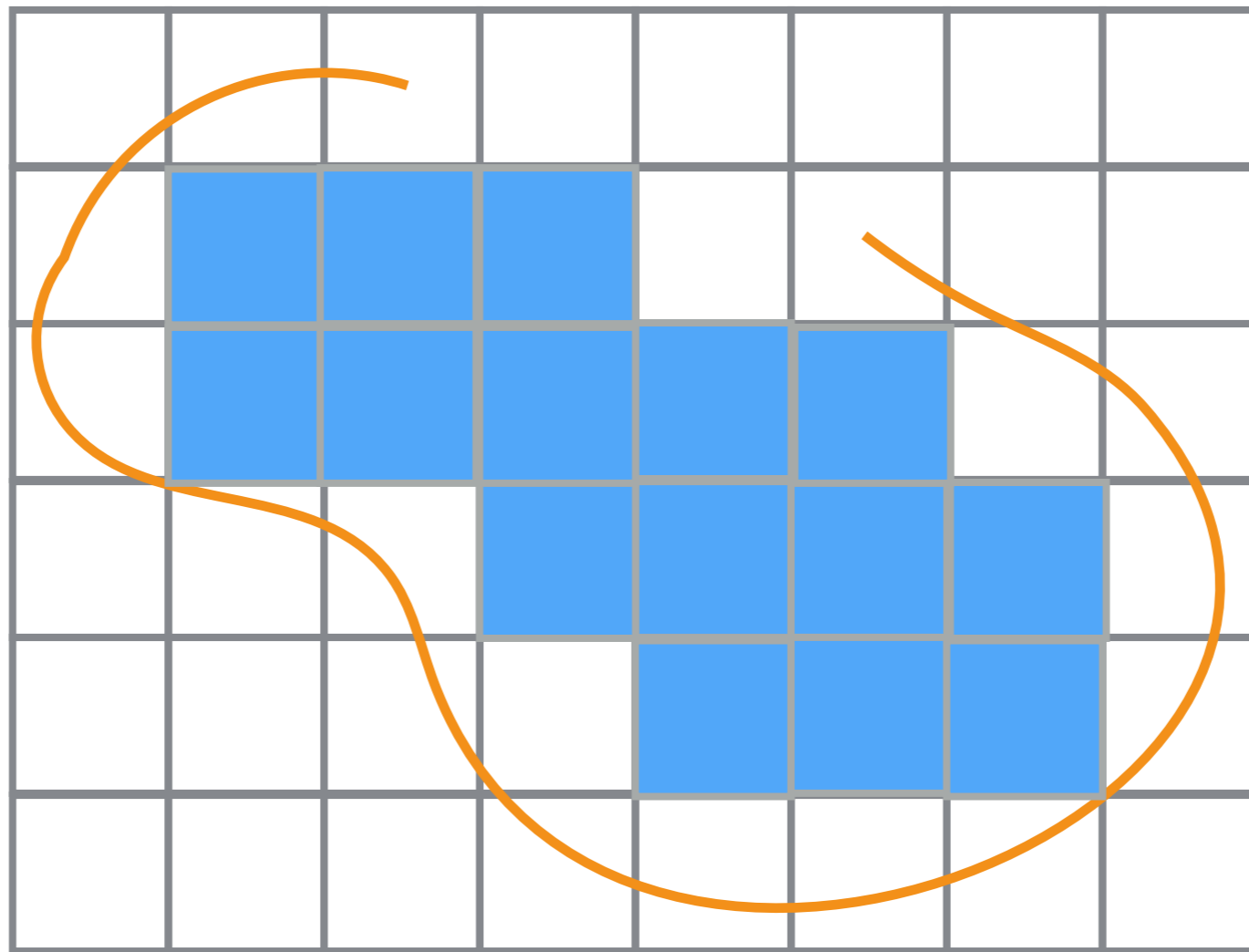
Region Growing



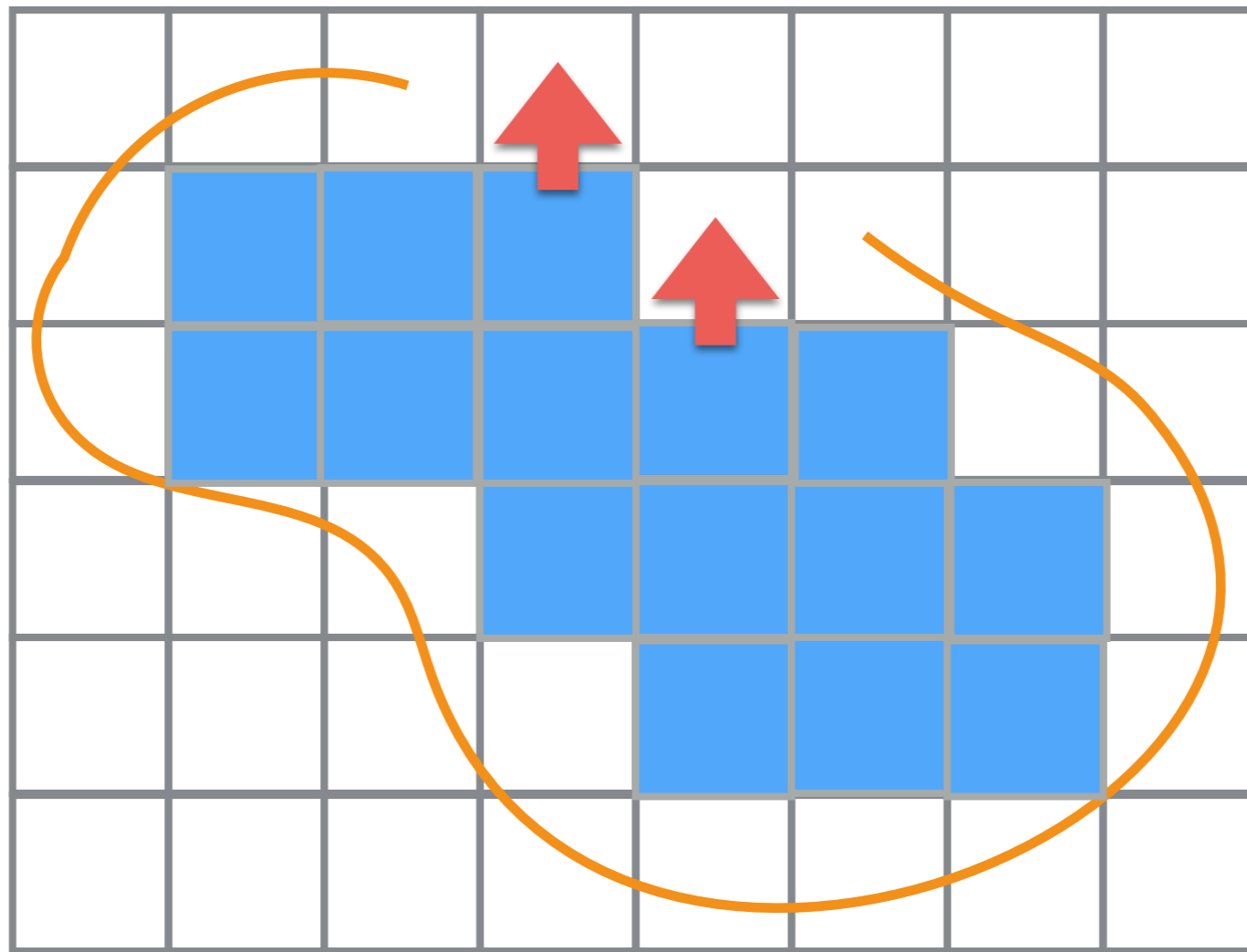
Region Growing

- It is straightforward to extend to 3D!
- This algorithm depends on:
 - The threshold of edge detection.
- It may be slow:
 - From an initial seed, the growing region needs to reach the farthest edge pixel/voxel.
 - Computational complexity is a function of the area/volume of the object we want to segment.

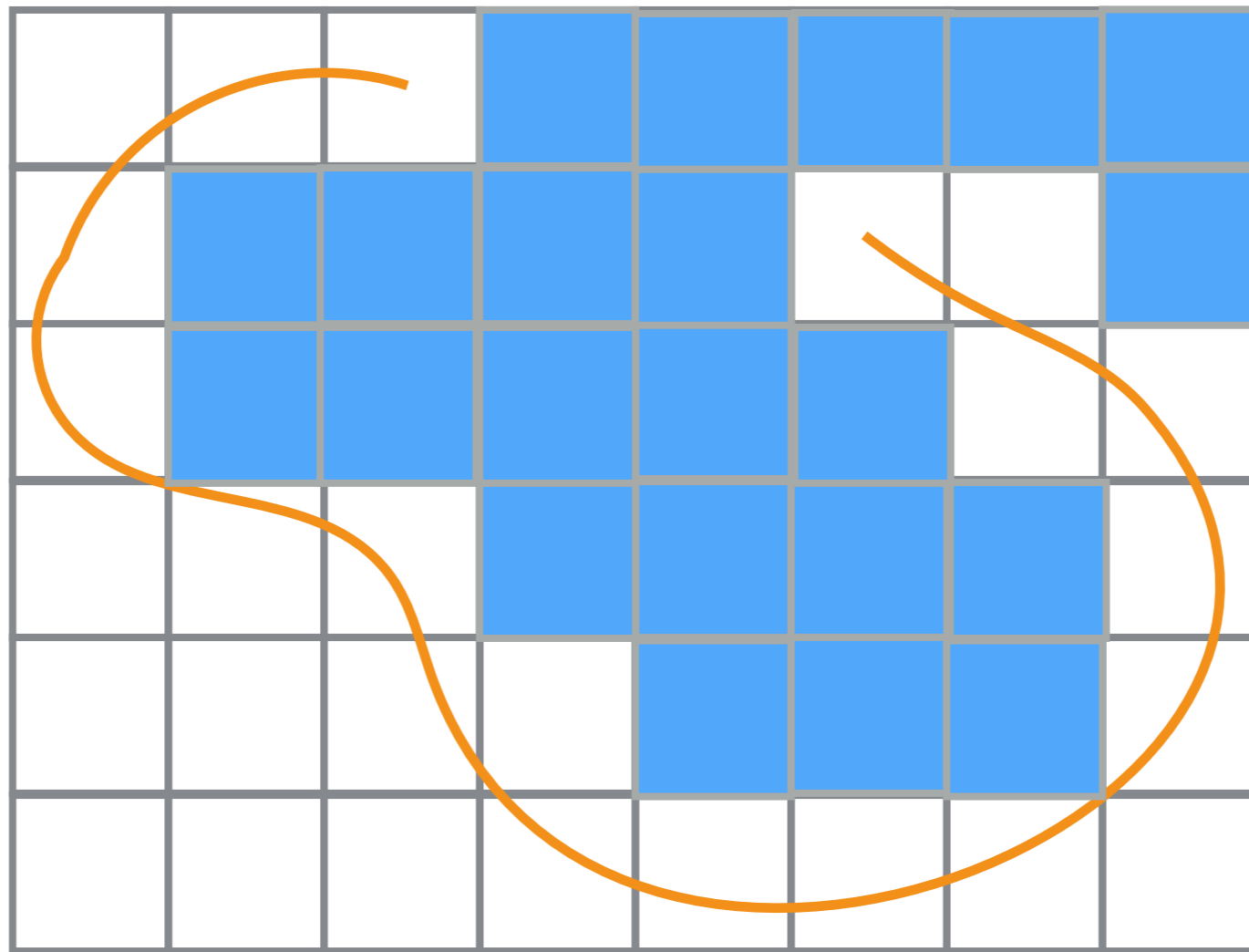
Region Growing: Epic Fail



Region Growing: Epic Fail



Region Growing: Epic Fail



Active Contour Model aka Snakes

Snakes

- A snake is a parametric curve:

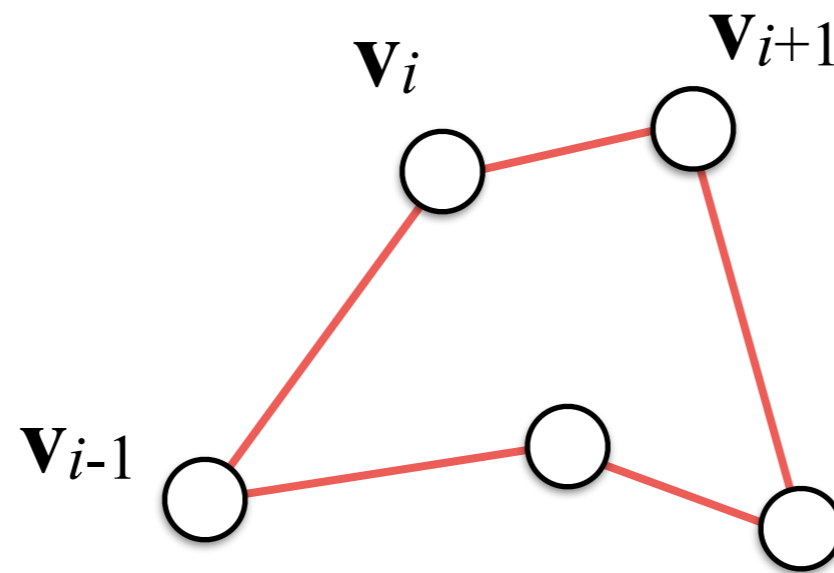
$$\mathbf{v}(t) = [x(t); y(t)] \quad t \in [0,1]$$

- Typically, it is a spline (original paper), but for sake of simplicity let's assume a piecewise linear curve.

Snakes

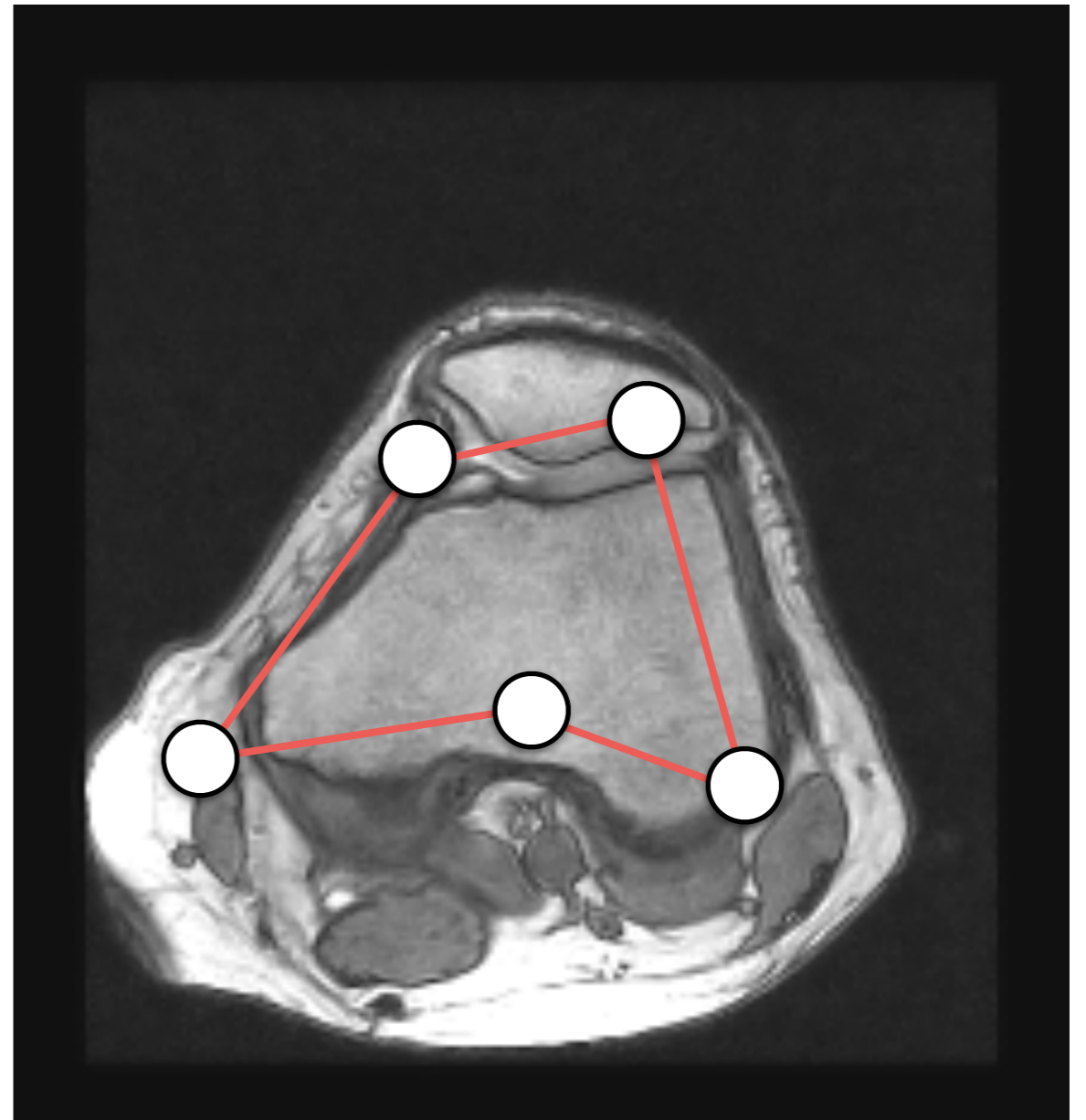
- The snake curve is defined by a set of control point that is defined as:

$$\mathbf{C} = \{\mathbf{v}_i | i \in [1, n]\} \quad \mathbf{v}_i = [x_i, y_i]$$



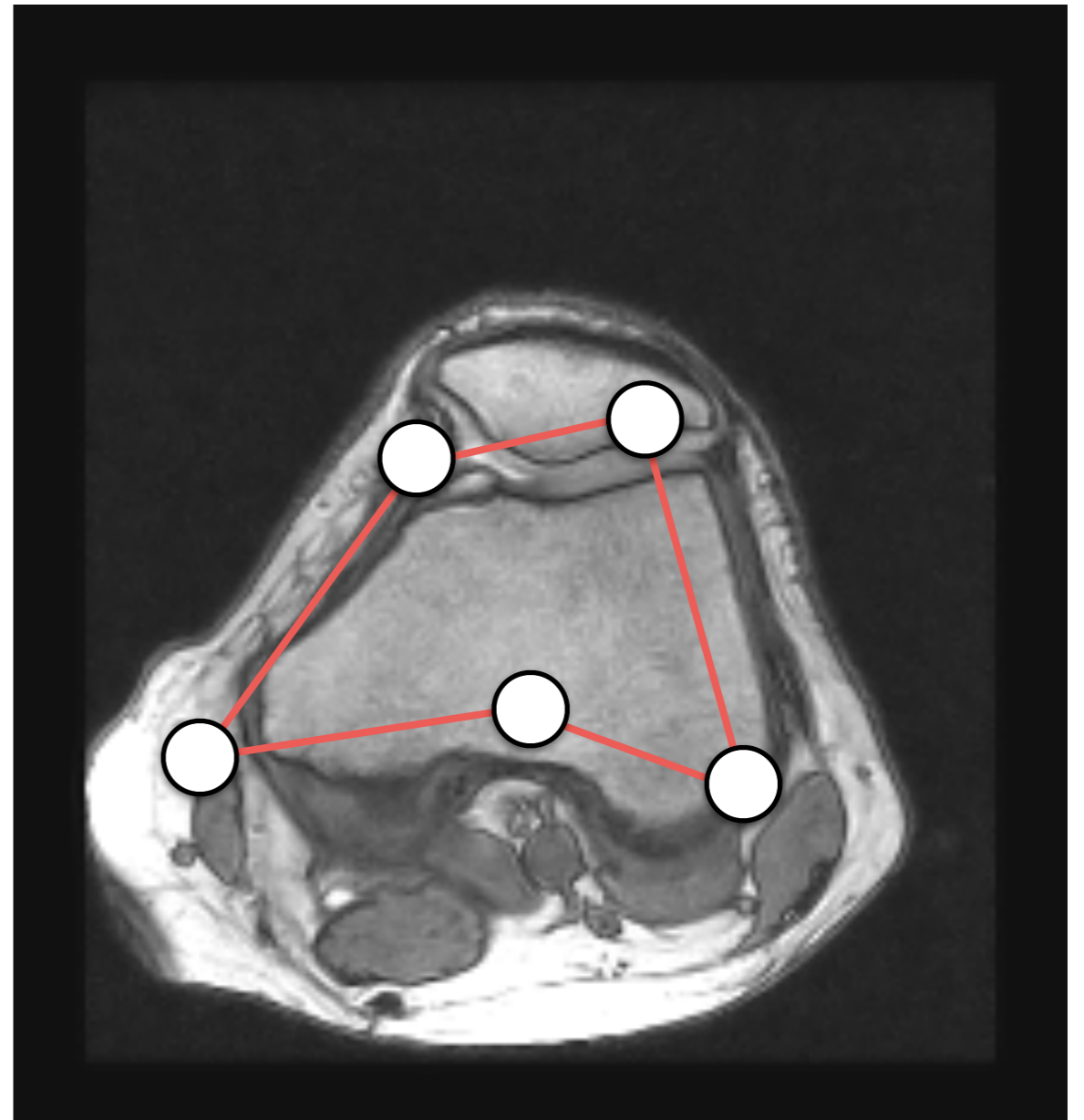
Snakes

- A first step, we draw a snake close to the boundary of the object we want to segment.



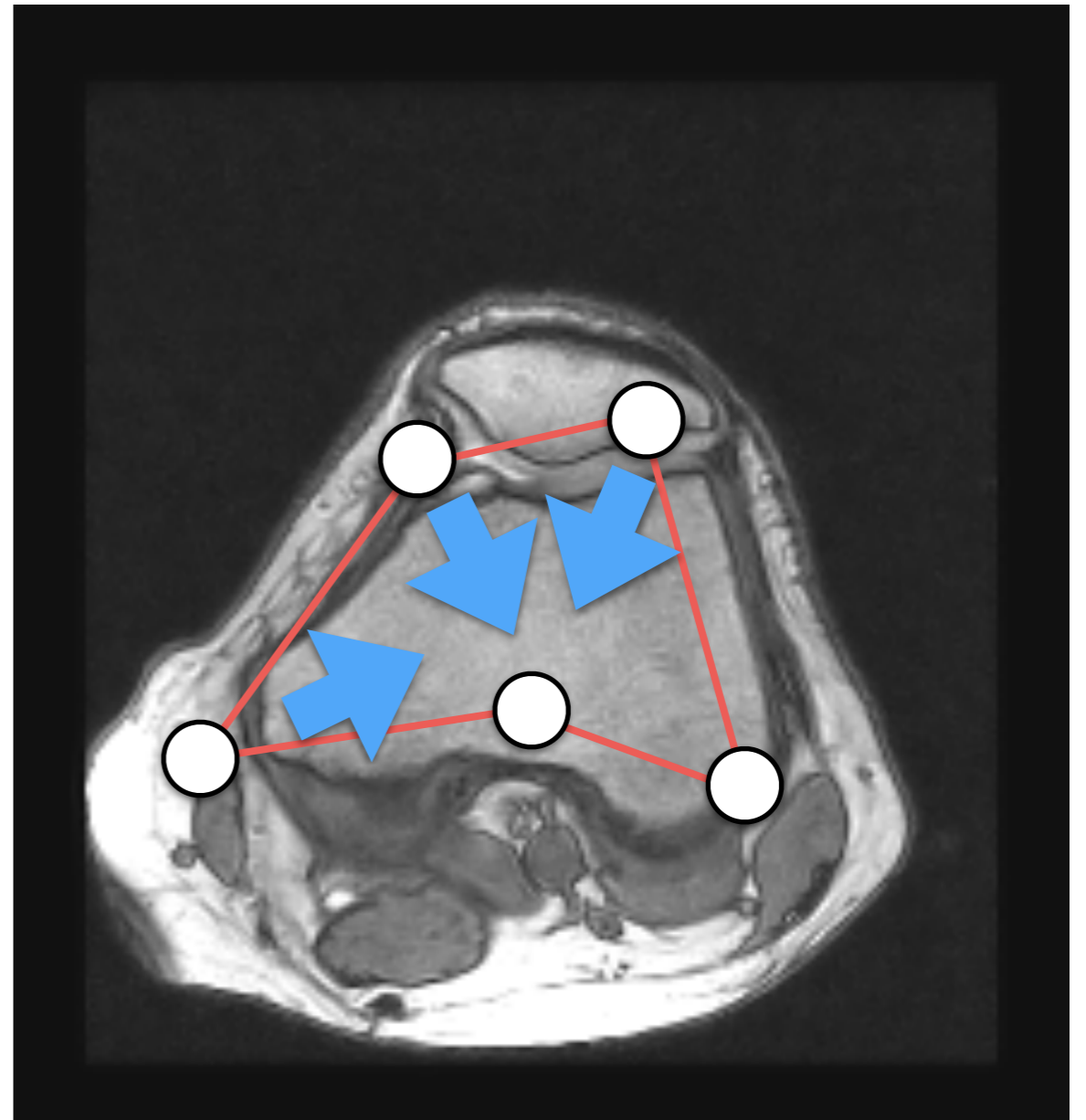
Snakes

- Then, we *deform* its control points in order to move them towards the object's boundary.



