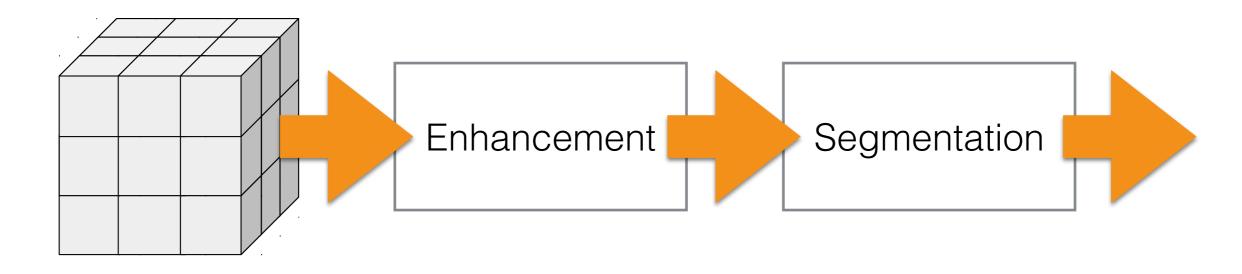
3D from Volume: Part II

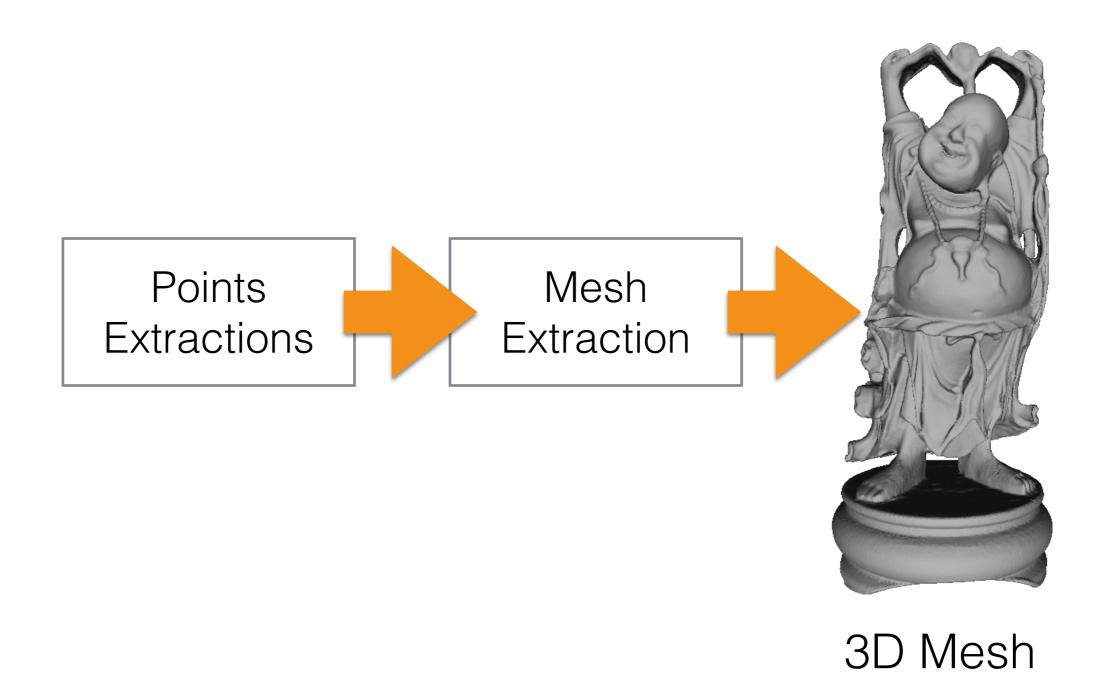
Francesco Banterle, Ph.D. francesco.banterle@isti.cnr.it