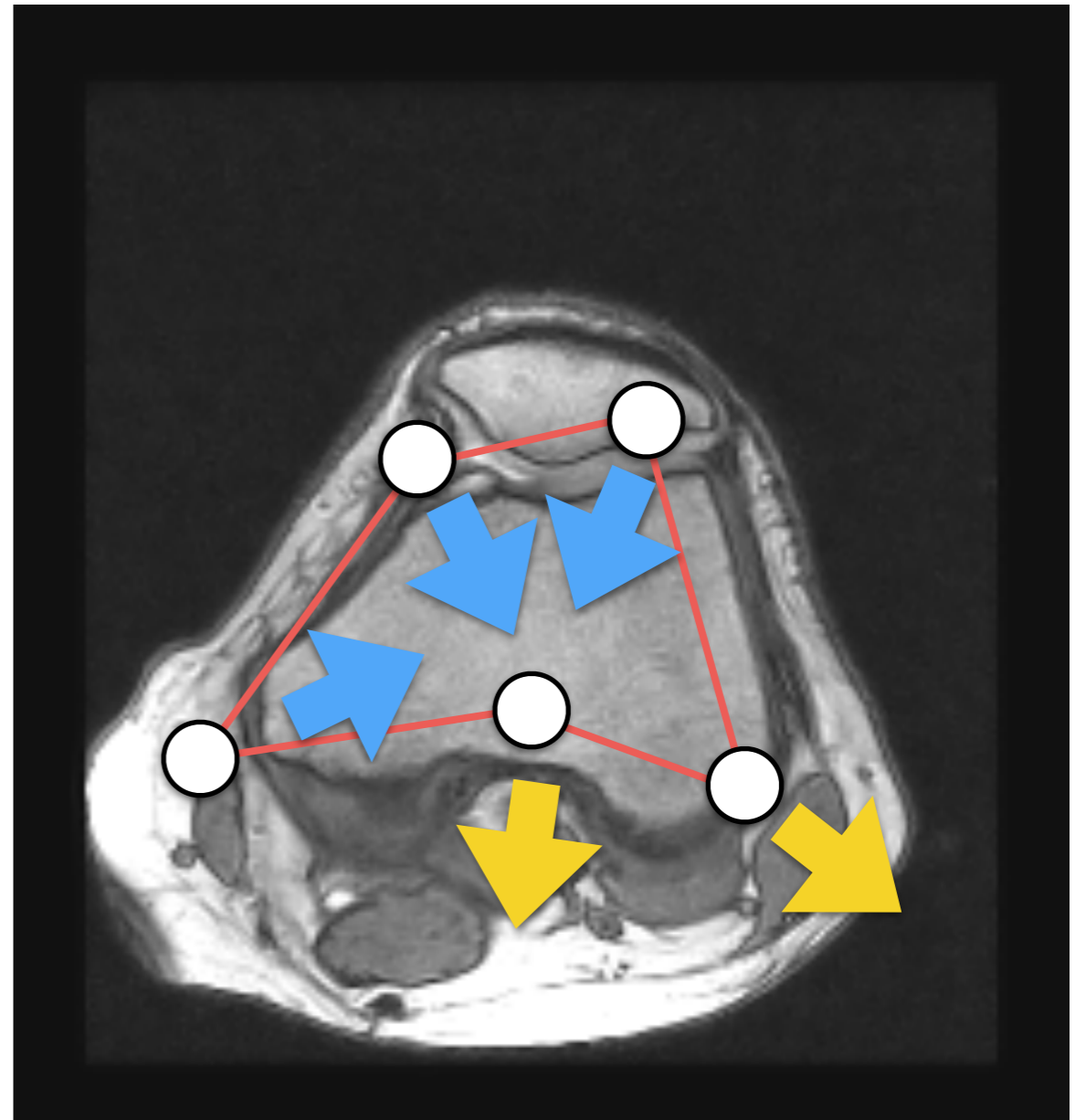
Snakes

- Then, we *deform* its control points in order to move them towards the object's boundary.



Snakes

- Then, we *deform* its control points in order to move them towards the object's boundary.



Snakes

- How do we deform the control points?
- An energy function E is associated with the curve.
- We deform control points by minimizing E ; i.e., we solve an optimization problem.

Snakes

- How do we define the energy function?
- The energy of a snake has three components:

$$E = E_{\text{internal}} + E_{\text{external}} + E_{\text{constraint}}$$

Snakes: Internal Energy

- This energy represents the internal energy of the curve due to bending. It is defined per point as

$$E_{\text{internal}}(\mathbf{v}(t)) = \frac{1}{2} \left(\alpha(t) \left| \frac{d\mathbf{v}(t)}{dt} \right|^2 + \beta(t) \left| \frac{d^2\mathbf{v}(t)}{d^2t} \right|^2 \right)$$

- The total energy is defined as

$$E_{\text{internal}} = \int_0^1 E_{\text{internal}}(\mathbf{v}(t)) dt$$

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Elasticity

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Elasticity Stiffness

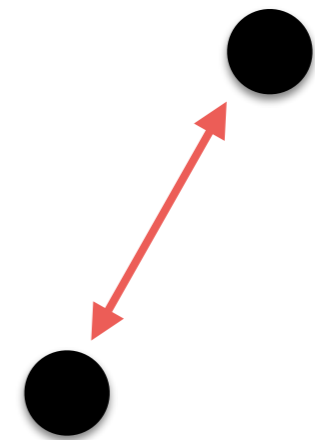
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Snakes: Internal Energy

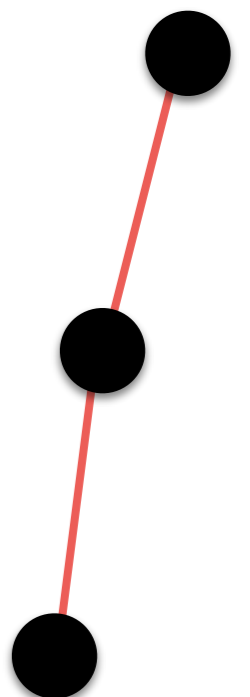
- The first term is an elastic energy:

$$\frac{d\mathbf{v}(t)}{dt} \approx \mathbf{v}_{i+1} - \mathbf{v}_i$$



- The second term is a bending energy:

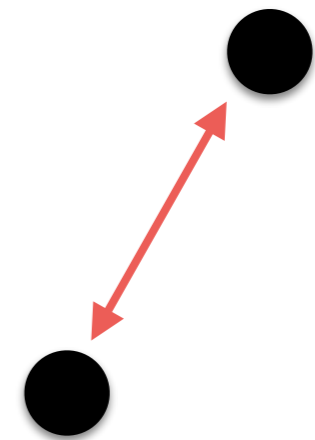
$$\frac{d^2\mathbf{v}(t)}{d^2t} \approx \mathbf{v}_{i+1} - 2\mathbf{v}_i + \mathbf{v}_{i-1}$$



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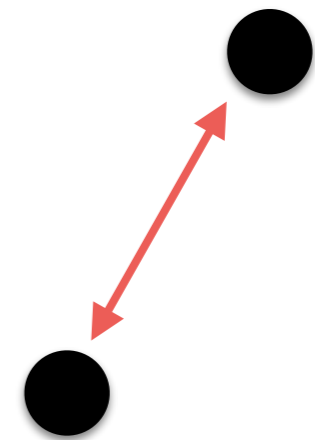
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Snakes: Internal Energy

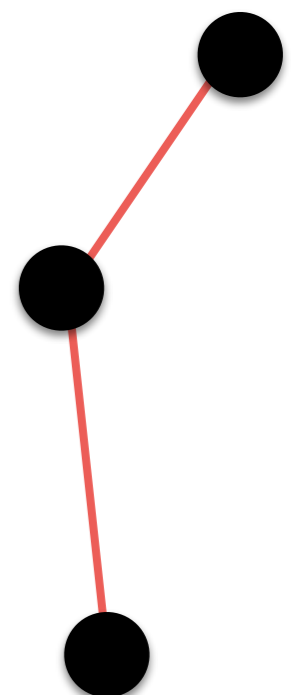
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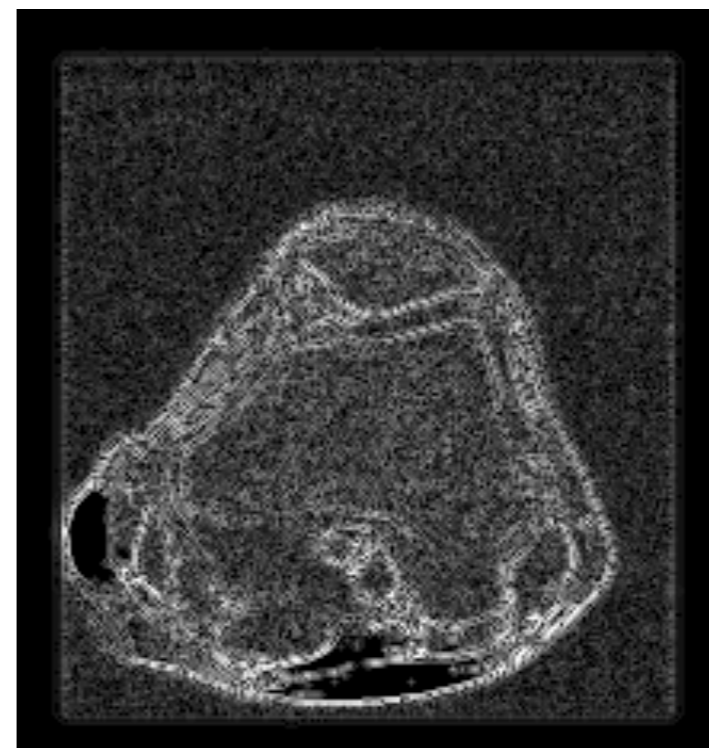
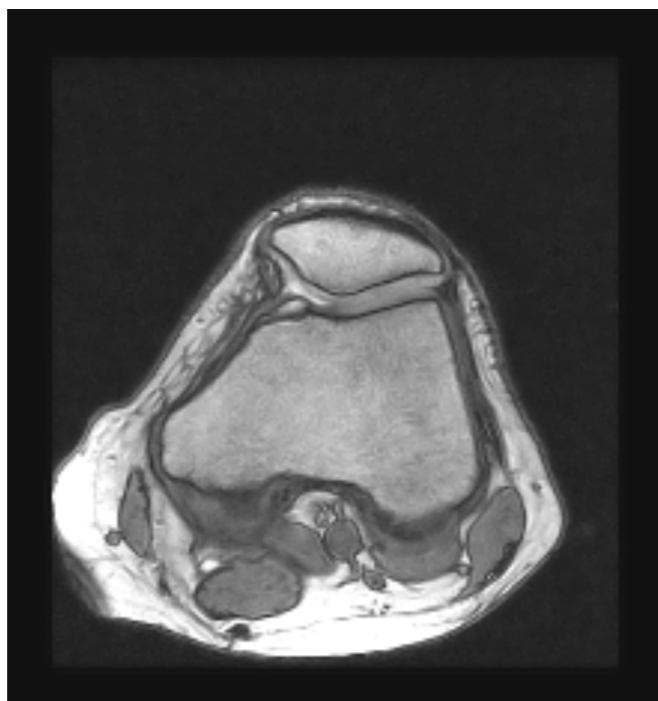
- The second term is a bending energy:

$$\frac{d^2\mathbf{v}(t)}{d^2t} \approx \mathbf{v}_{i+1} - 2\mathbf{v}_i + \mathbf{v}_{i-1}$$



Snakes: External Energy

- This energy determines how well the snake matches with the image locally!
- How can we achieve this?
 - Gradients magnitude



Snakes: External Energy

- It is defined per point as

$$E_{\text{external}}(\mathbf{v}(t)) = -\|\nabla I(\mathbf{v}(t))\|^2$$

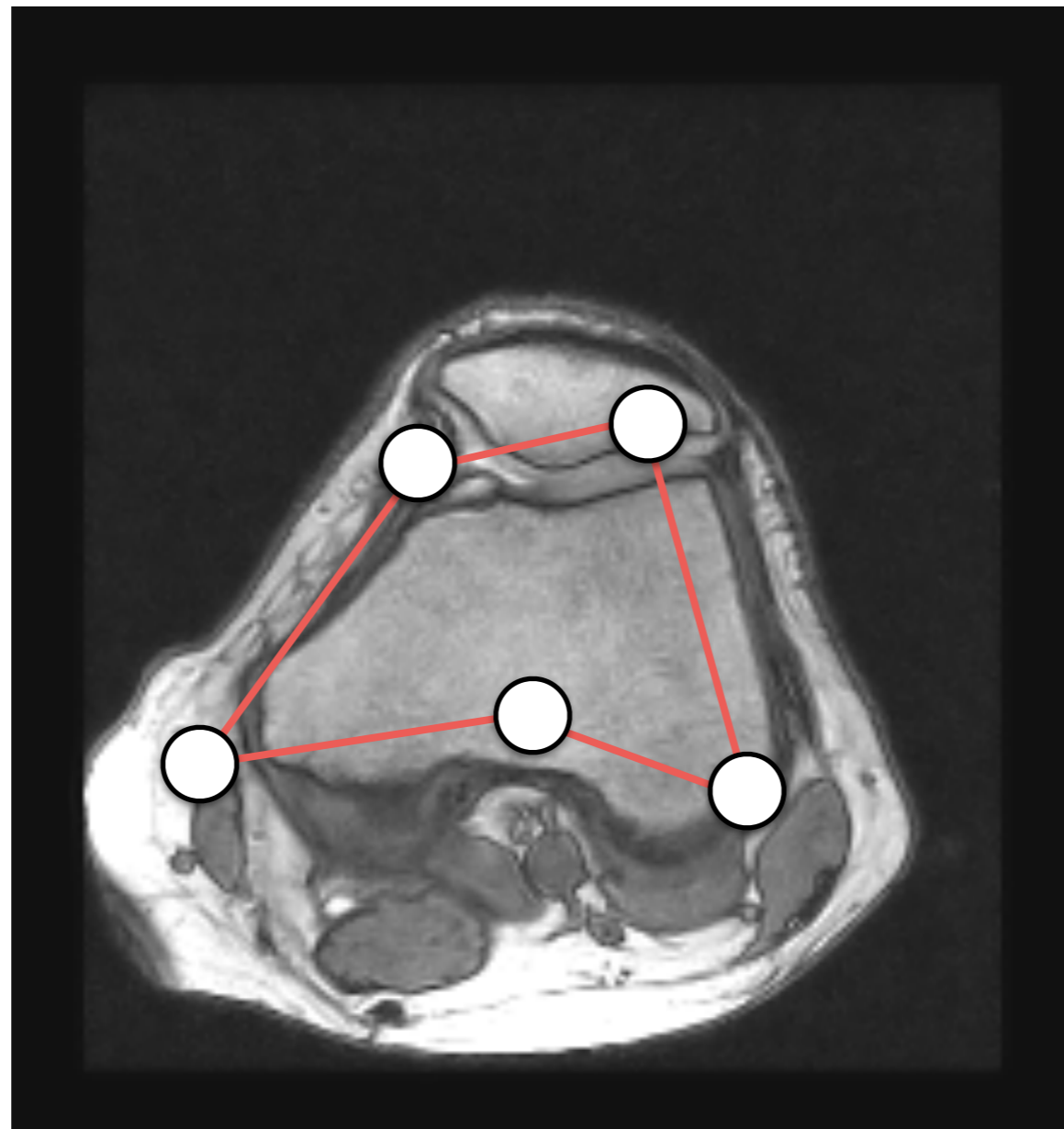
- The total energy is defined as

$$E_{\text{external}} = \int_0^1 E_{\text{external}}(\mathbf{v}(t)) dt$$

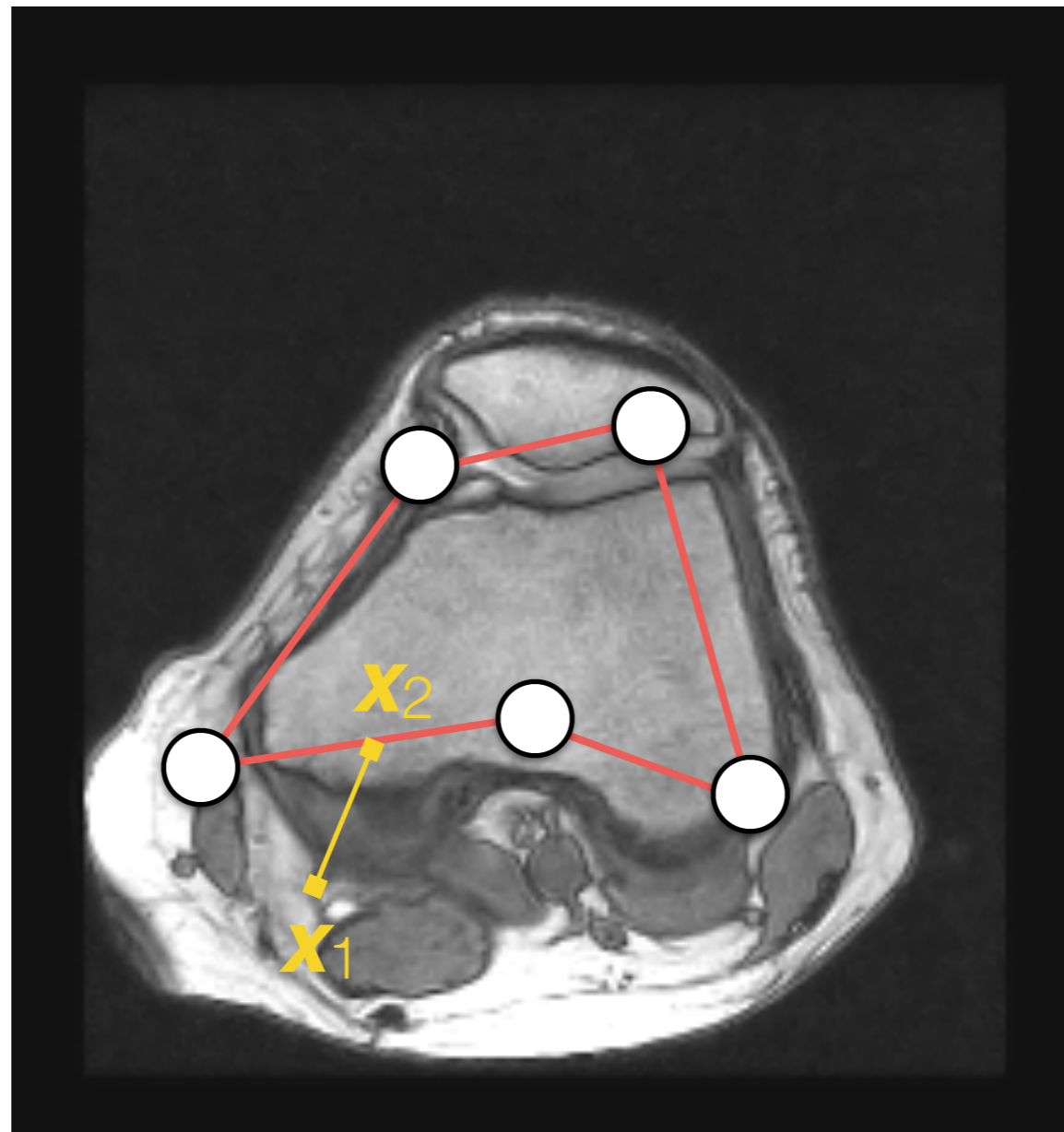
Snakes: Constraint Energy

- This energy is meant for interactive systems.
- The user interactively monitors the minimization, and she/he can push/pull vertices using the mouse cursor's position:
 - Repulsion forces or “vulcano”: $\frac{1}{r^2}$
 - Spring forces: $-k\|\mathbf{x}_1 - \mathbf{x}_2\|^2$

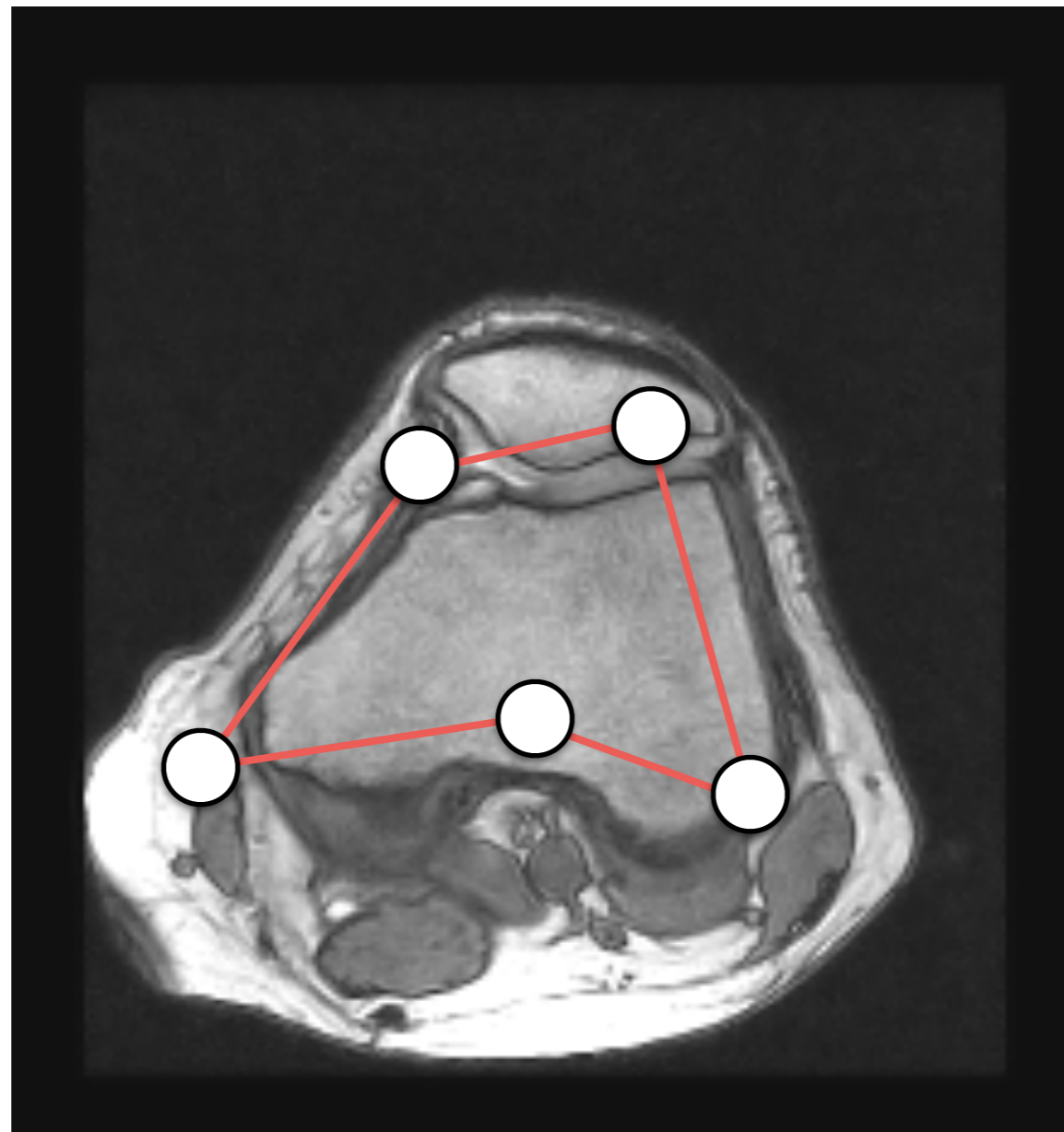
Snakes: Constraint Energy



Snakes: Constraint Energy



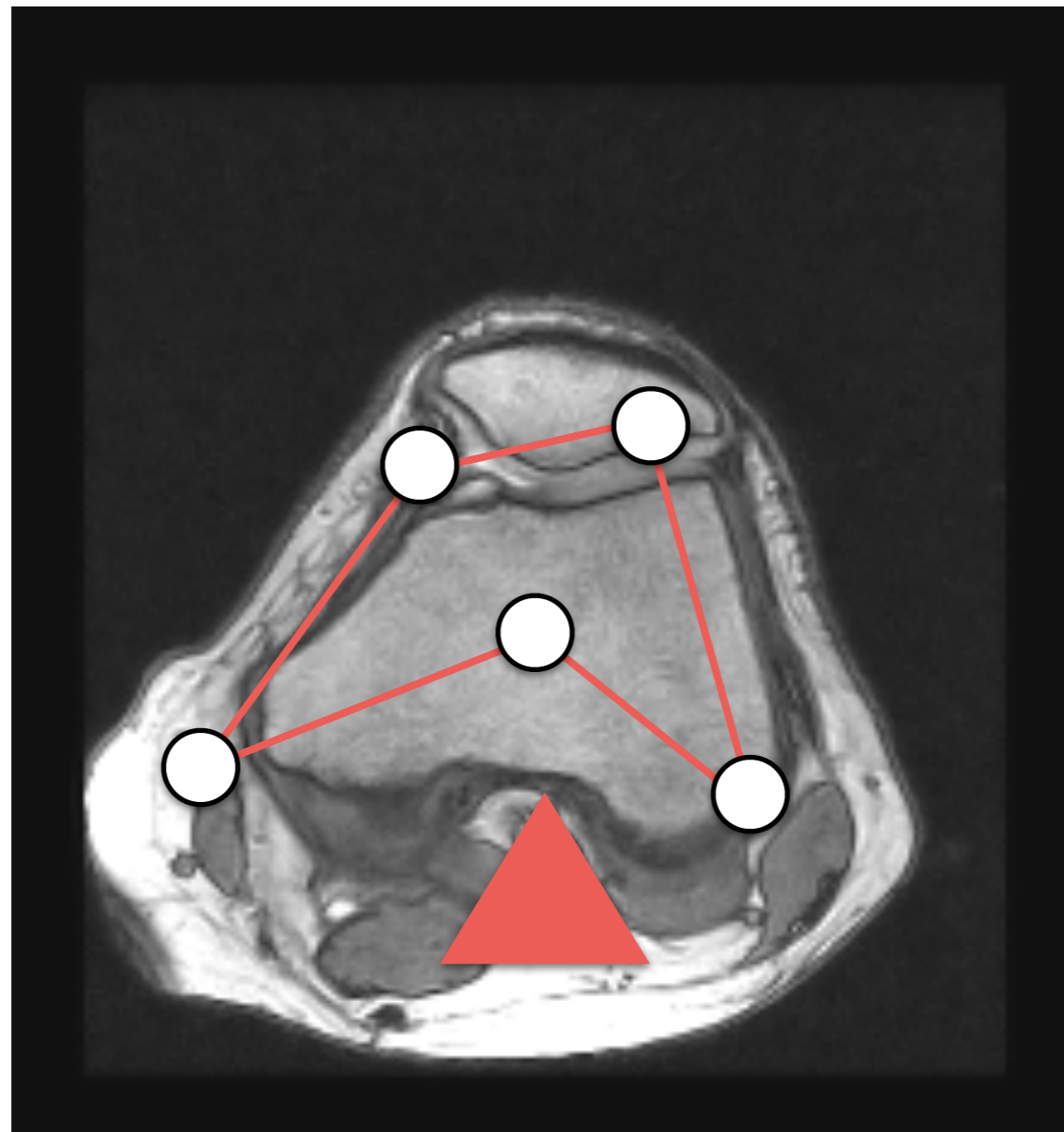
Snakes: Constraint Energy



Snakes: Constraint Energy



Snakes: Constraint Energy



How do we solve E ?