The Processing Pipeline

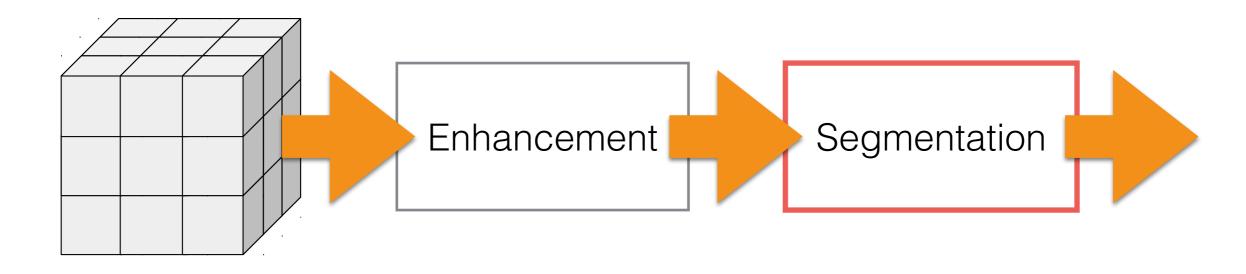


RAW Volume

The Processing Pipeline



The Processing Pipeline

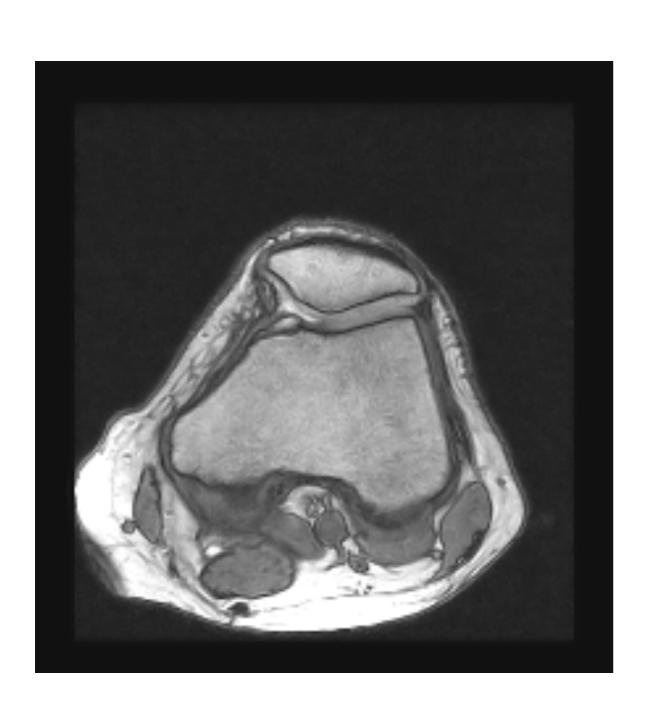


RAW Volume

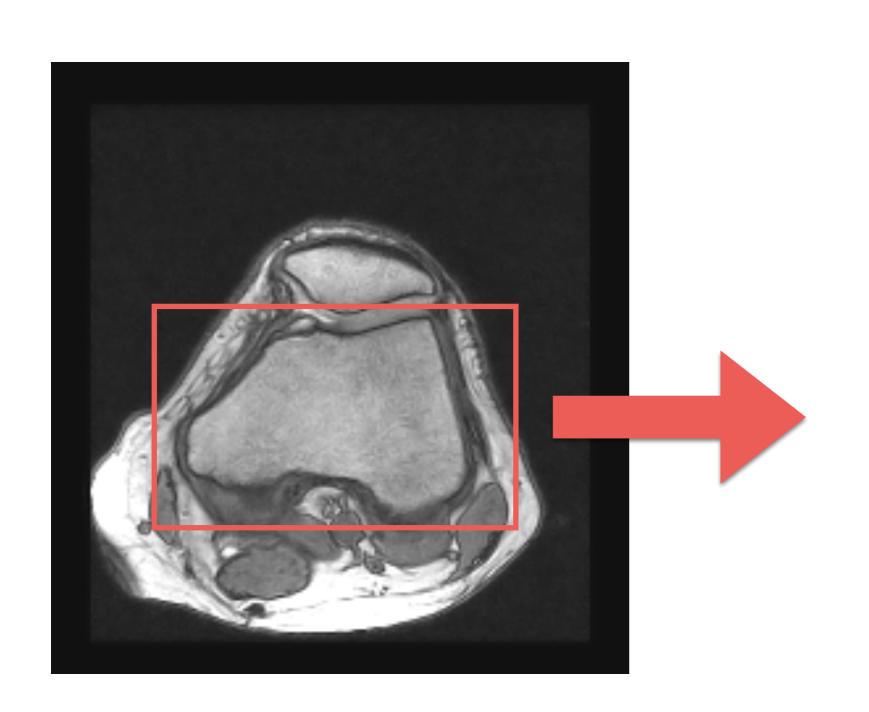
2D/3D Segmentation

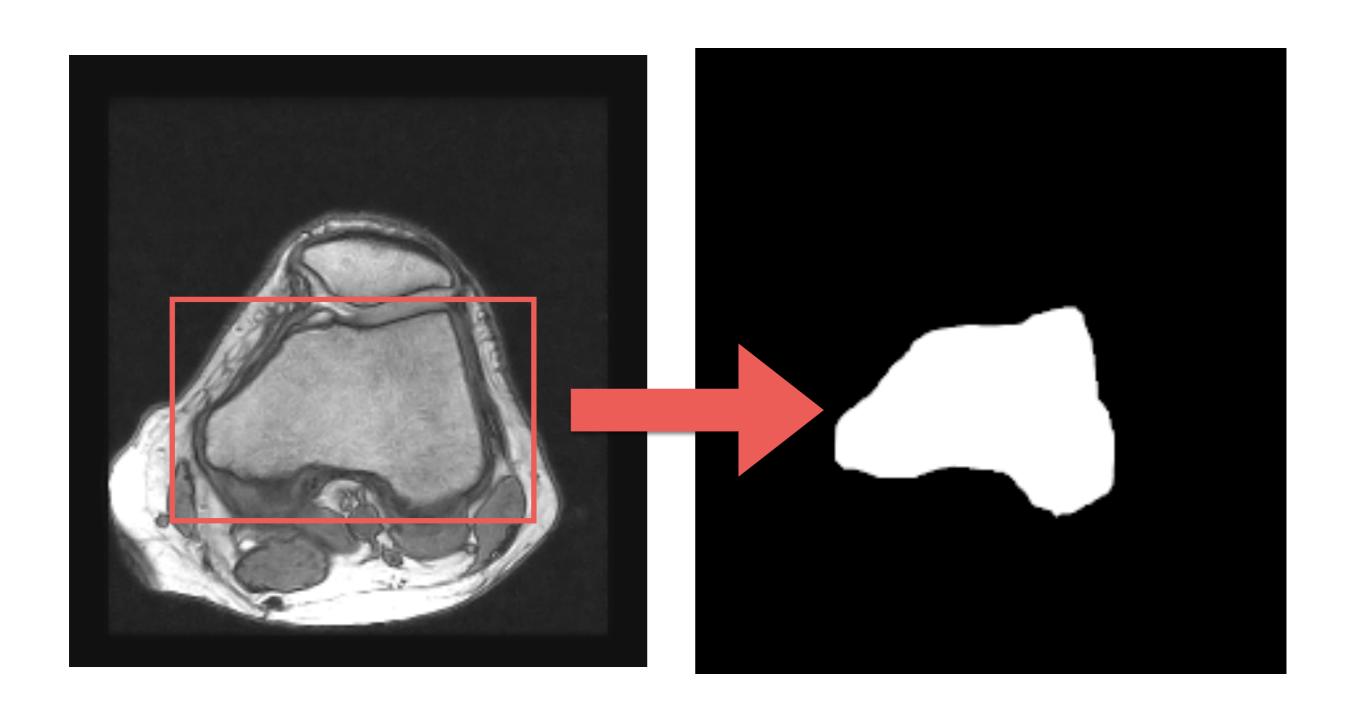
Segmentation

- 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 —> the pixel/voxel belongs to a structure of our interest