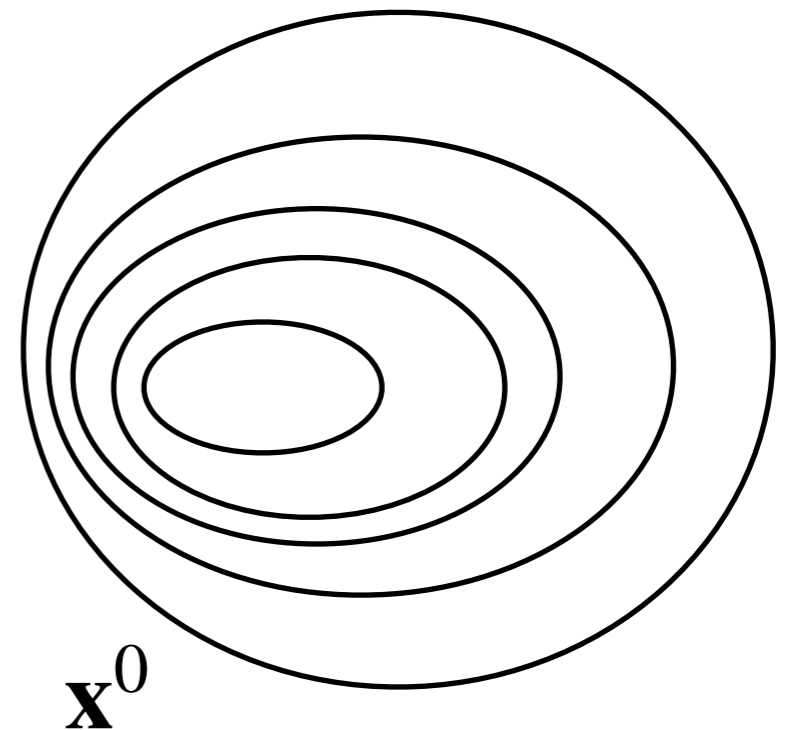
$$E = E_{\text{internal}} + E_{\text{external}} + E_{\text{constraint}}$$

Gradient Descent

- It is a first-order iterative optimization method:

$$\mathbf{x}_j^{i+1} = \mathbf{x}_j^i - \alpha \frac{\partial}{\partial \mathbf{x}_j} f(\mathbf{x}^i)$$

- \mathbf{x}^0 is a “**good**” guess.
- It will find a **local minimum**!
- f has to be differentiable.

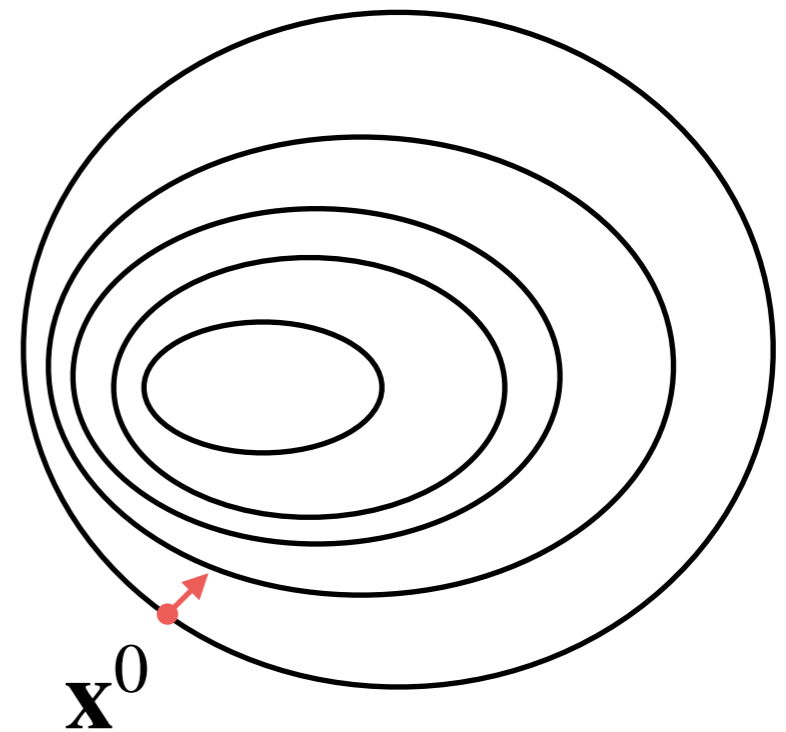


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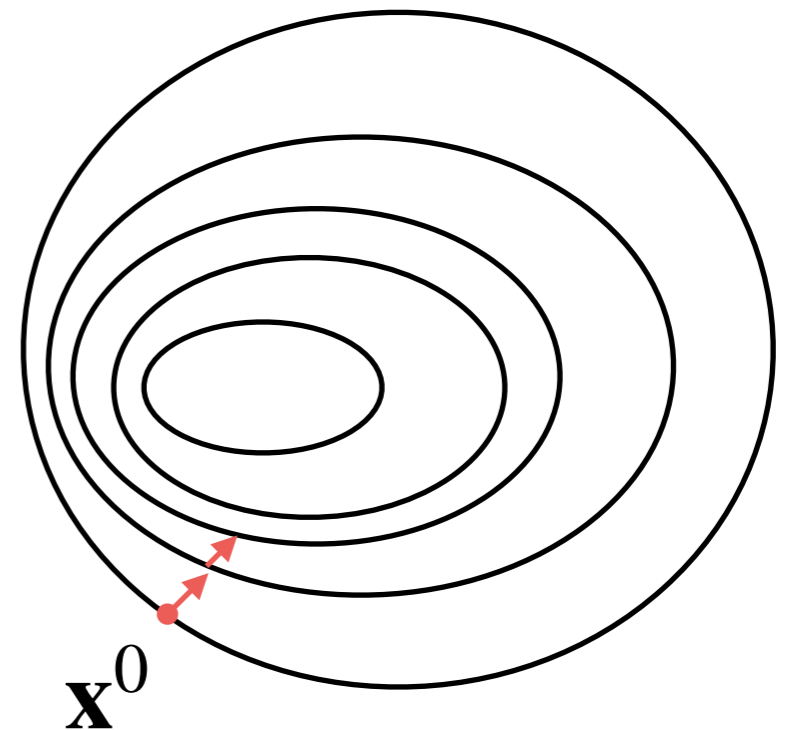


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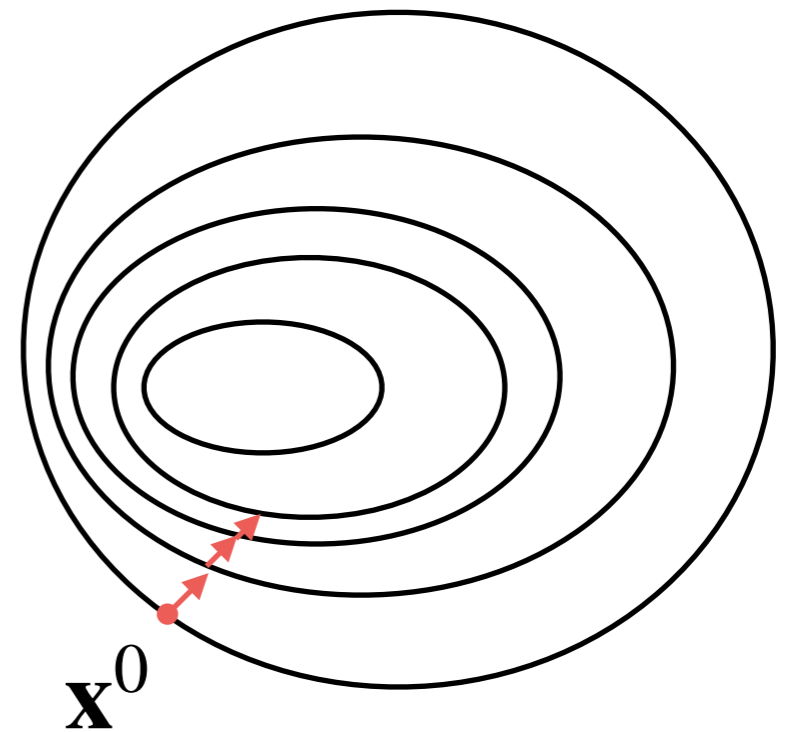


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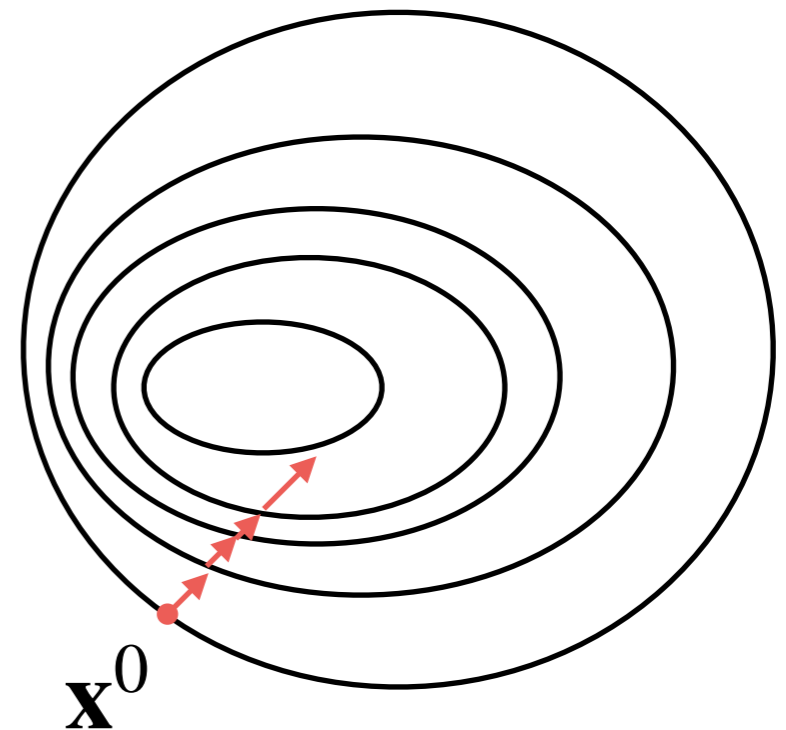


Gradient Descent

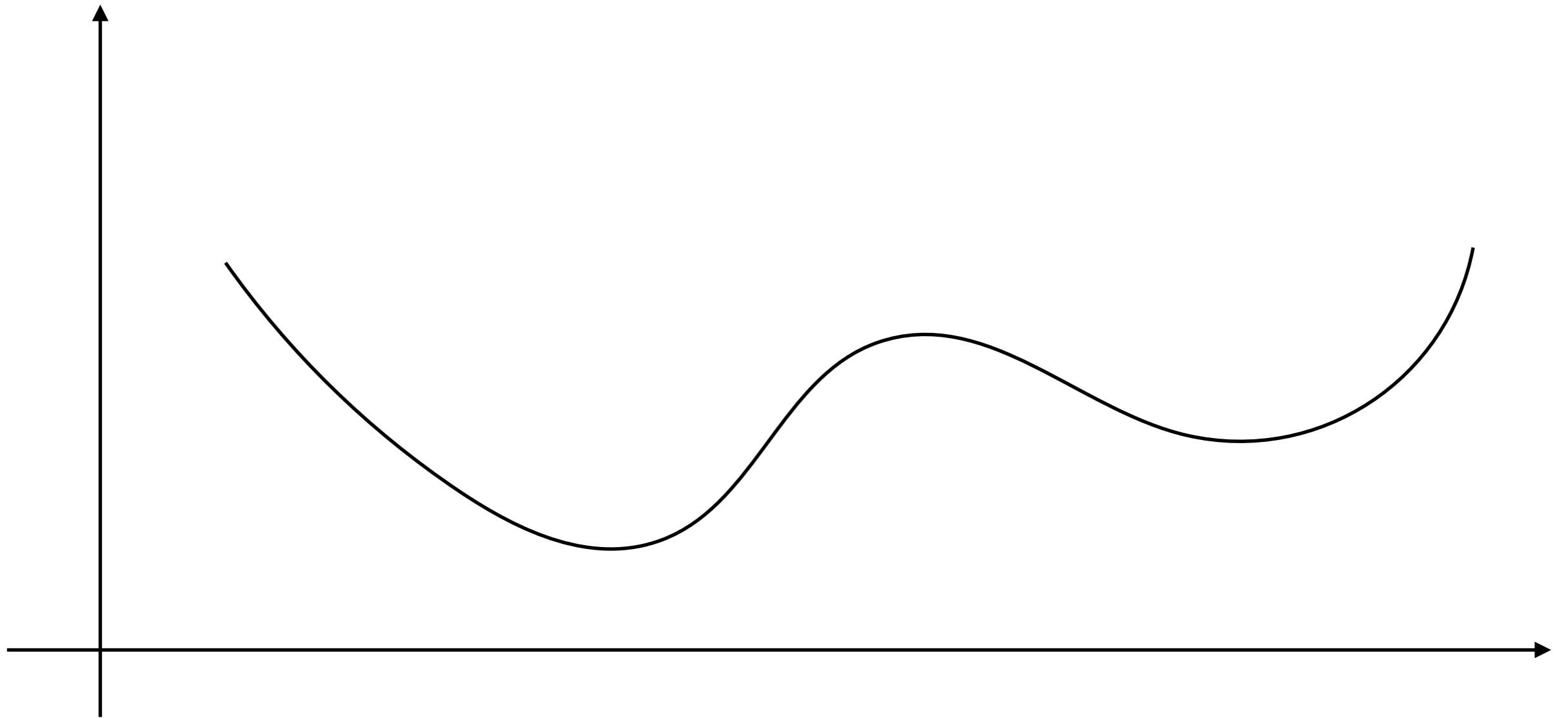
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$$\mathbf{x}_j^{i+1} = \mathbf{x}_j^i - \alpha \frac{\partial}{\partial \mathbf{x}_j} f(\mathbf{x}^i)$$

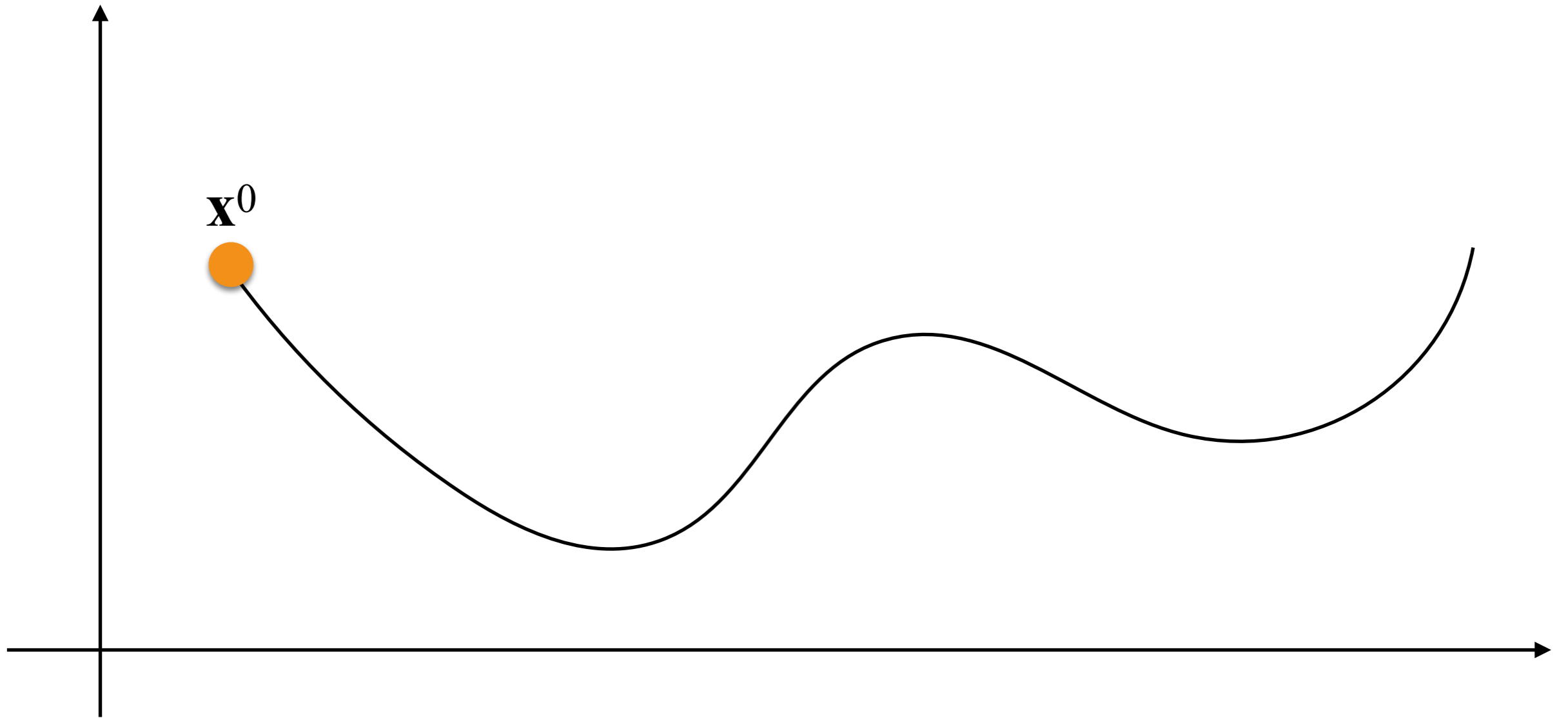
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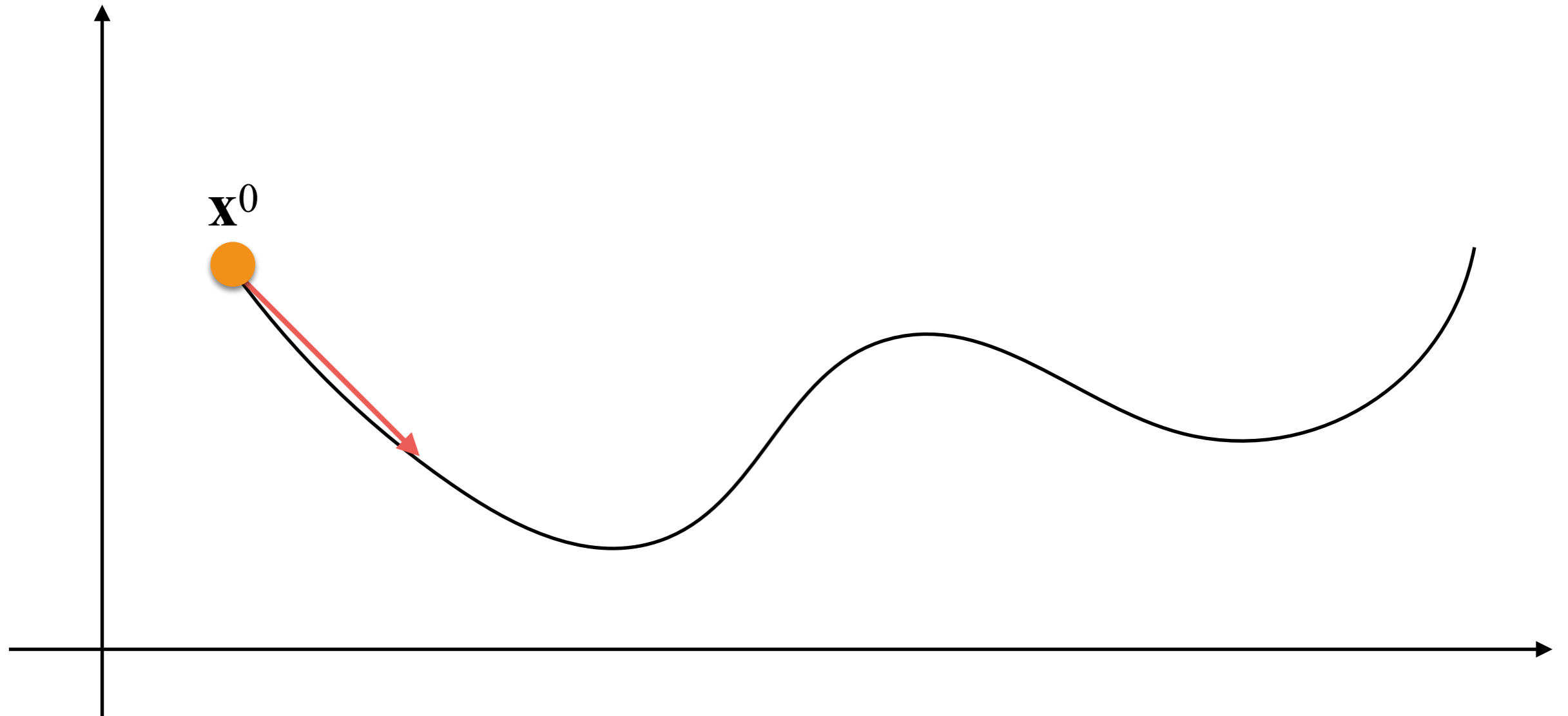
Gradient Descent



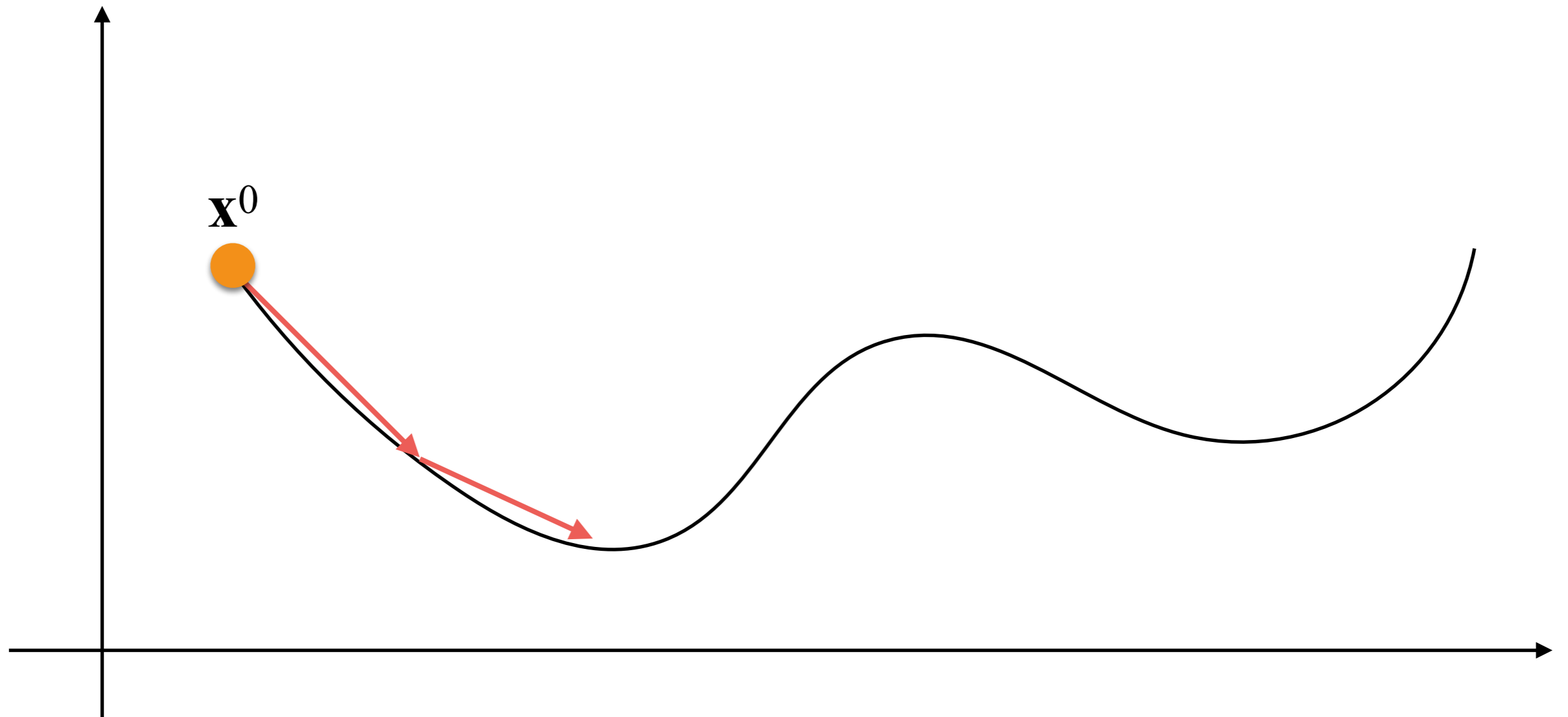
Gradient Descent



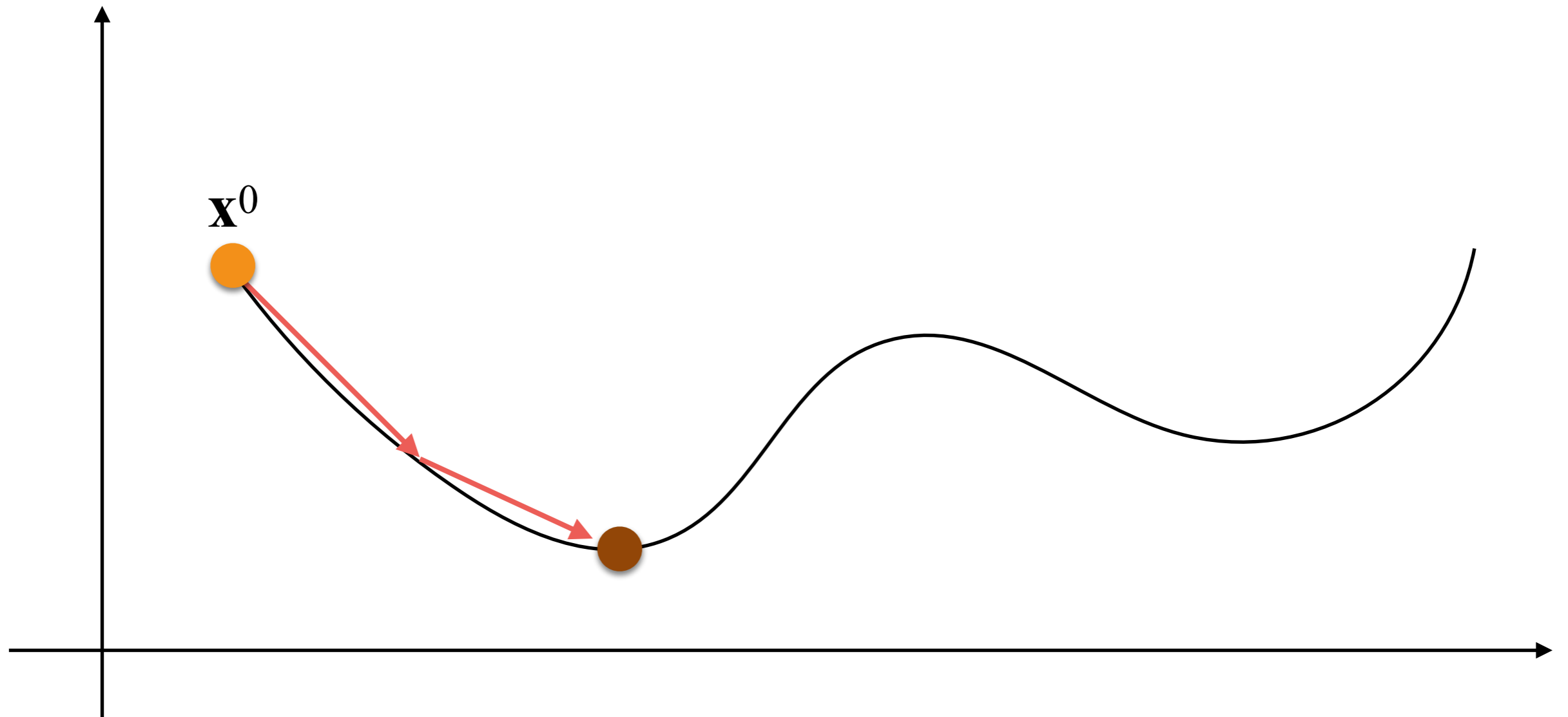
Gradient Descent



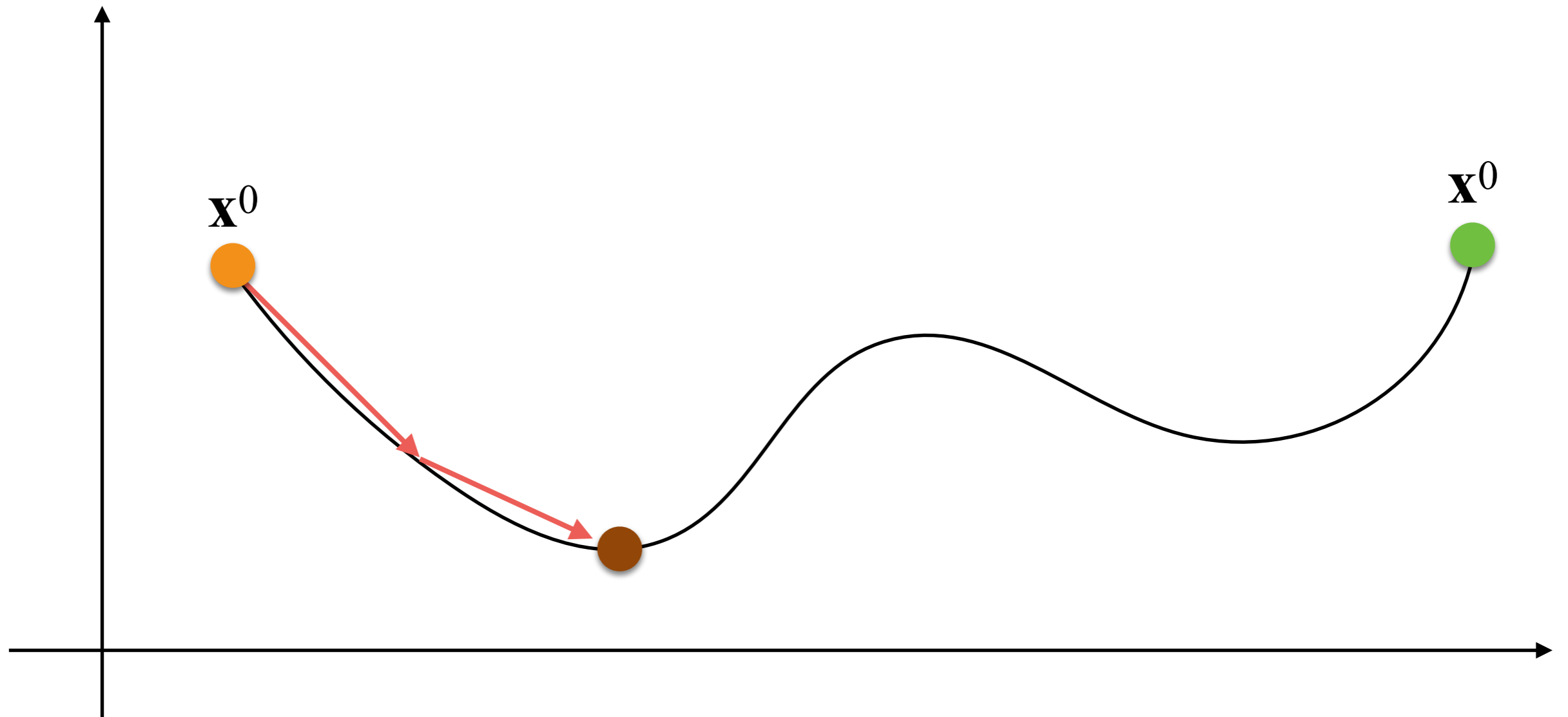
Gradient Descent



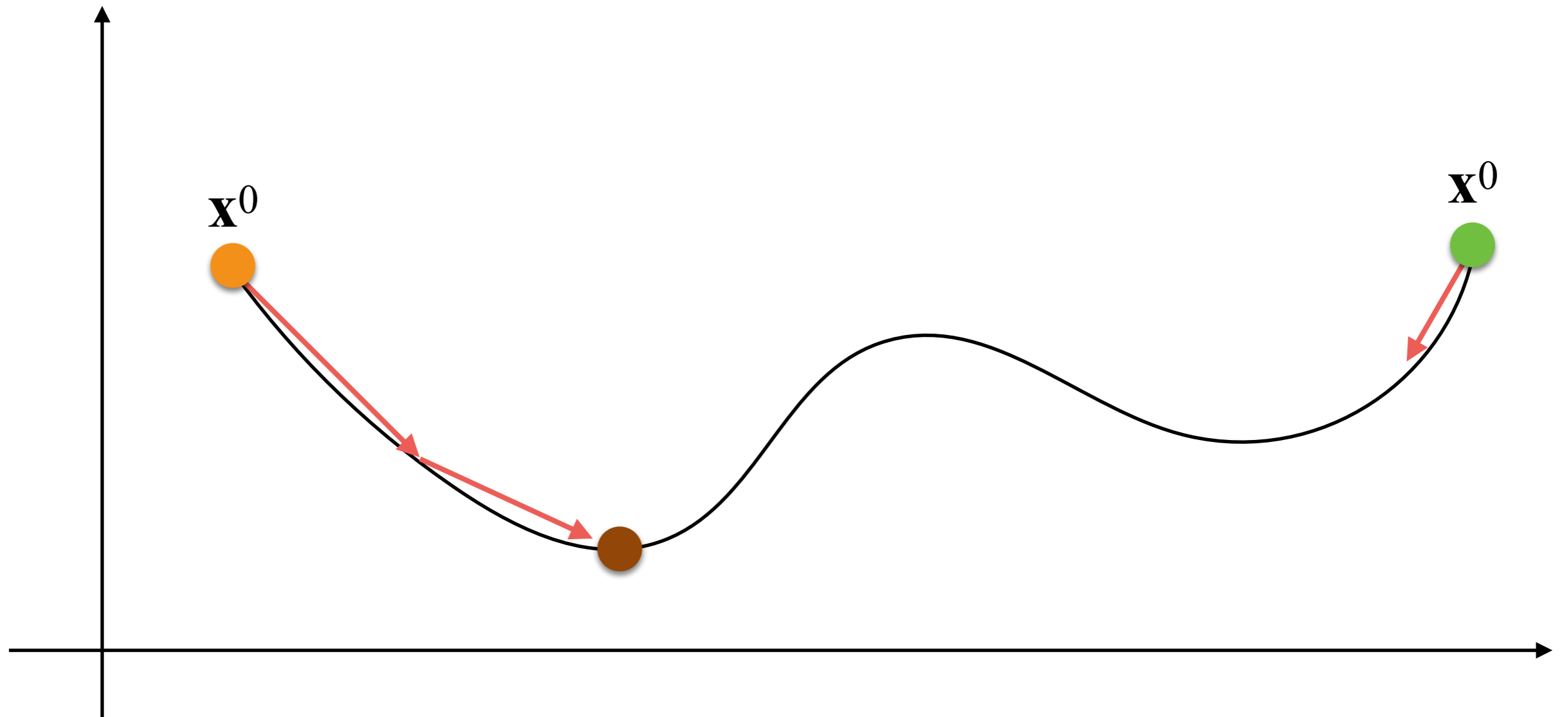
Gradient Descent



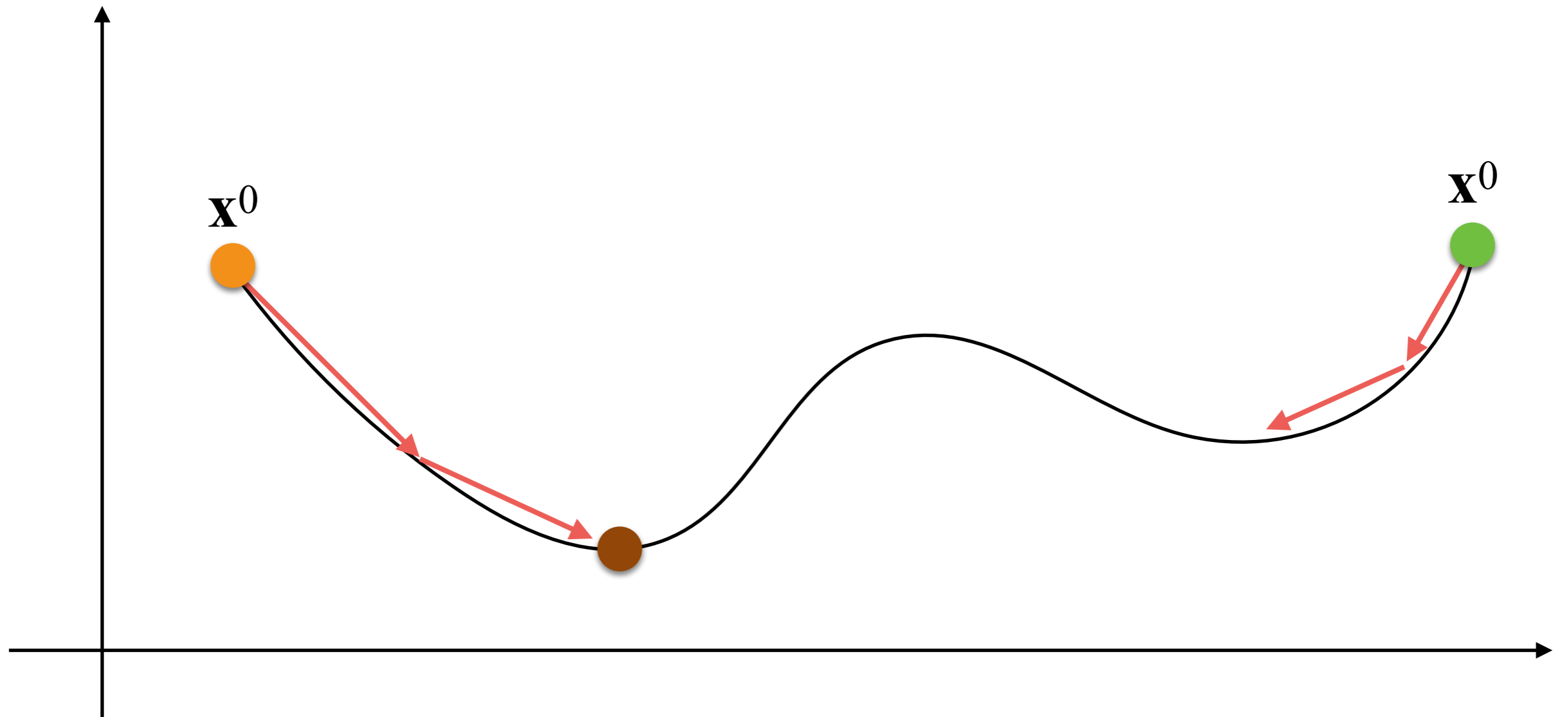
Gradient Descent



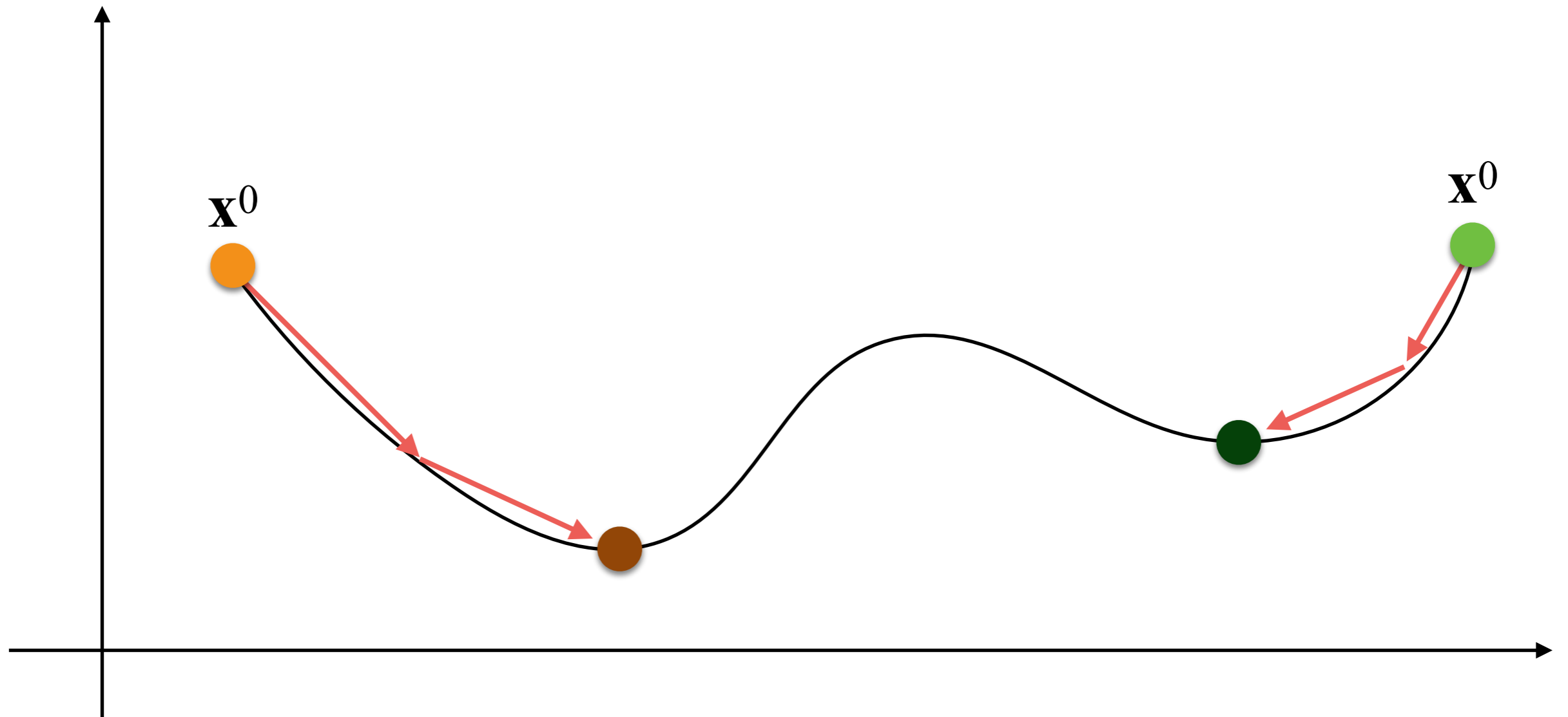
Gradient Descent



Gradient Descent



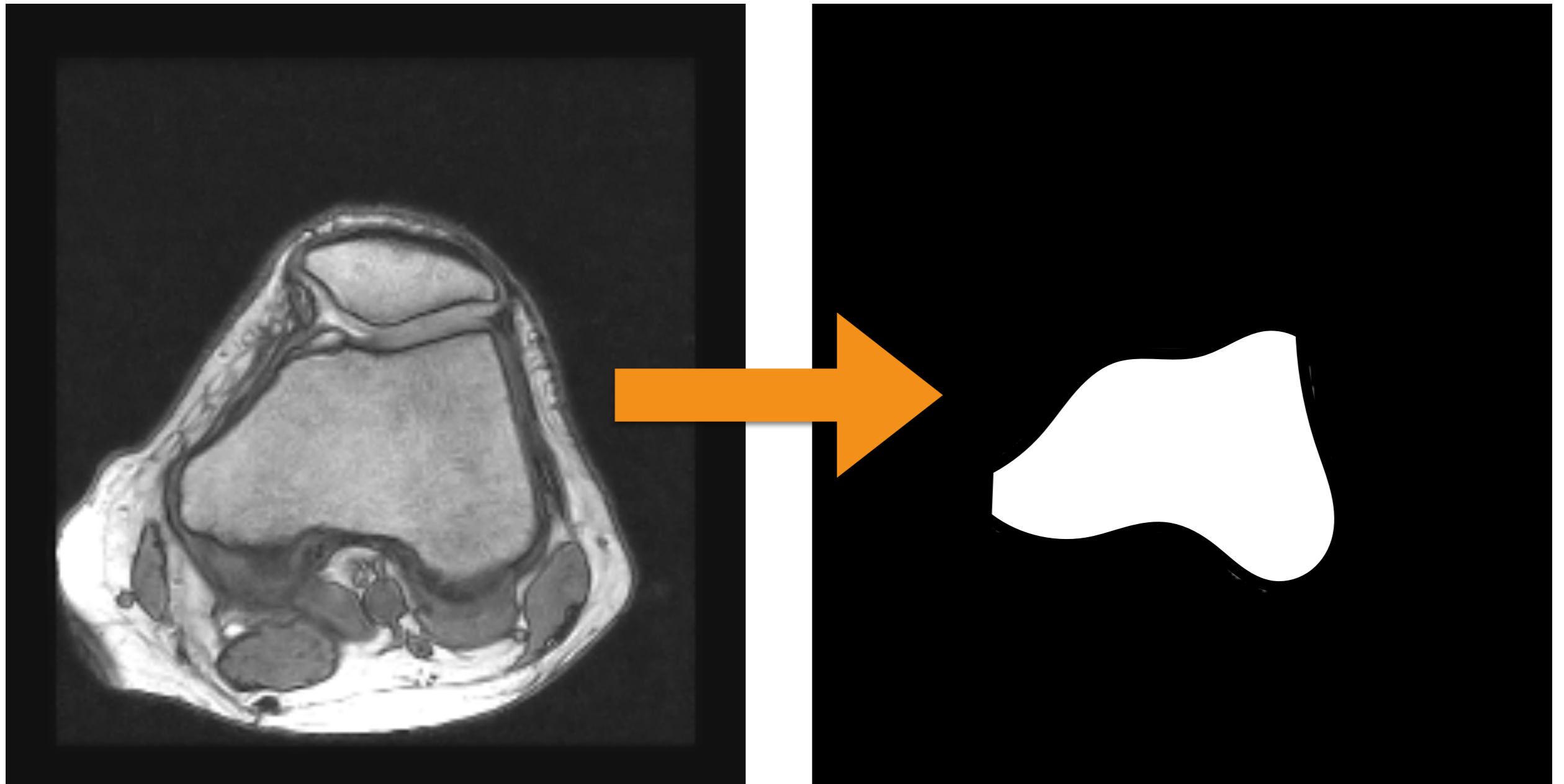
Gradient Descent



Snakes: Gradient Descent

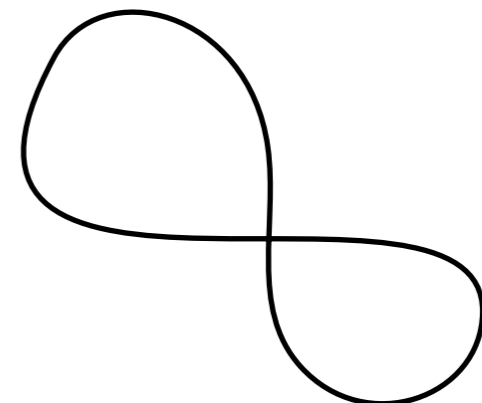
- What is our \mathbf{x}^0 in the snake minimization?
- We need to click a few points in the image around our object of interest!