- 0 —> the pixel/voxel does not!









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 as label a number in [1,k]. Background is always 0!

3D Segmentation

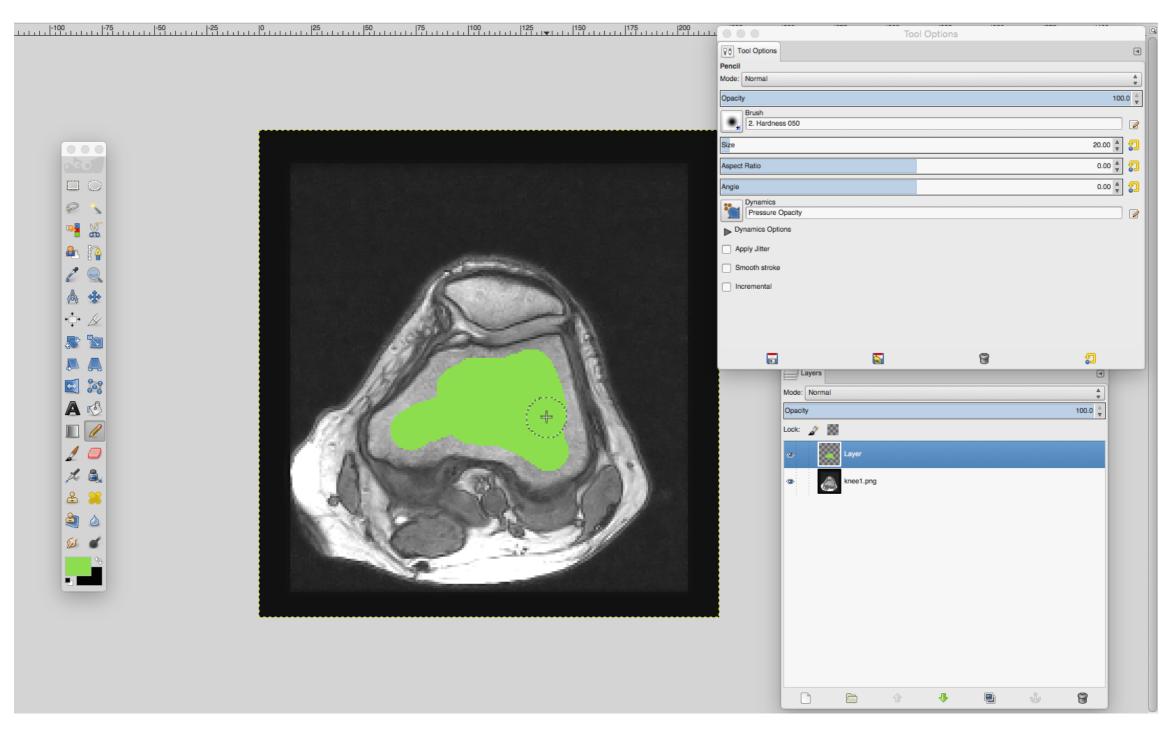
- There are typically two approaches:
 - 2D segmentation for each slice
 - 2D segmentation of a slice and propagation of the segmentation