Snakes An Example



Snakes

- Extension to the 3D case:
 - Instead of a curve we have a parametric surface; e.g., we can start using a sphere.
- Disadvantages:
 - We may have an over-smooth boundaries when using splines
 - How many n control points?
 - Not trivial to avoid self-intersection!



Stroke-Based

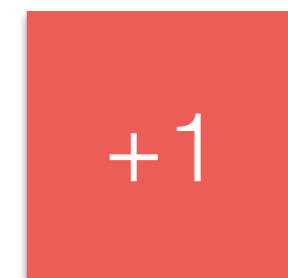
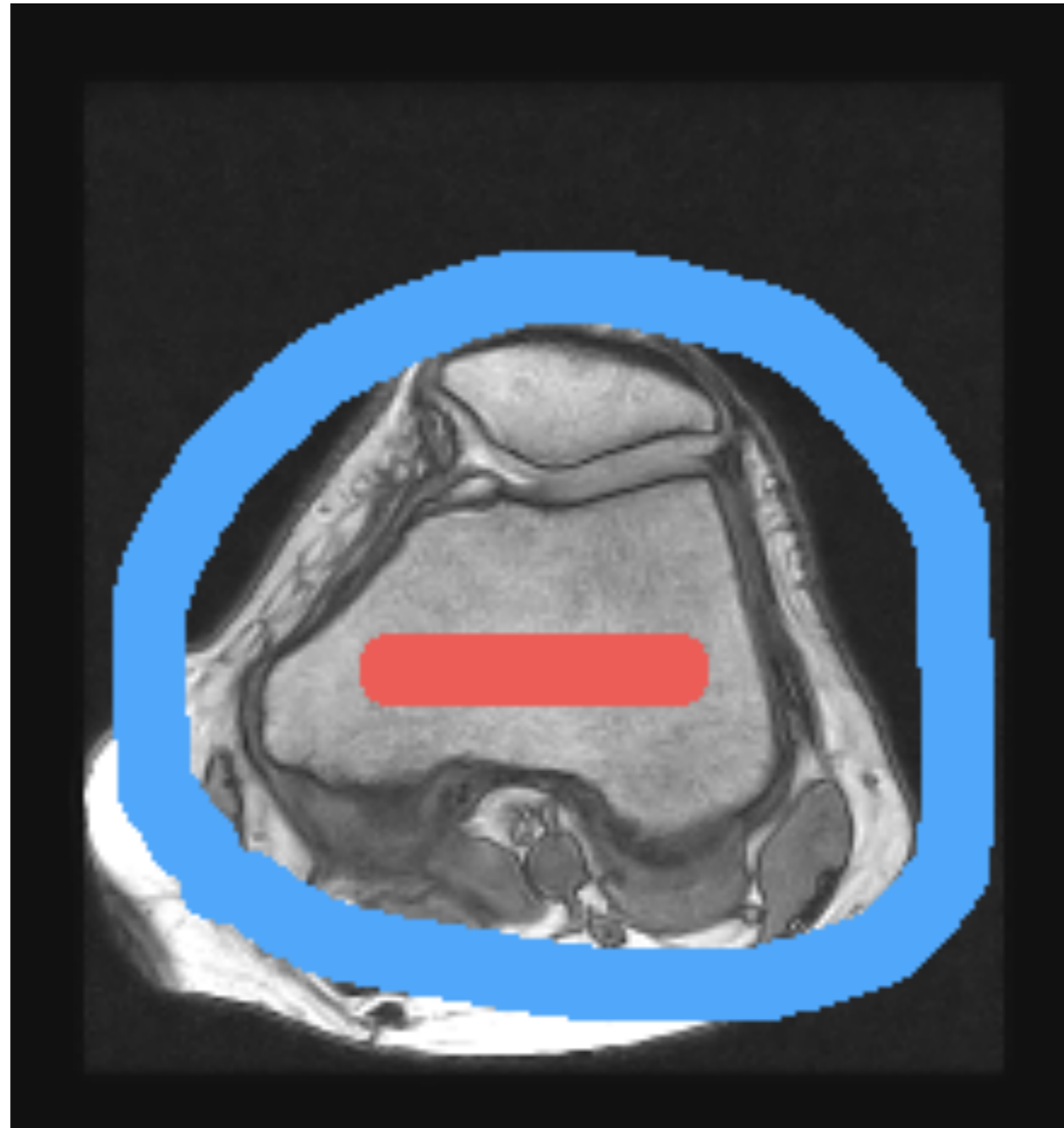
Stroke-Based

- Stroke-based algorithms are based on the idea to define with a stroke what is foreground (i.e., our object of interest) and what is background.
- These strokes are roughly painted.
 - However, they have to be placed in areas where we are 100% sure how to classify the image.

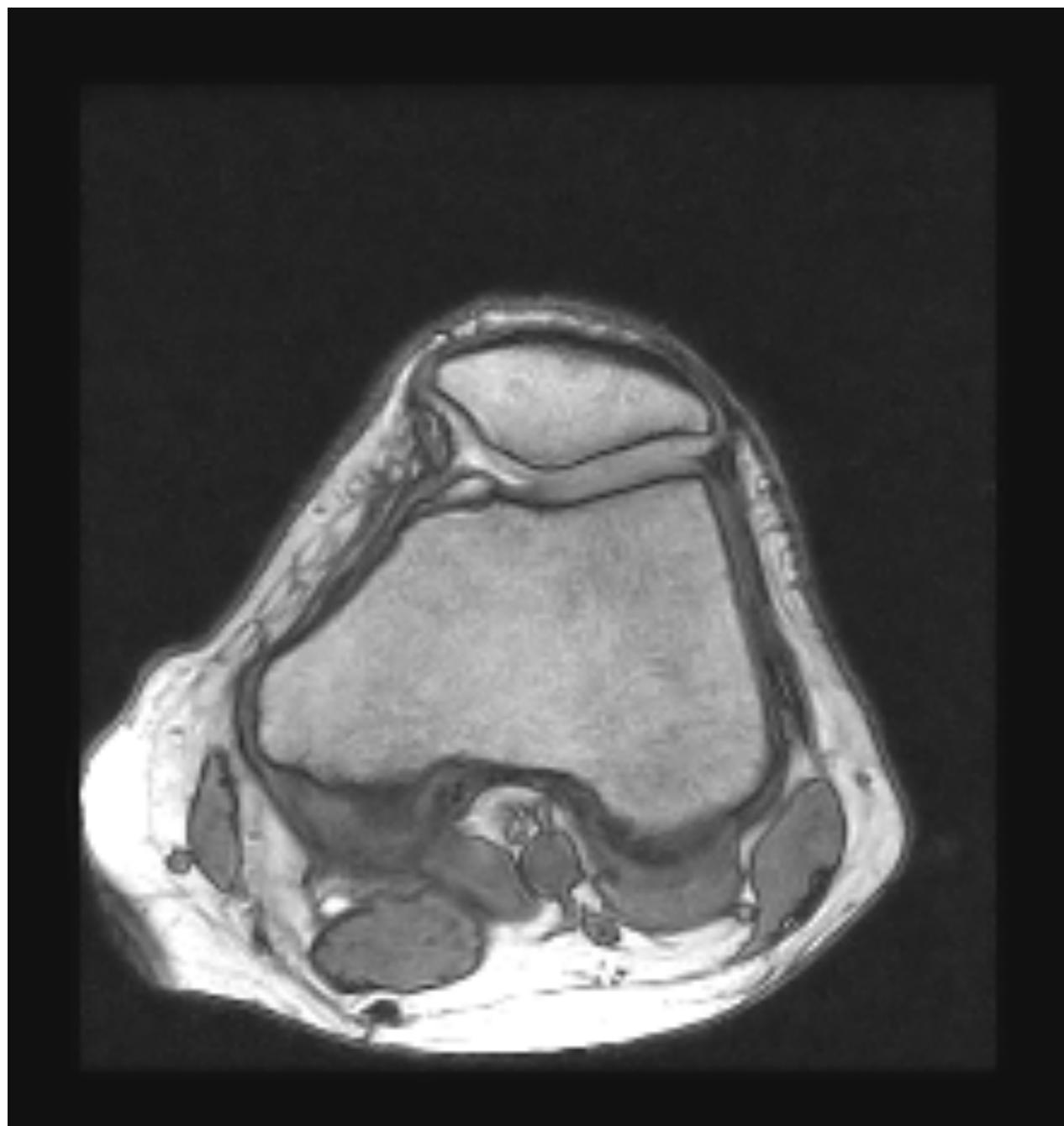
Stroke-Based



Stroke-Based



Stroke-Based



C



S

Stroke-Based: Grow-Cut

- Grow-cut is a stroke-based method.
- The idea is to propagate the label of the current pixels if its neighbors are “similar”.

Stroke-Based: Grow-Cut

- For each pixel, we have:

$$\langle l_i; \theta_i; C_i \rangle$$

- Initialization for pixels **not covered** by a stroke (s):

$$\langle l_i = 0; \theta_i = 0; C_i = I(x_i, y_i) \rangle \quad \forall_i s(x_i, y_i) = 0$$

- Initialization for pixels **covered** by a stroke (s):

$$\langle l_i = s(x_i, y_i); \theta_i = 1; C_i = I(x_i, y_i) \rangle \quad \forall_i s(x_i, y_i) \neq 0$$

Stroke-Based: Grow-Cut

- For each pixel, we have:

$$\text{Label} \langle \overset{\text{Strength}}{\boxed{l_i; \theta_i; C_i}} \rangle \text{ Intensity}$$

- Initialization for pixels **not covered** by a stroke (s):

$$\langle l_i = 0; \theta_i = 0; C_i = I(x_i, y_i) \rangle \quad \forall_i s(x_i, y_i) = 0$$

- Initialization for pixels **covered** by a stroke (s):

$$\langle l_i = s(x_i, y_i); \theta_i = 1; C_i = I(x_i, y_i) \rangle \quad \forall_i s(x_i, y_i) \neq 0$$

Stroke-Based: A Single Grow-Cut Pass

- For each pixel I in the image:

- We copy the previous status:

$$\langle l_i^{t+1}, \theta_i^{t+1}, C_i^{t+1} \rangle = \langle l_i^t, \theta_i^t, C_i^t \rangle$$

- For each neighbor j of i :

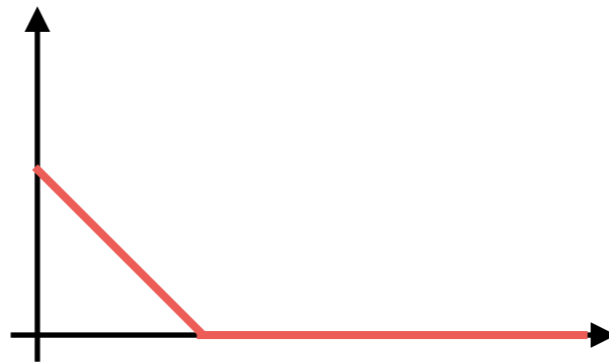
- if $g(C_i^t - C_j^t) \cdot \theta_j^t > \theta_i^t$ then

$$l_i^{t+1} = l_j^t$$
$$\theta_i^{t+1} = g(\|C_i^t - C_j^t\|_2) \cdot \theta_j^t$$

Stroke-Based: A Single Grow-Cut Pass

- Note that g is a decreasing function. For example:

$$g(x) = \begin{cases} 1 - |x| & \text{if } |x| \leq 1 \\ 0 & \text{otherwise} \end{cases}$$



- This means that if the two pixels, which we compare, are close in intensity/color values they should have the same label l .
- They should also share the same label the neighbors have a higher strength!

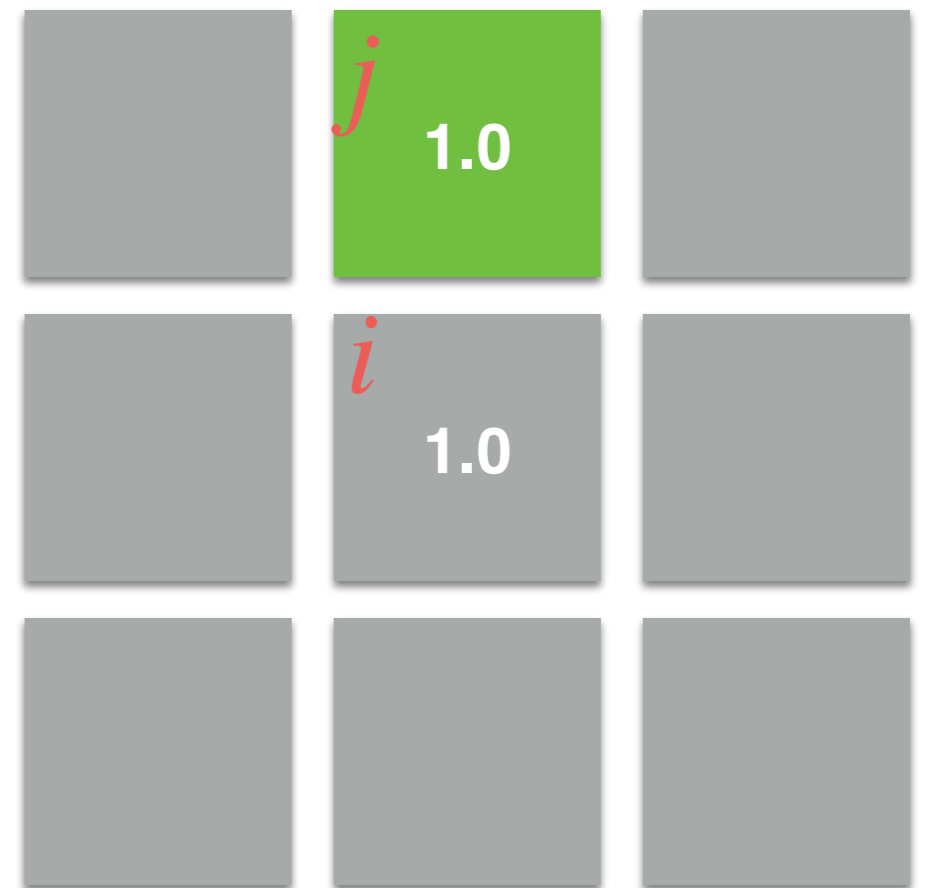
Example 1 Flat Area: Switching Labels

Stroke-Based: Example 1 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



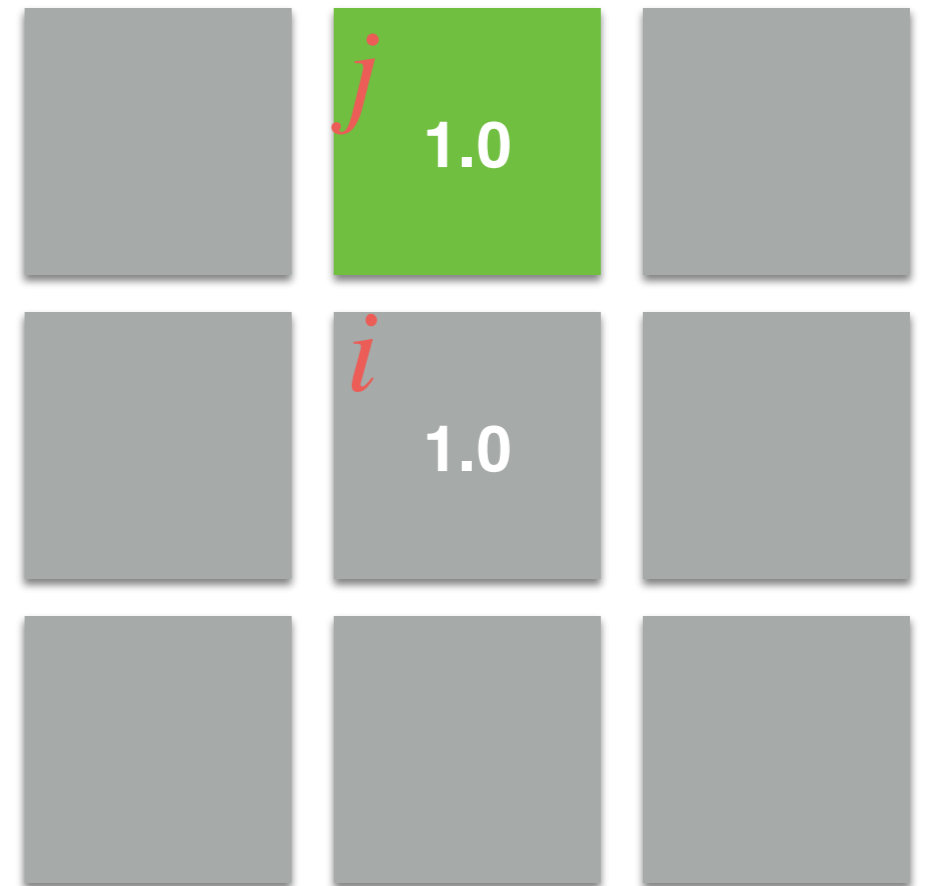
$$g(C_i^t - C_j^t) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 1 - Flat Areas

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neighbor

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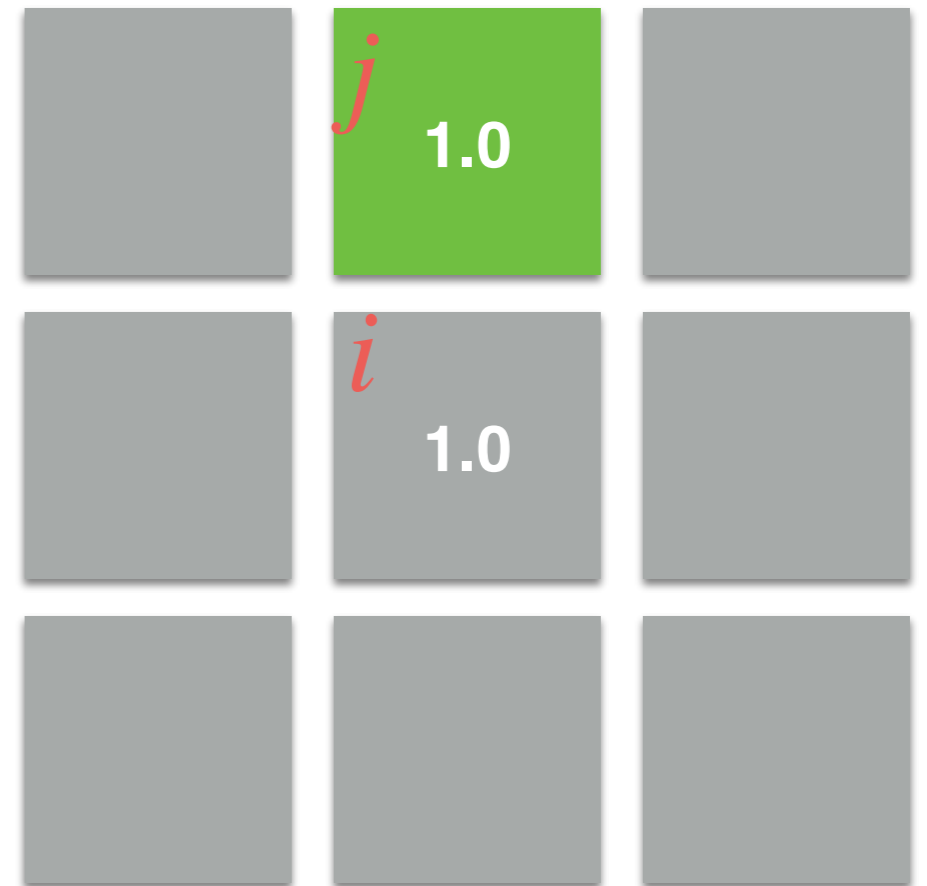
$$g(1.0 - 1.0) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 1 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



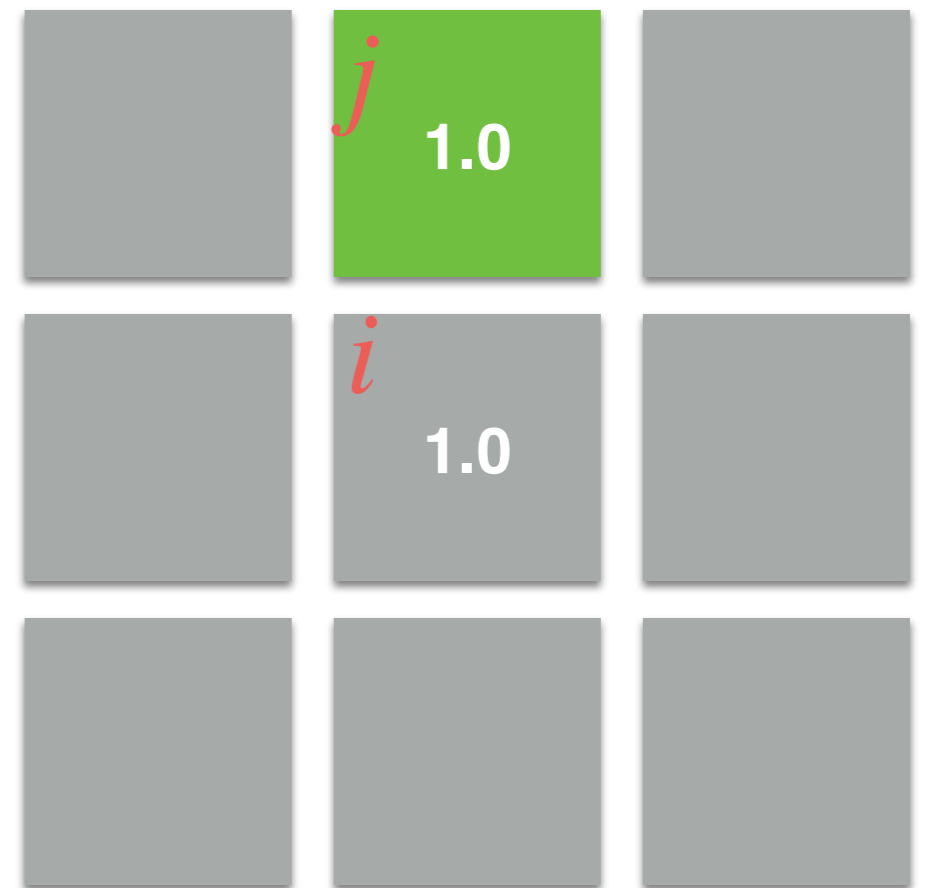
$$g(0.0) \cdot \theta_j^t > \theta_i^t$$