Manual Segmentation

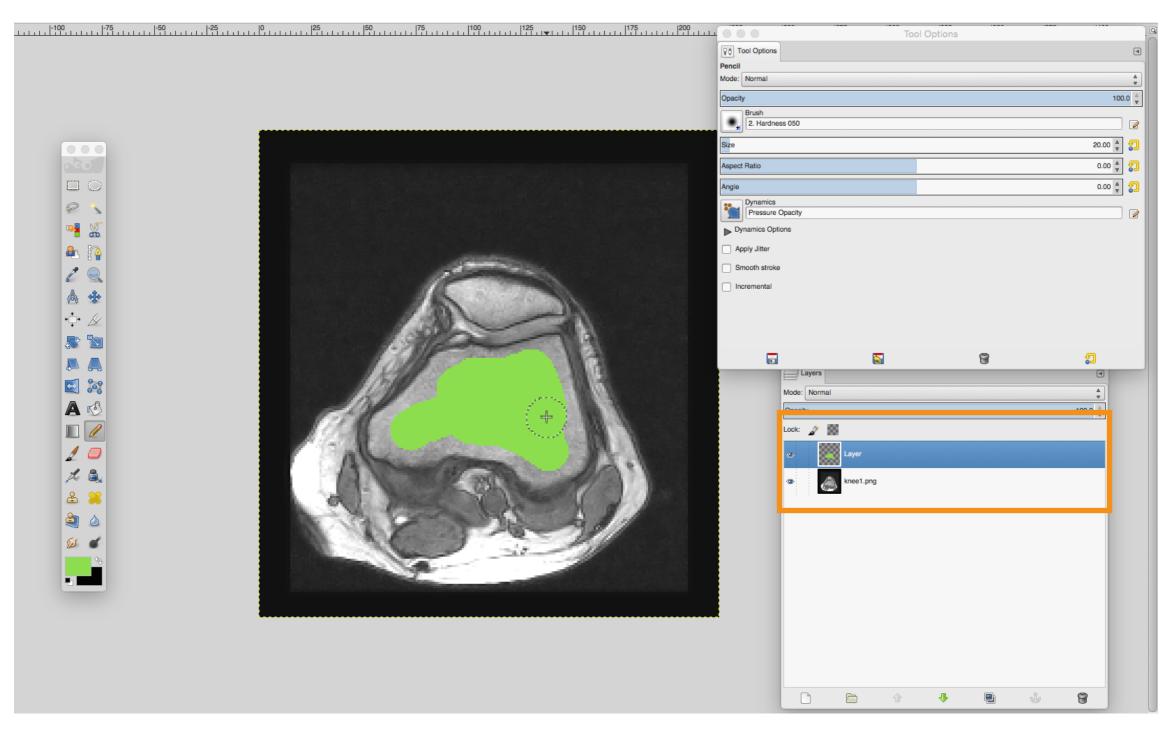
Manual Segmentation: Painting Approach

- 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



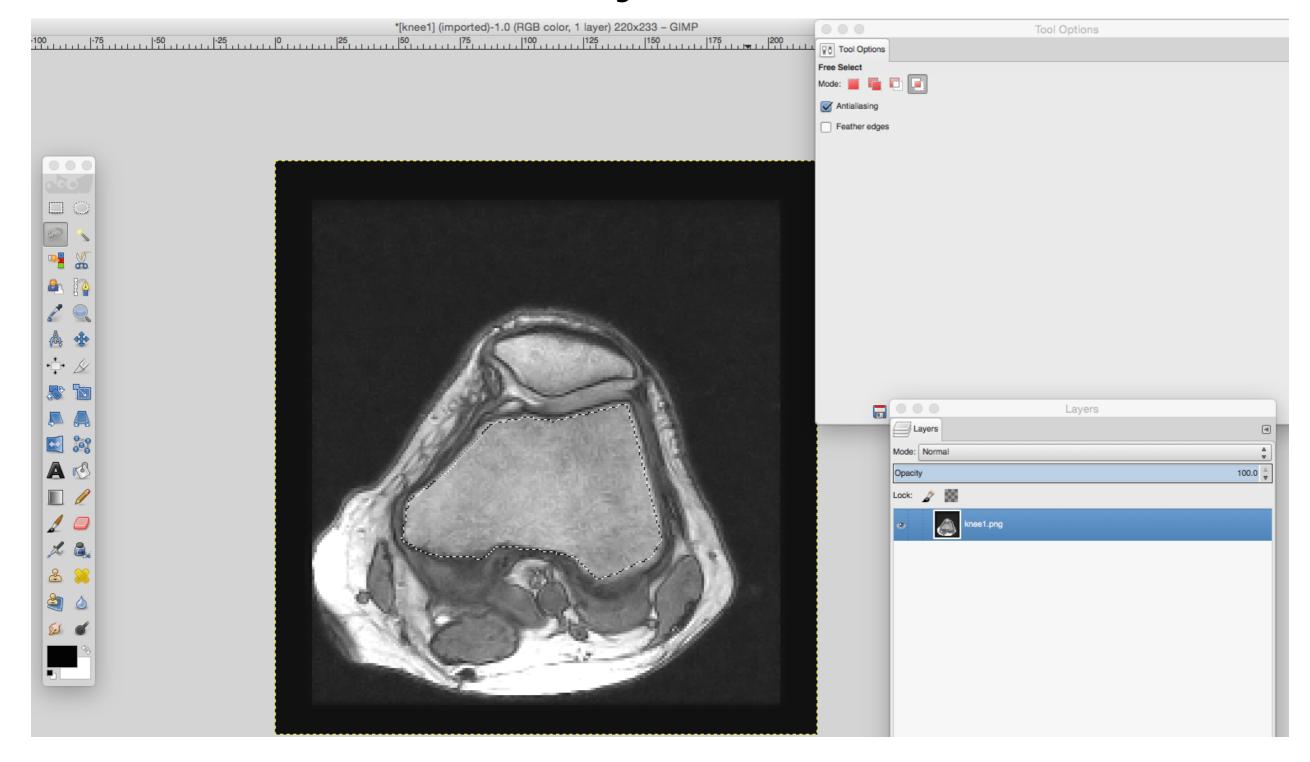
Manual Segmentation: Painting Approach



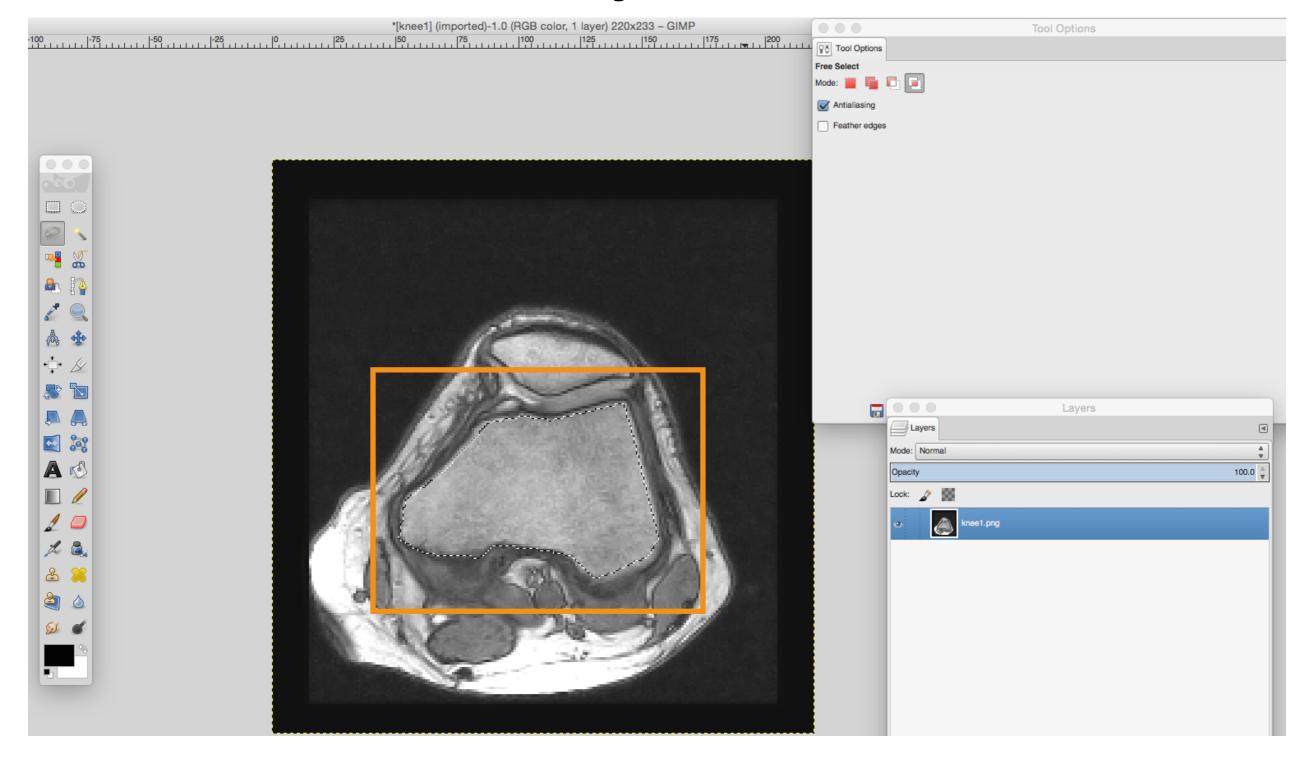
Manual Segmentation: Boundary Definition

- 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.

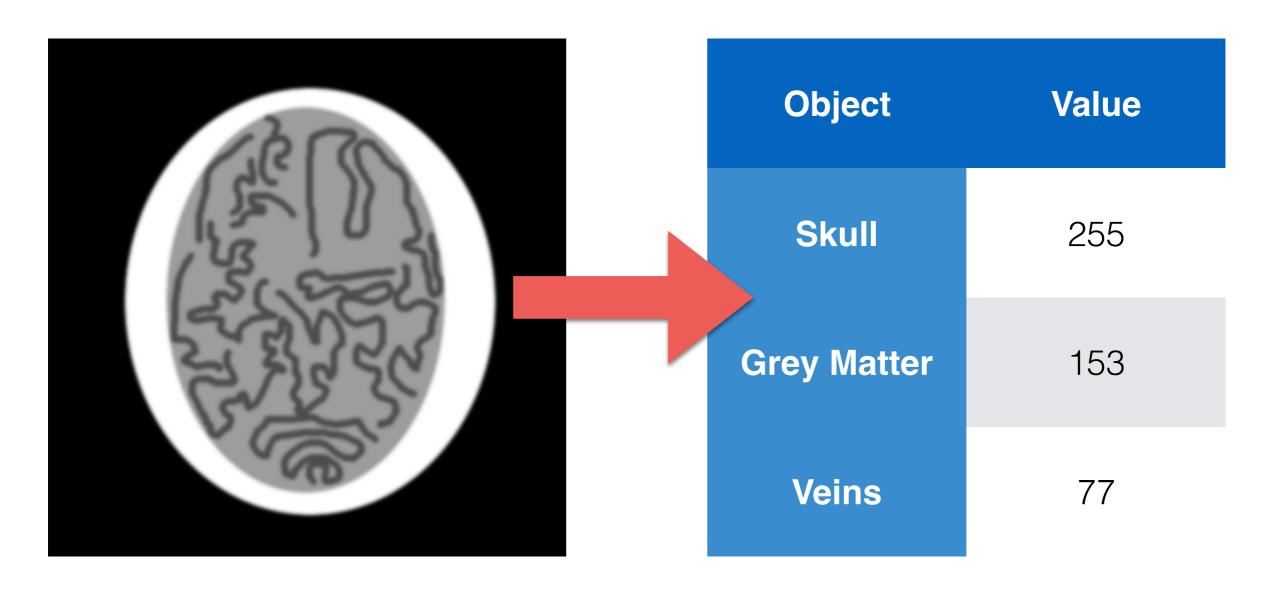
Manual Segmentation: Boundary Definition



Manual Segmentation: Boundary Definition



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



• 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:

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

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

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:

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

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

This means:

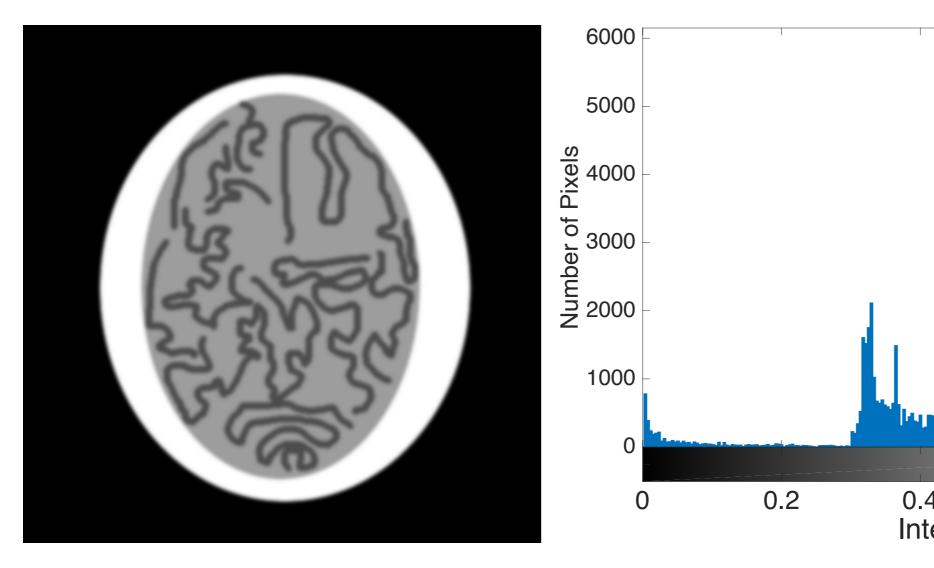
Reference Value

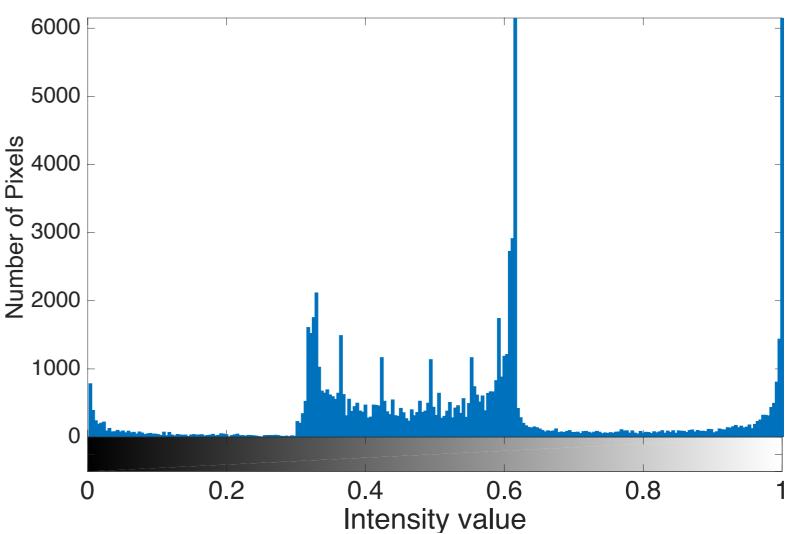
$$M(i,j) = \begin{cases} 1 & \text{if } d(I(i,j), I_t) < t \\ 0 & \text{otherwise} \end{cases}$$
 Threshold

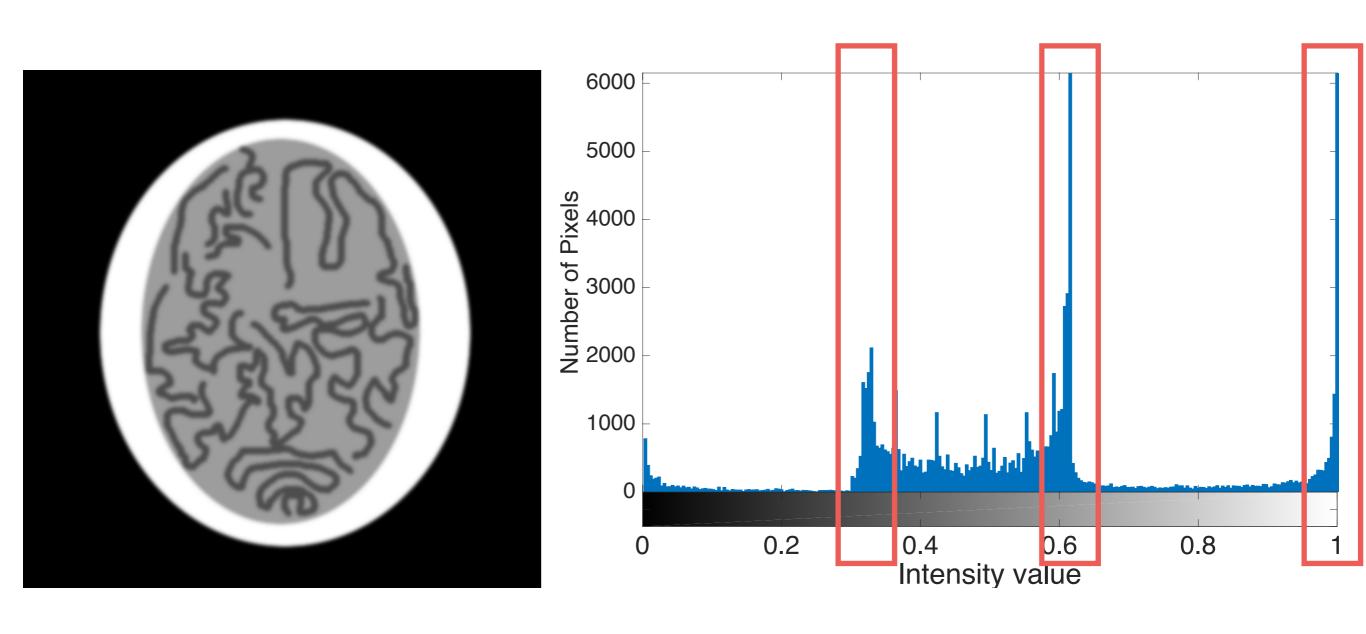
We can have different distance functions:

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

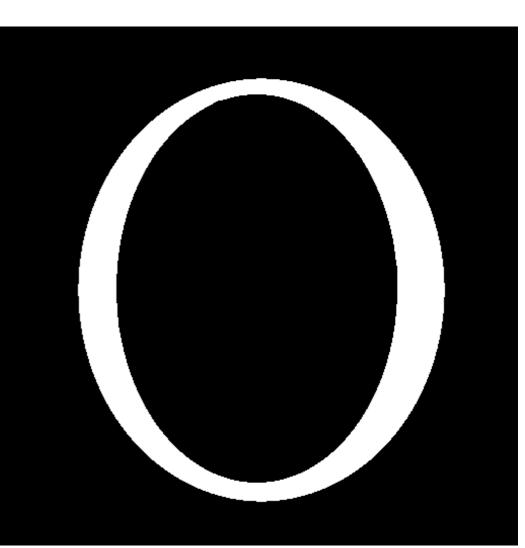
$$d(x, y, \sigma) = \exp\left(-\frac{(x - y)^2}{2\sigma^2}\right)^2$$





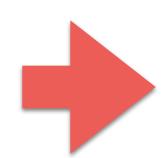






$$I_{t} = 1$$
 $t = 0.1$

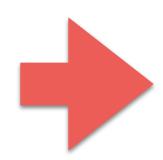






$$I_t = 0.6$$
 $t = 0.1$







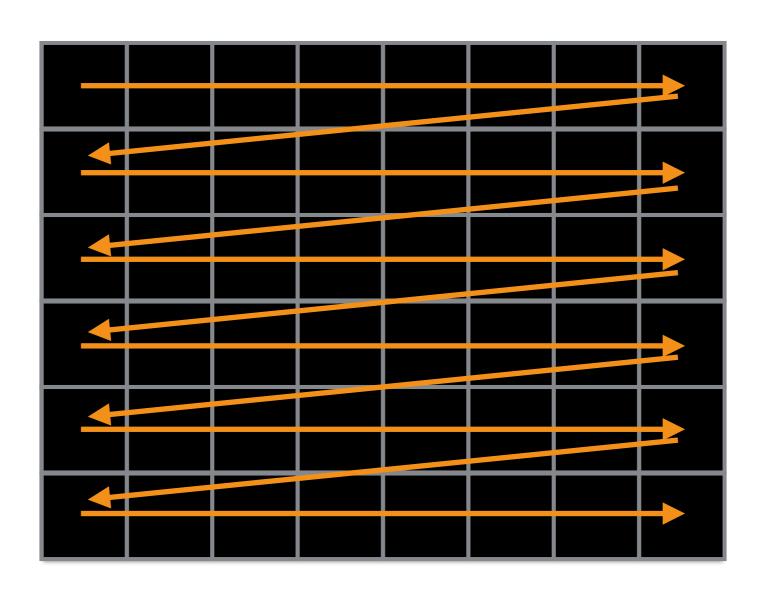
$$I_t = 0.6$$
 $t = 0.1$

$$t = 0.1$$

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

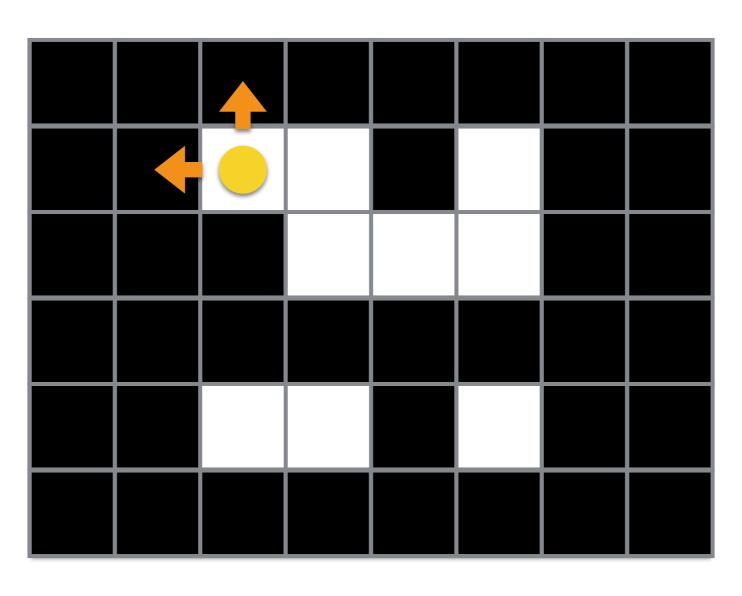


- A two-pass algorithm that works in scan order (from left to right and from top to bottom).
- 1-Pass: it creates labels to groups of pixel.
- 2-Pass: it merges groups that are connected.