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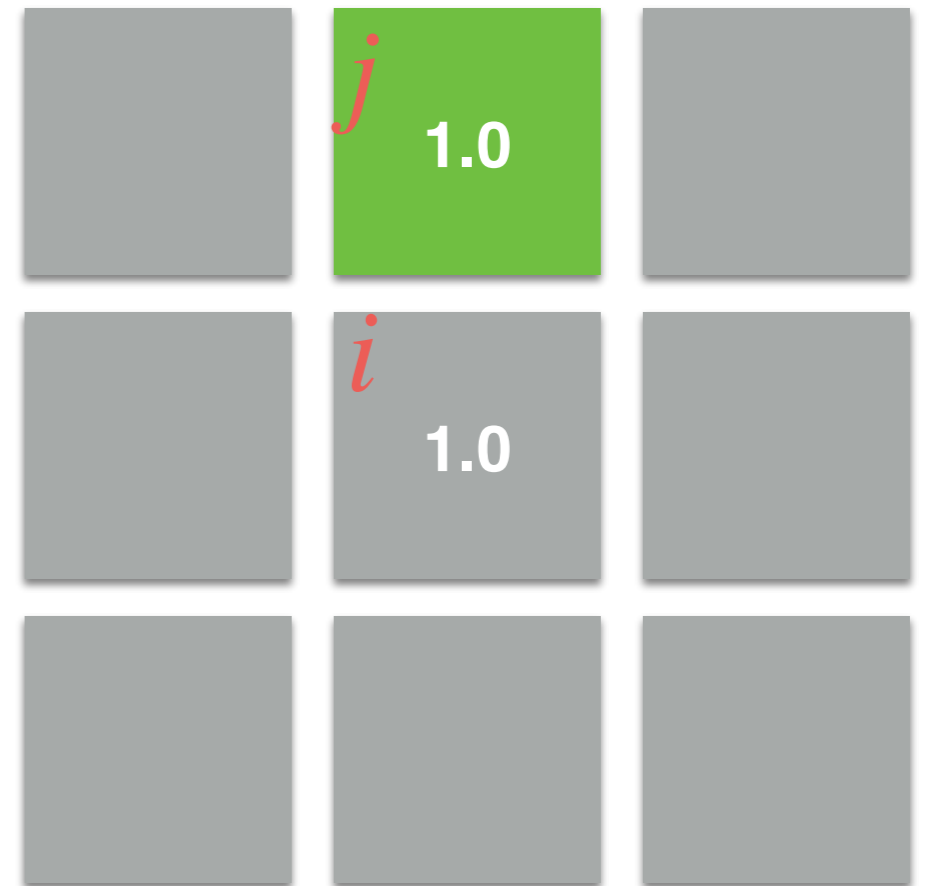
$$1.0 \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 1 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



$$0.9 > 0.8$$

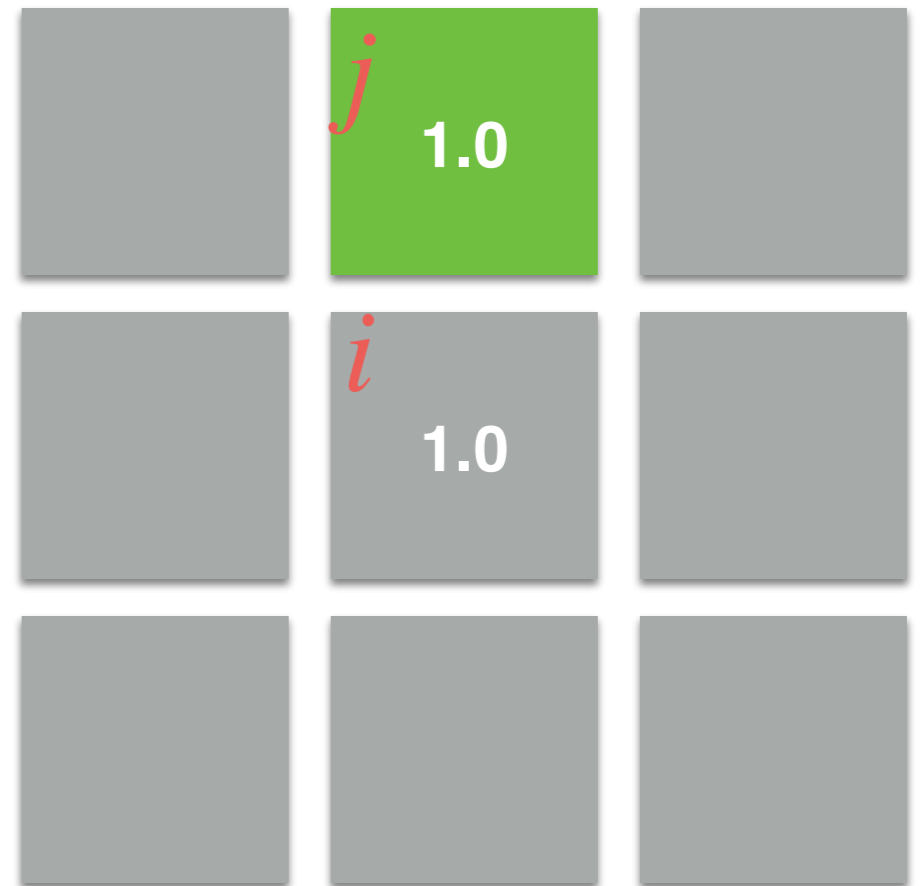
Stroke-Based: Example 1 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$

0.9 is greater than 0.8. So, so we assign
the value l_j^t to l_i^{t+1} .



$$0.9 > 0.8$$

$$l_i^{t+1} = l_j^t = 0!$$

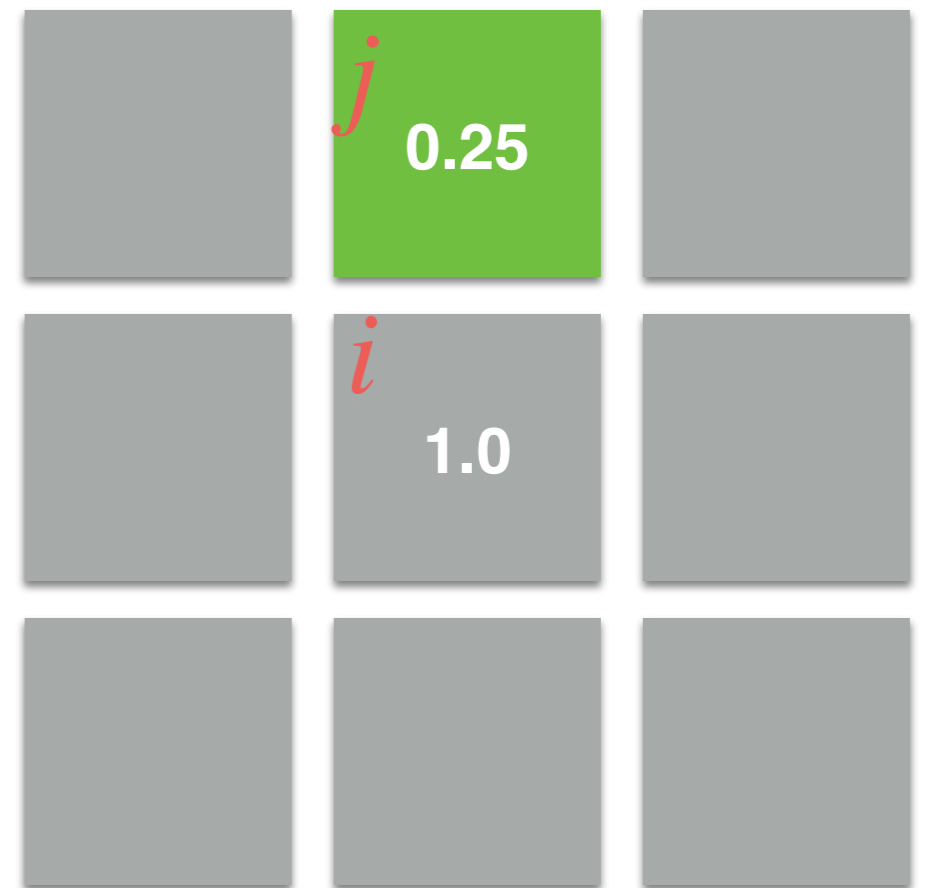
Example 2 Edge:
Maintaining Labels

Stroke-Based: Example 2 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



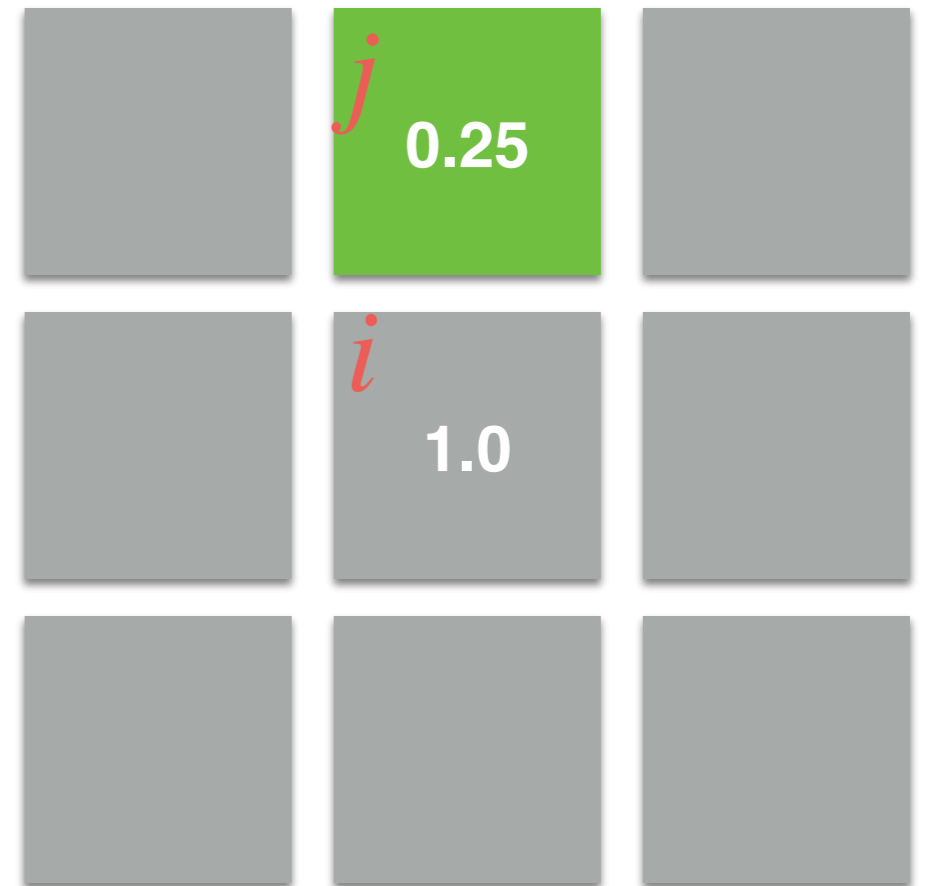
$$g(C_i^t - C_j^t) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 2 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



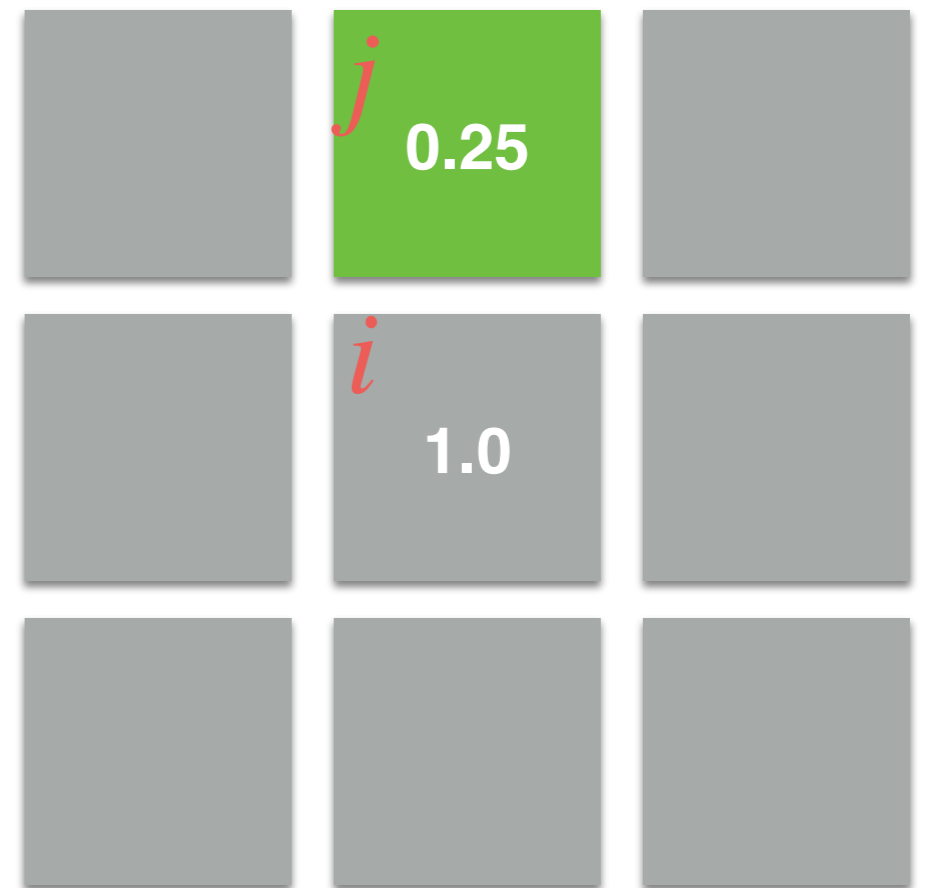
$$g(1.0 - 0.25) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 2 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



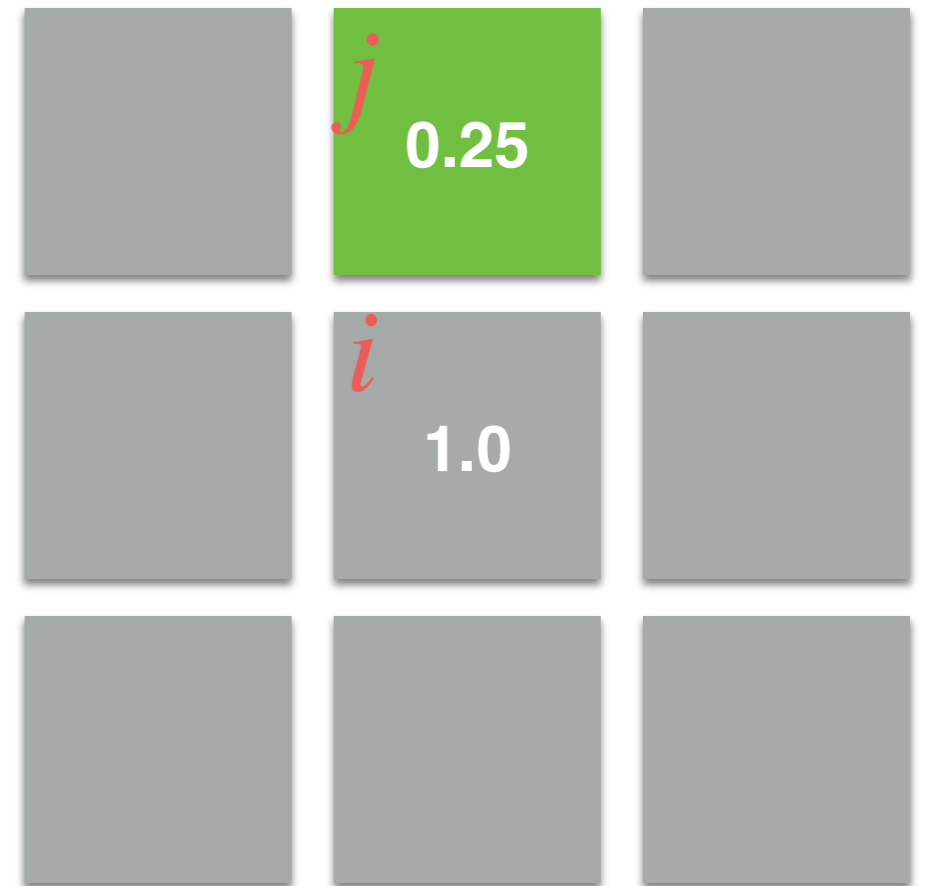
$$g(0.75) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 2 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



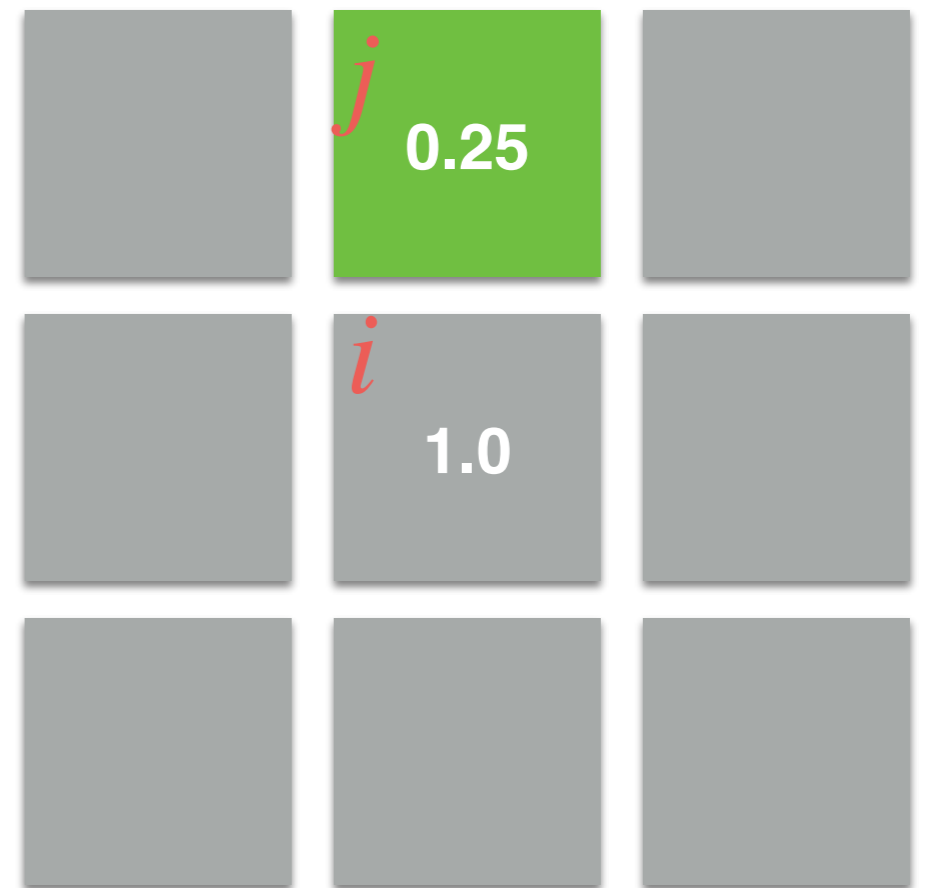
$$0.25 \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 2 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



$$0.25 \cdot 0.9 > 0.8$$

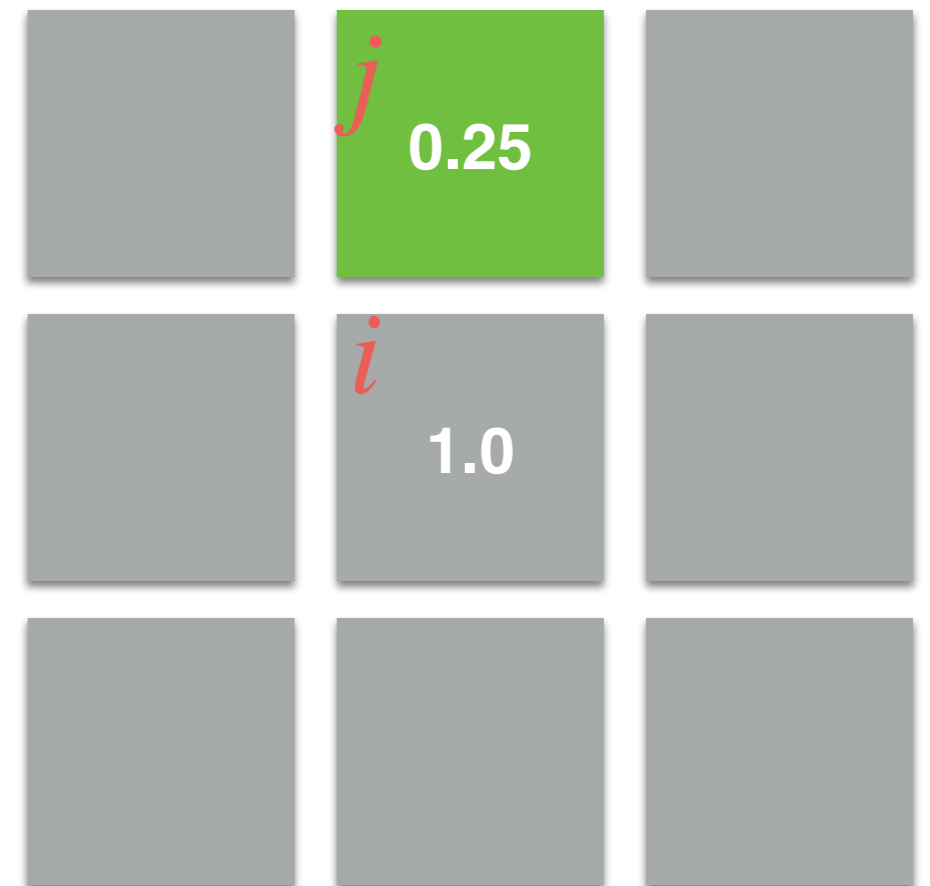
$$0.225 > 0.8$$

Stroke-Based: Example 2 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



0.225 is not greater than 0.8. So, l_i^{t+1} **for**
this neighbor remains the previous
value!

$$0.25 \cdot 0.9 > 0.8$$

$$0.225 > 0.8$$

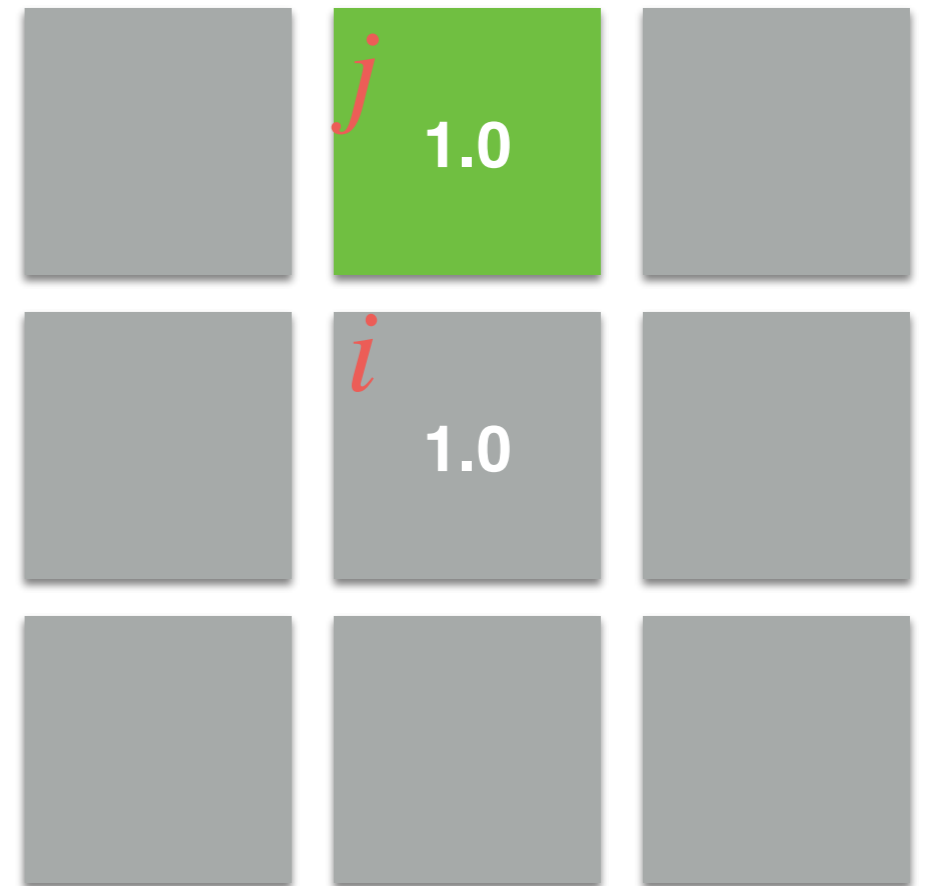
Example 3 Flat Area: Maintaining Labels

Stroke-Based: Example 3 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.7, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



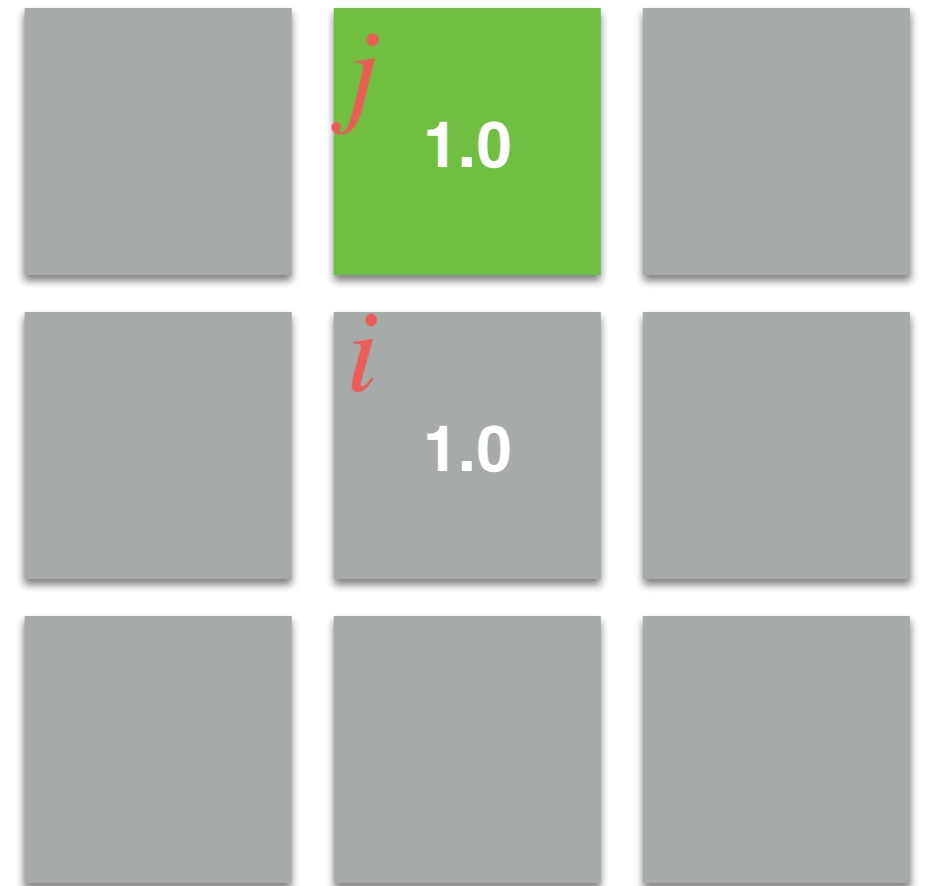
$$g(C_i^t - C_j^t) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 3 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.7, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



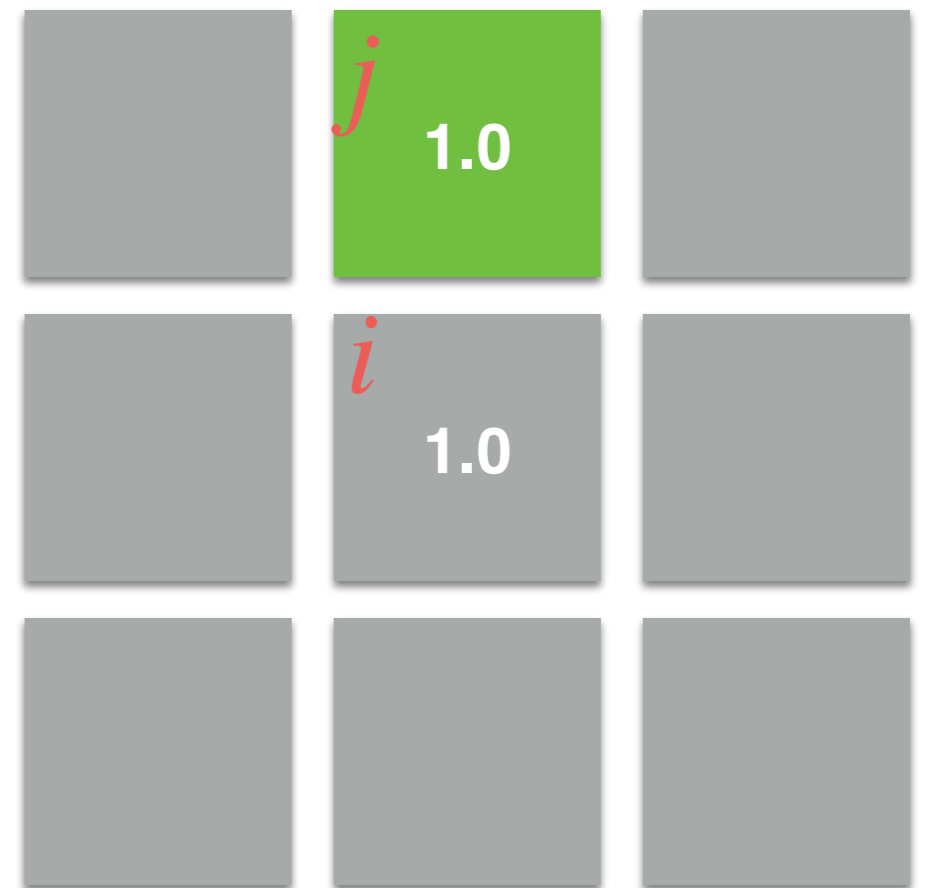
$$g(1.0 - 1.0) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 3 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.7, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



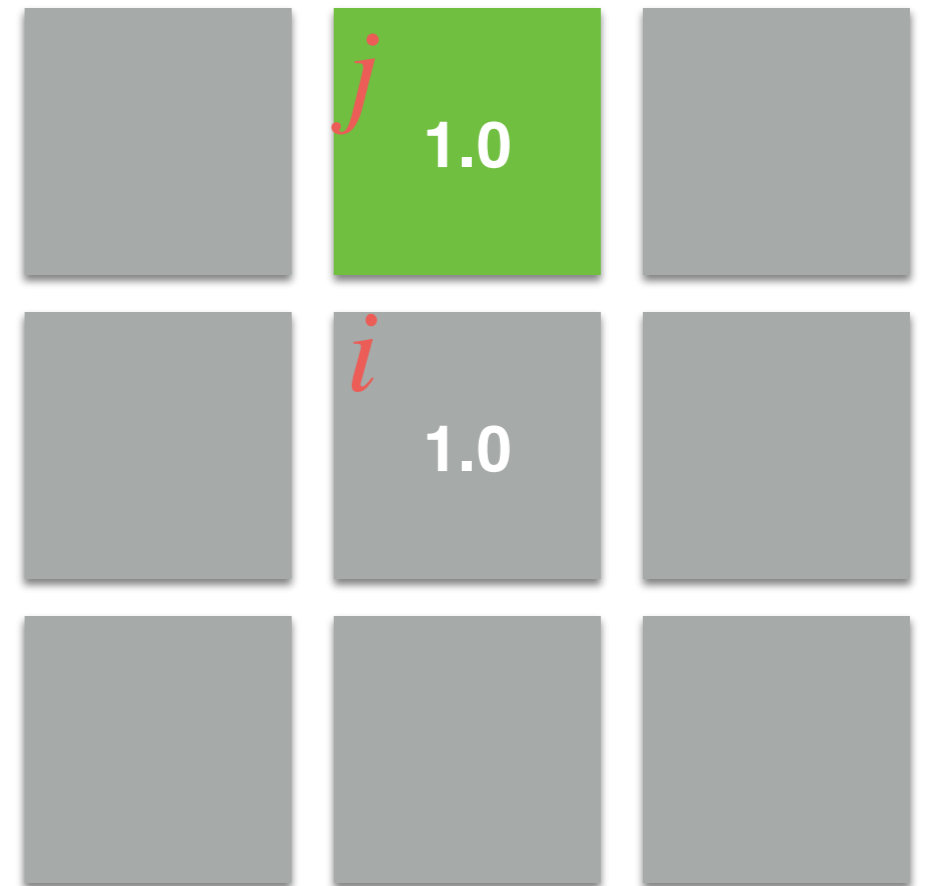
$$g(0.0) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 3 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.7, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



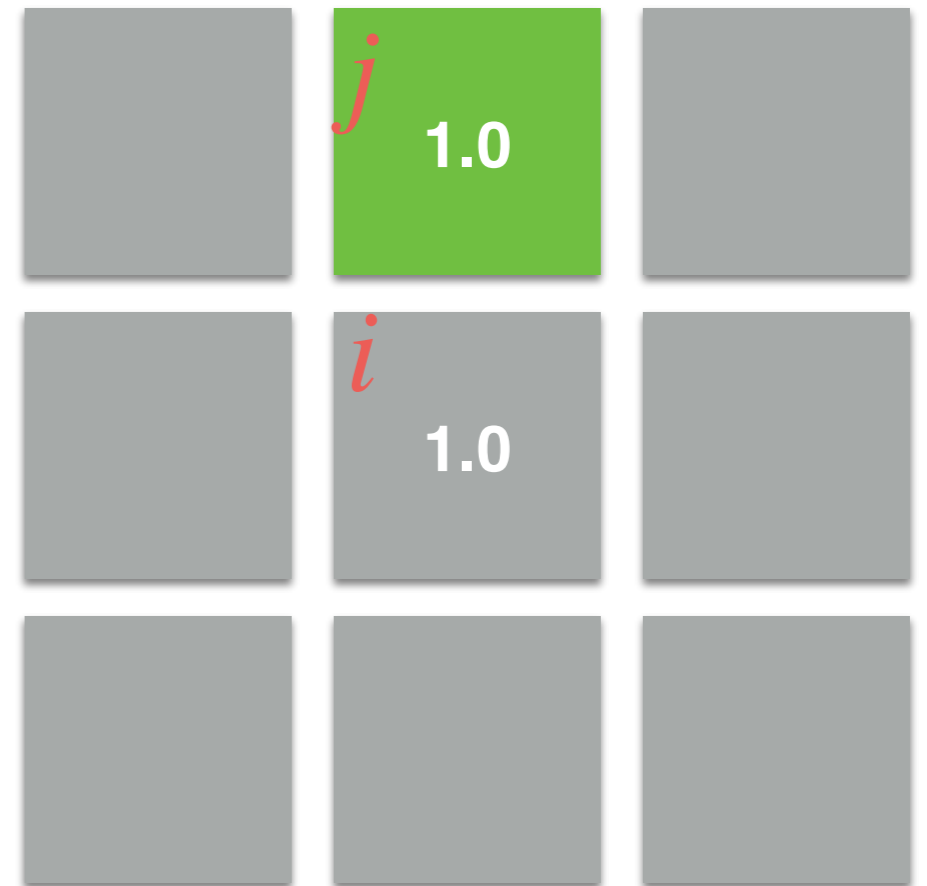
$$1.0 \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 3 - Flat Areas

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.7, 1.0 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.8, 1.0 \rangle$$



0.7 is not greater than 0.8. So, l_i^{t+1} **for this neighbor** remains the previous value!

$$1.0 \cdot 0.7 > 0.8$$

$$0.7 > 0.8$$

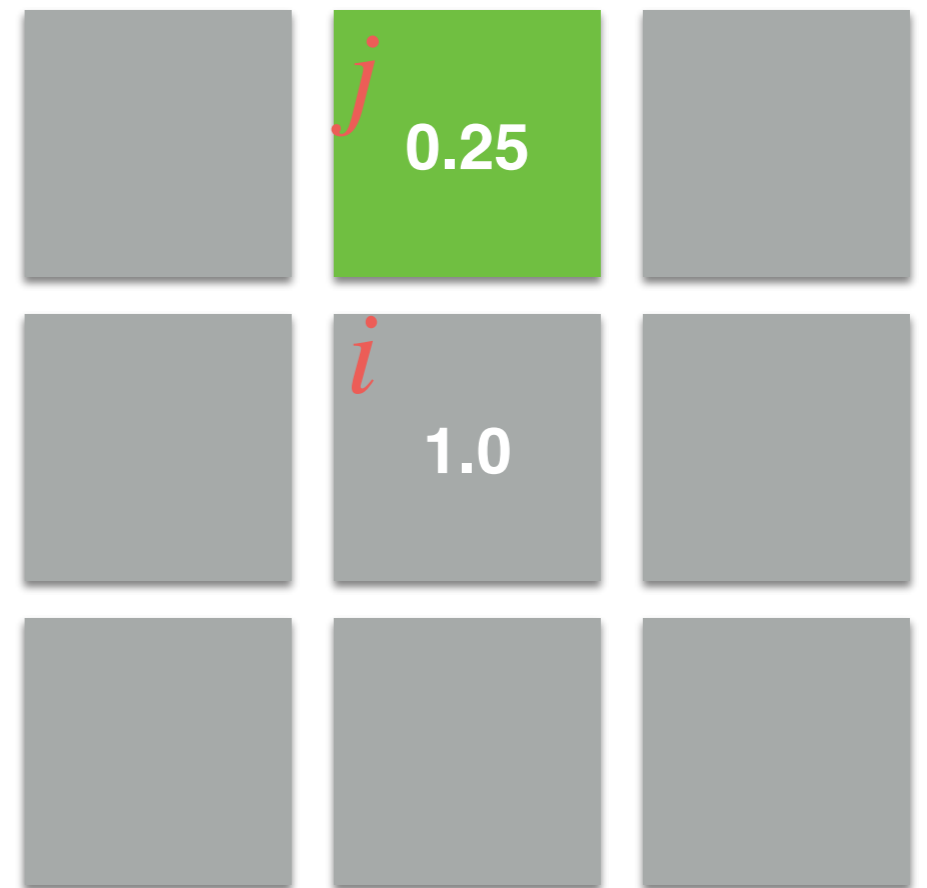
Example 4 Edges: Switching Labels

Stroke-Based: Example 4 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.1, 1.0 \rangle$$



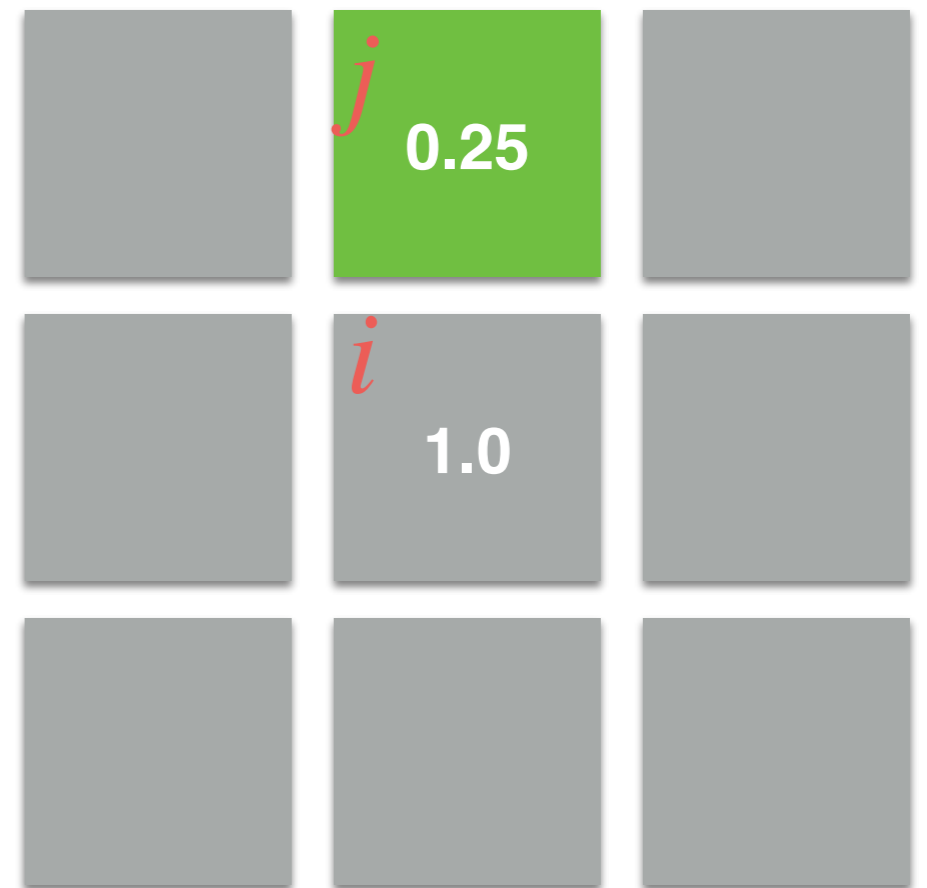
$$g(C_i^t - C_j^t) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 4 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.1, 1.0 \rangle$$



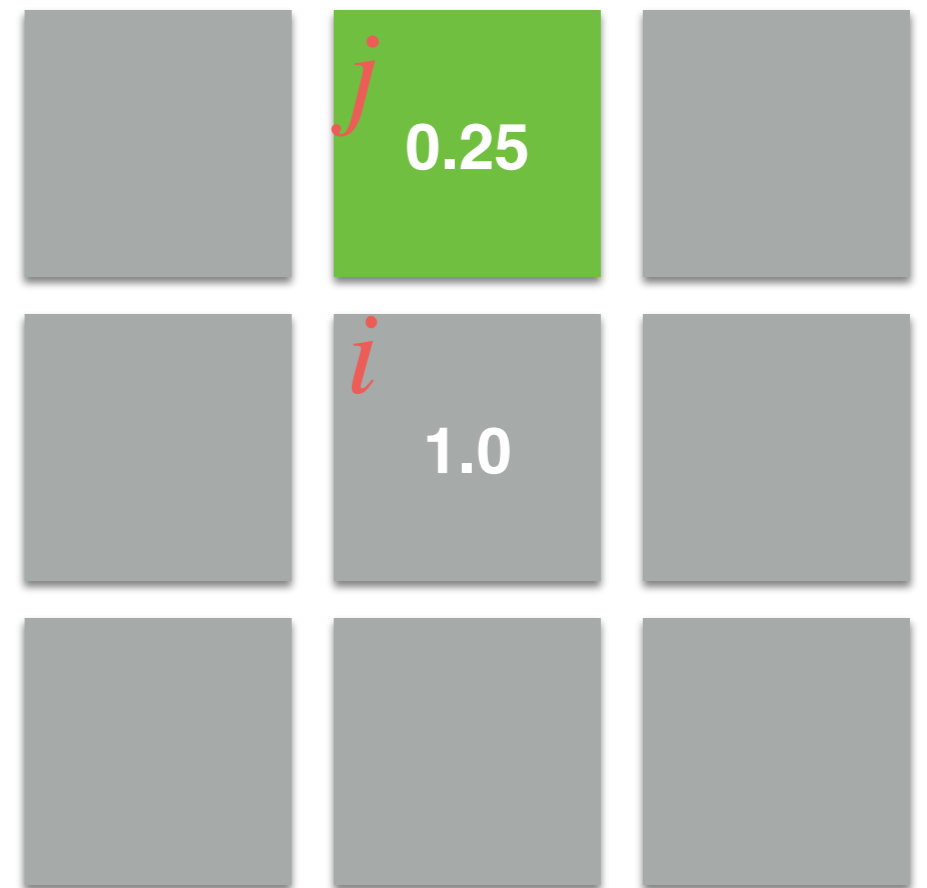
$$g(1.0 - 0.25) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 4 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.1, 1.0 \rangle$$



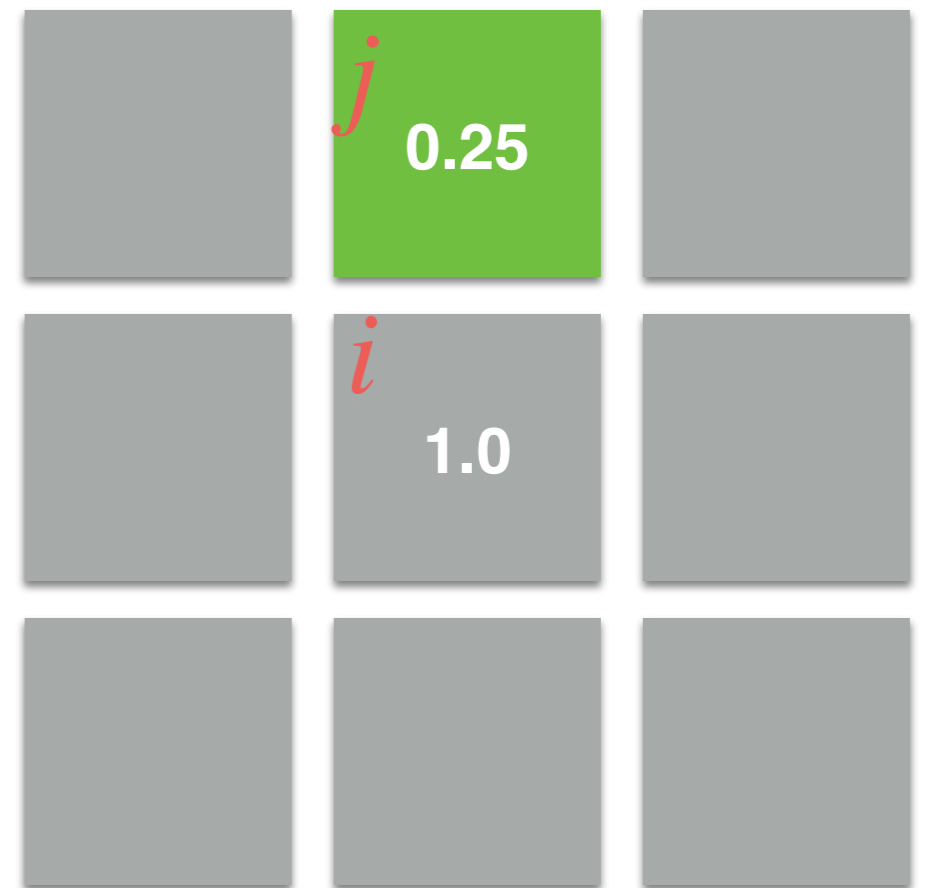
$$g(0.75) \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 4 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.1, 1.0 \rangle$$



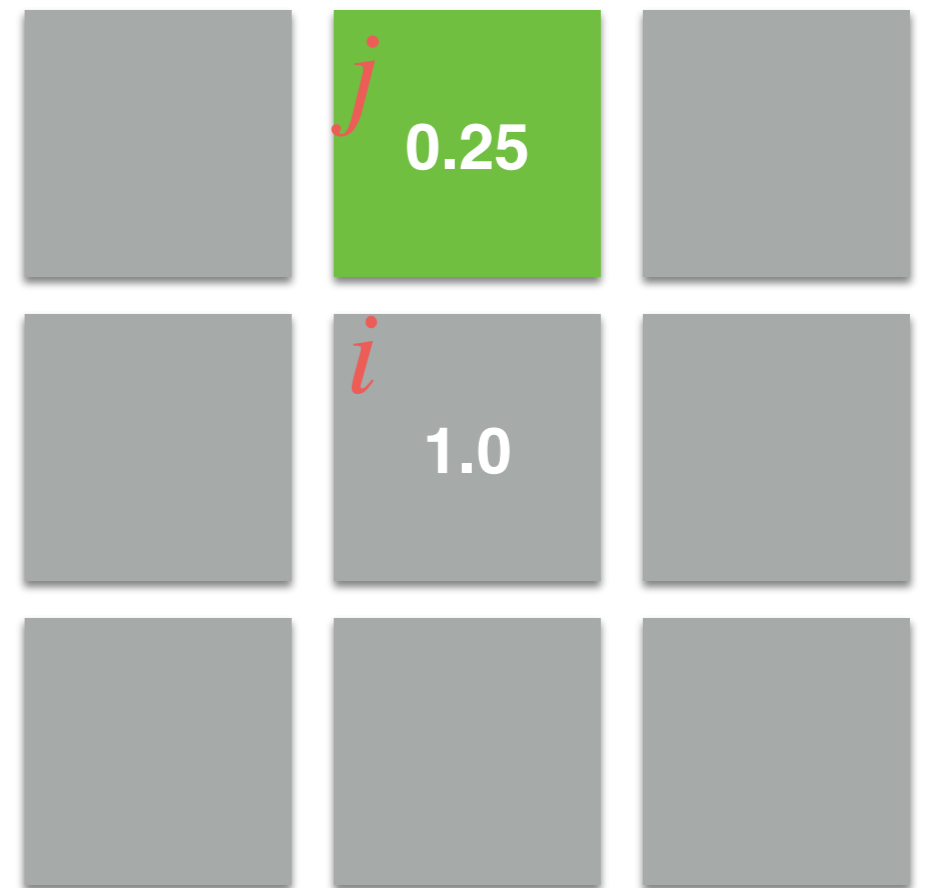
$$0.25 \cdot \theta_j^t > \theta_i^t$$

Stroke-Based: Example 4 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.1, 1.0 \rangle$$



$$0.25 \cdot 0.9 > 0.1$$

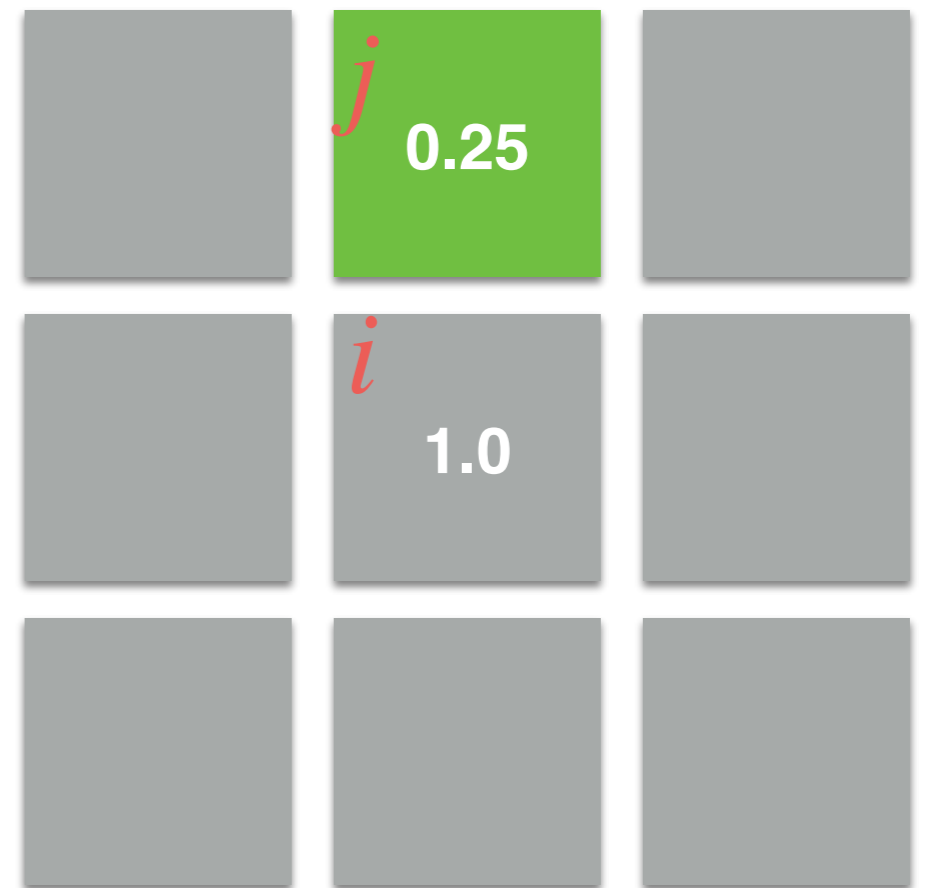
Stroke-Based: Example 4 - Edges

Current
neighbor

$$\langle l_j^t, \theta_j^t, C_j^t \rangle = \langle 0, 0.9, 0.25 \rangle$$

$$\langle l_i^t, \theta_i^t, C_i^t \rangle = \langle 1, 0.1, 1.0 \rangle$$

0.225 is greater than 0.1. So, so we
assign the value l_j^t to l_i^{t+1} .



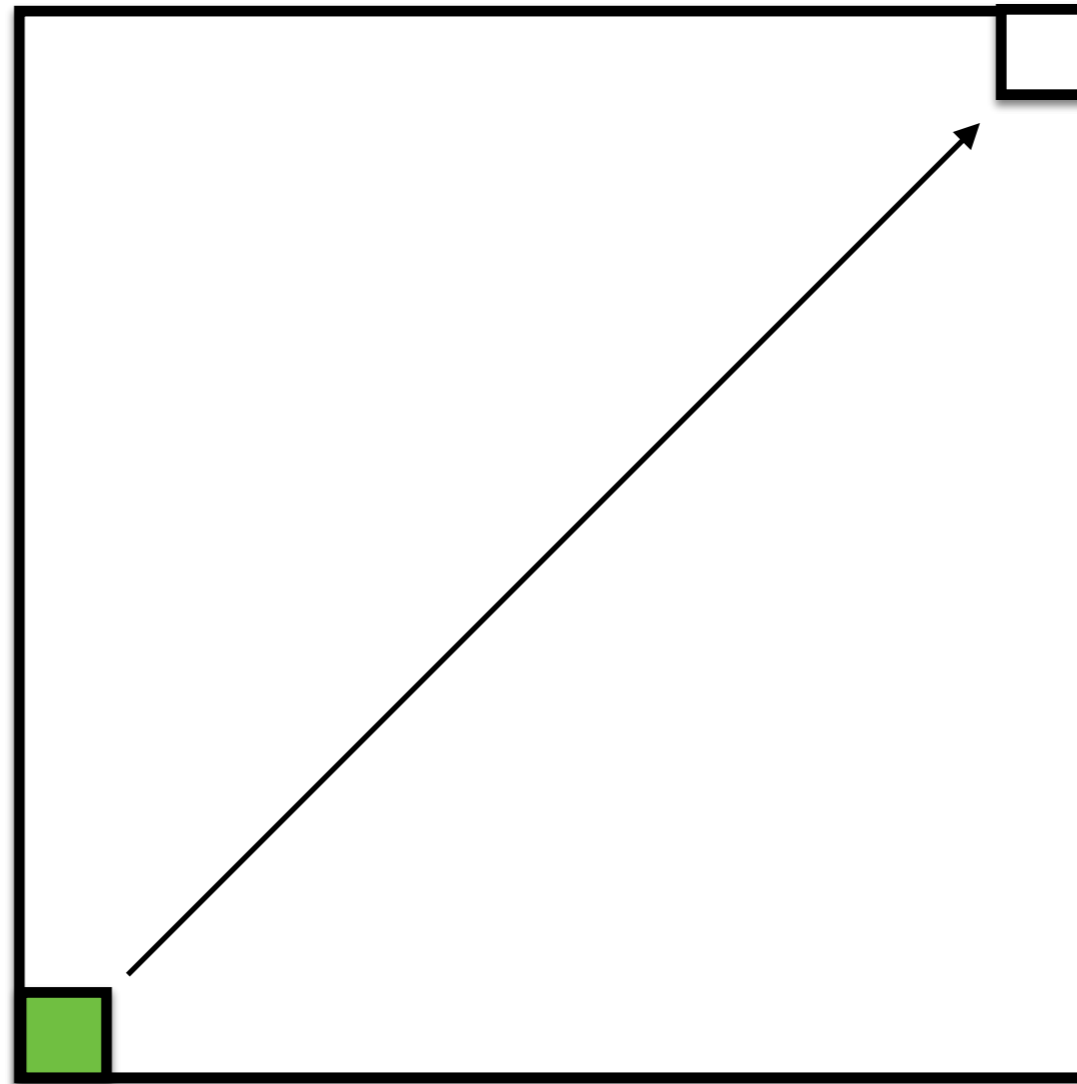
$$0.225 > 0.1$$

$$l_i^{t+1} = l_j^t = 0!$$

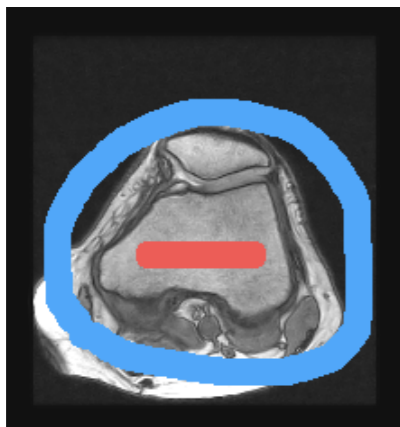
Stroke-Based: Grow-Cut

- Stopping criteria:
 - This process is iterated until either convergence; i.e., *no changes in the labels!*
 - Labels have been propagated for enough iterations; e.g., the *number of pixels of the diagonal*. This trick is helpful for reducing the total computational time.

Stroke-Based: Grow-Cut

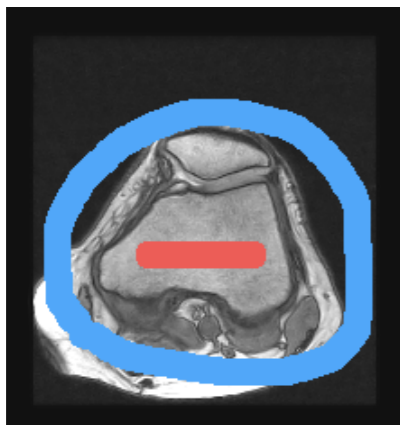


Stroke-Based: Grow-Cut Example



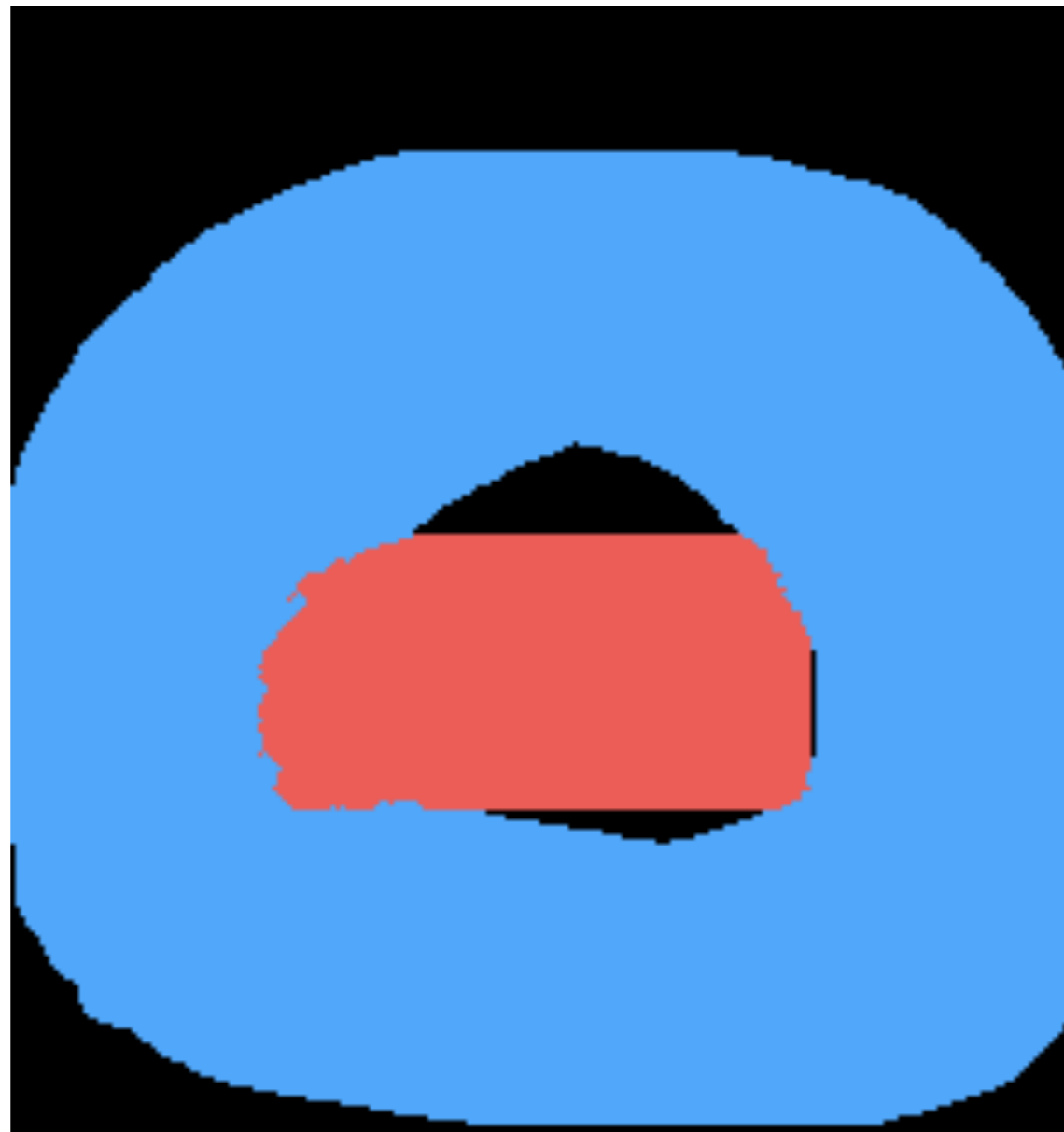
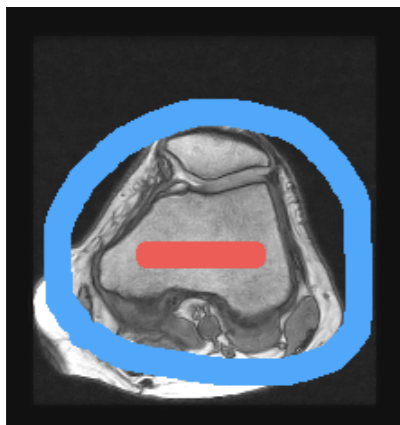
Iteration = 0

Stroke-Based: Grow-Cut Example



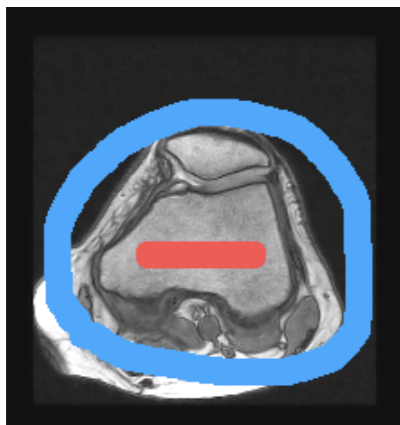
Iteration = 10

Stroke-Based: Grow-Cut Example



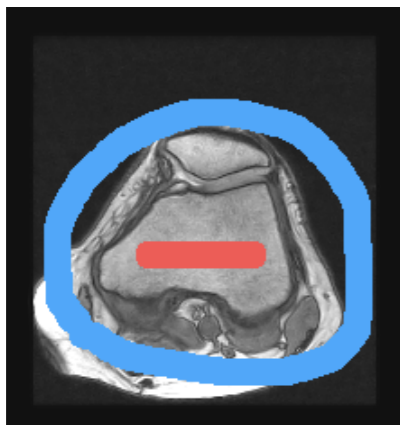
Iteration = 20

Stroke-Based: Grow-Cut Example



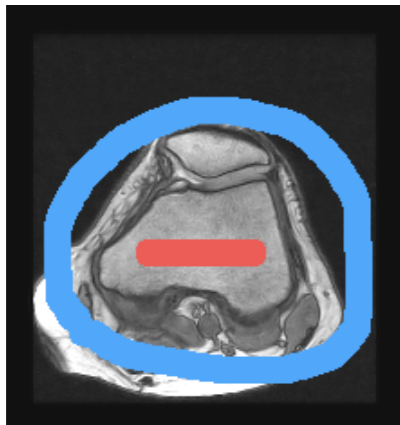
Iteration = 30

Stroke-Based: Grow-Cut Example



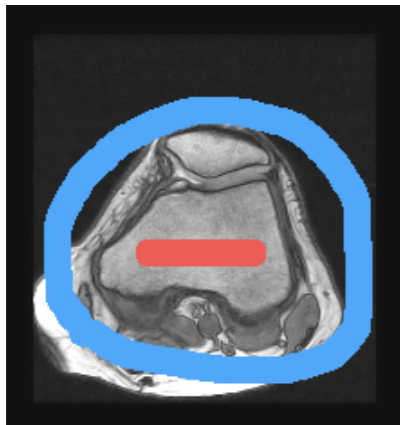
Iteration = 40

Stroke-Based: Grow-Cut Example



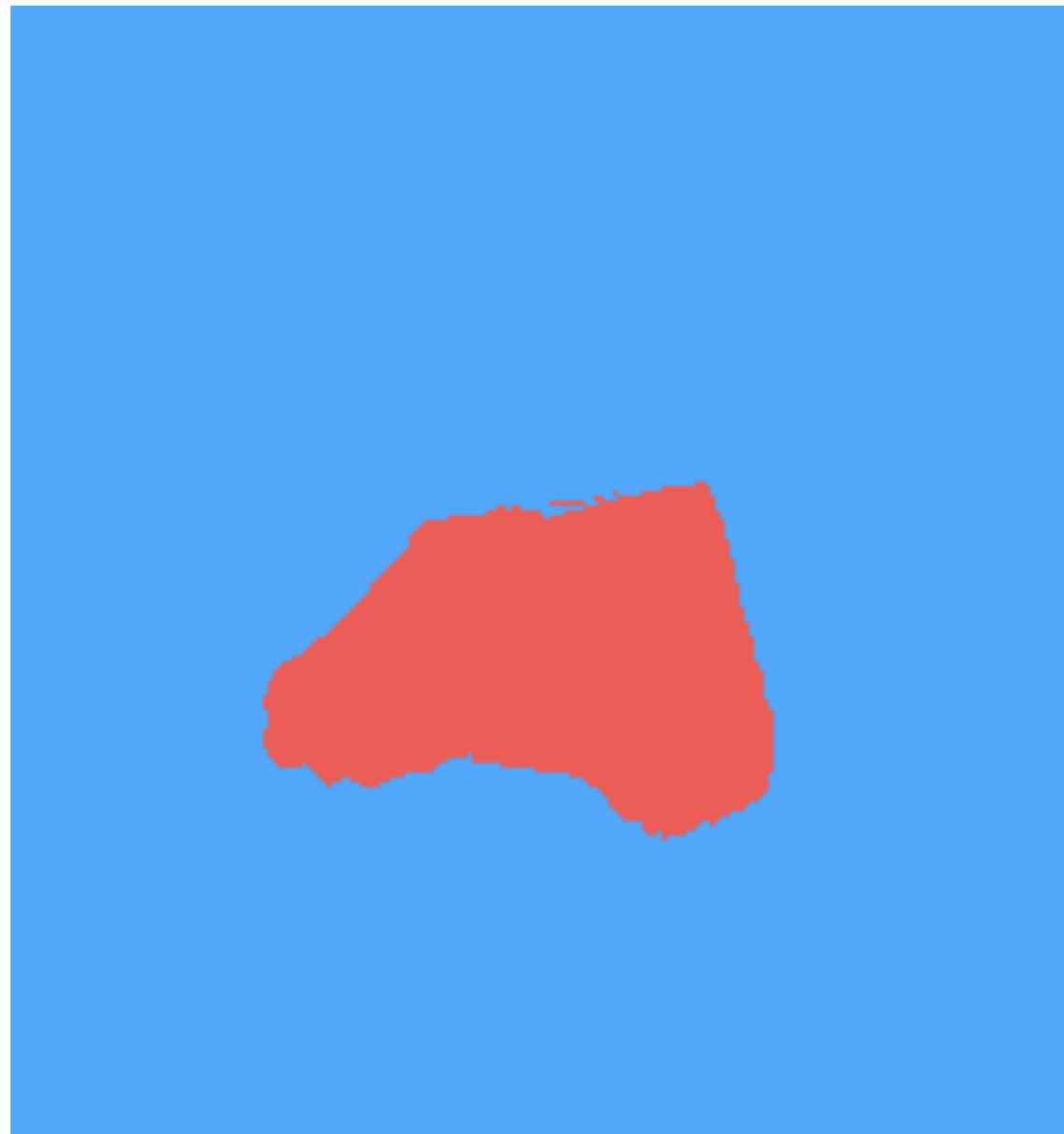
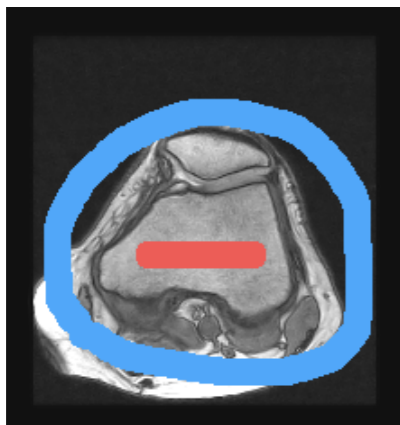
Iteration = 50

Stroke-Based: Grow-Cut Example



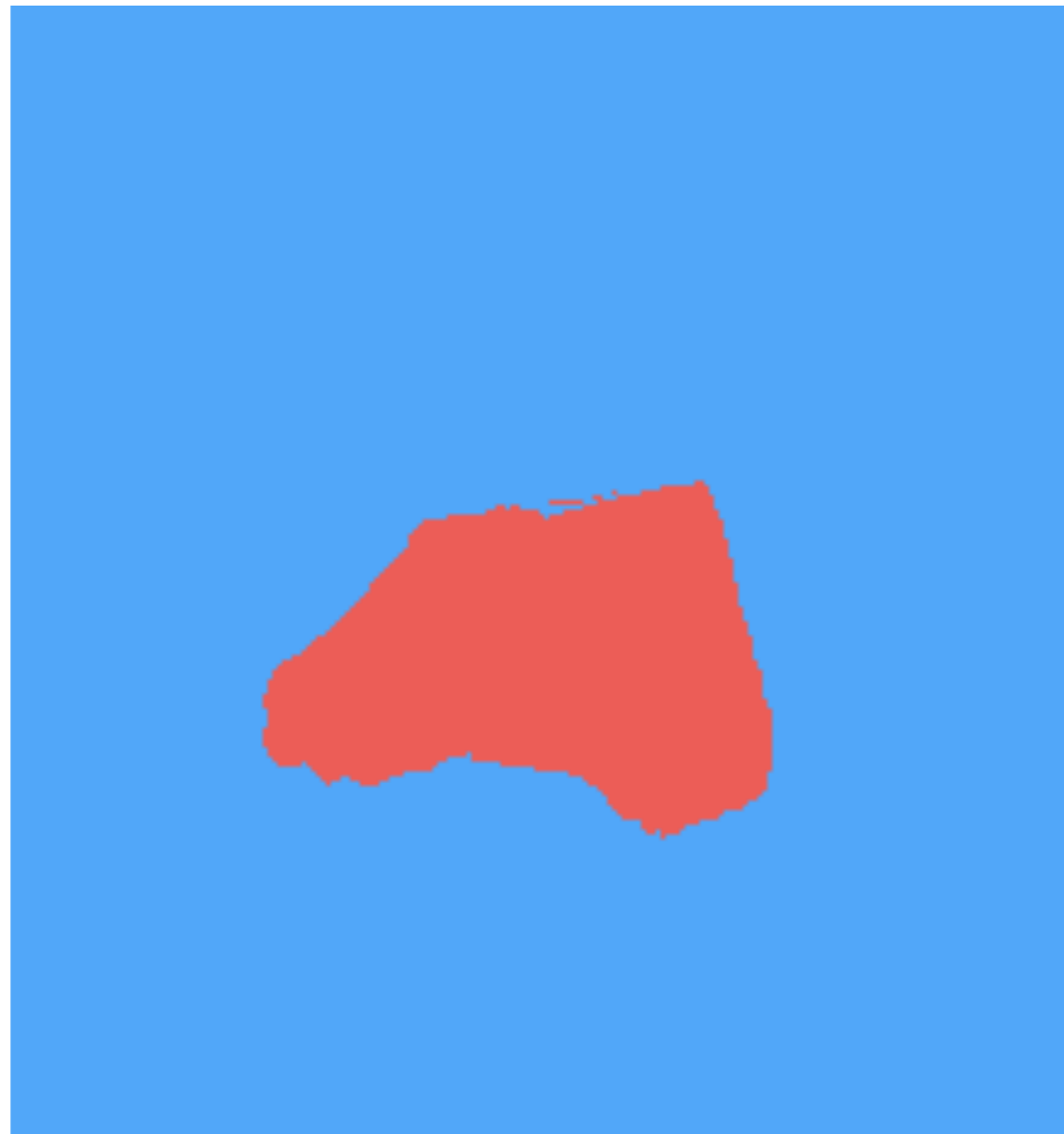
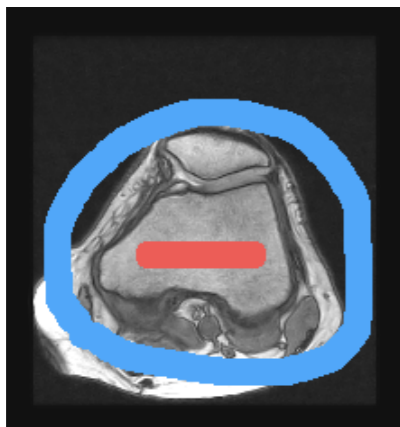
Iteration = 100

Stroke-Based: Grow-Cut Example



Iteration = 200

Stroke-Based: Grow-Cut Example



Iteration = 318

Stroke-Based: Grow-Cut

- This algorithm can be extended to 3D in a straightforward way, and it can be parallelized on the GPU.
- Disadvantages:
 - It is computationally slow!

that's all folks!