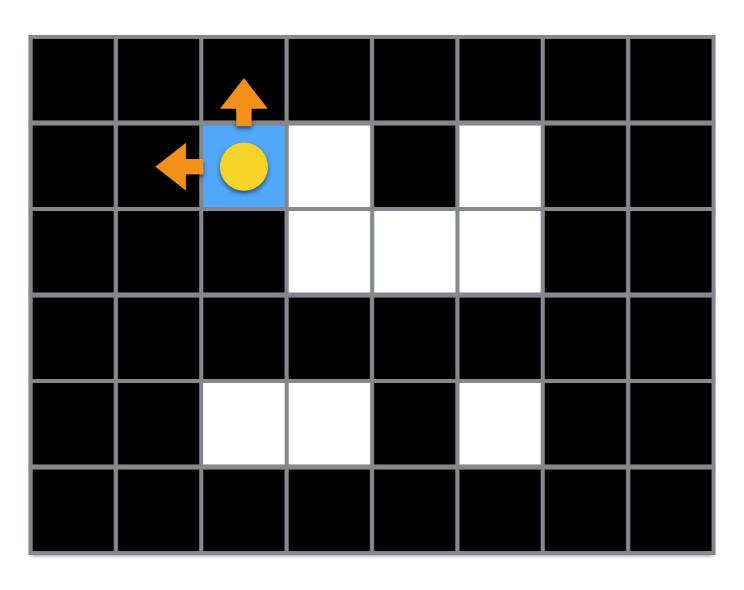


Scan order

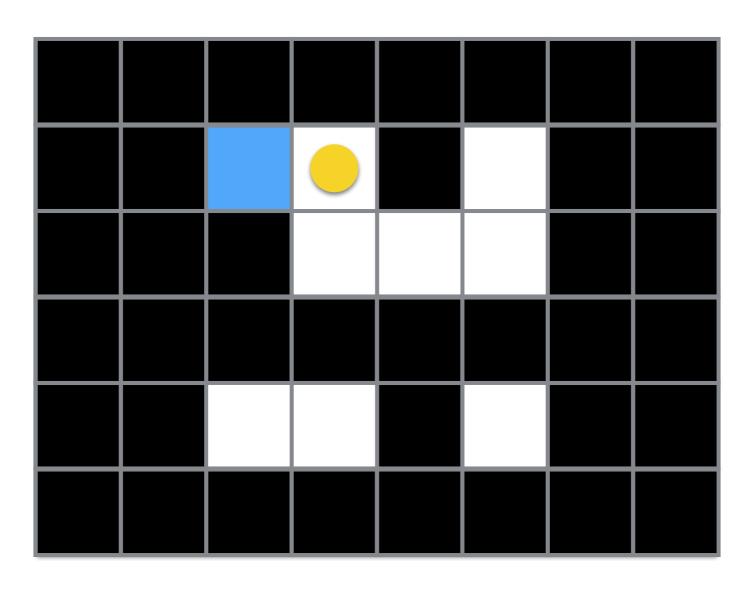
First Pass



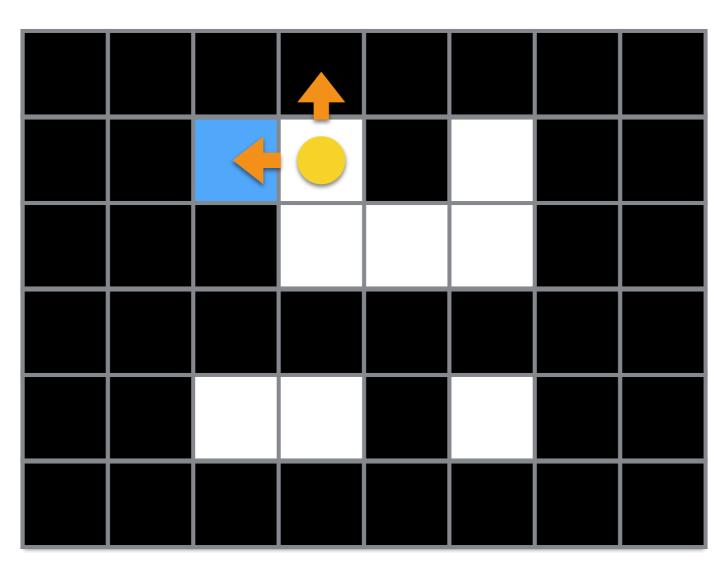
We check up and left neighbors to see if they have a label.



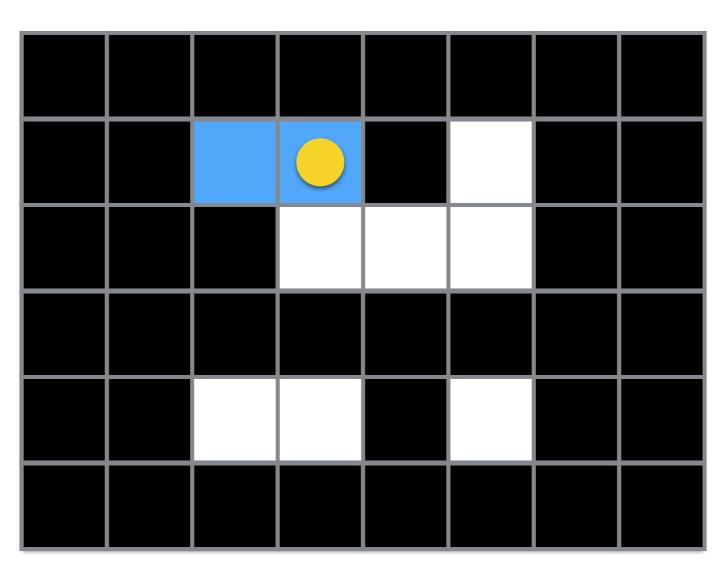
If not we create a new one.



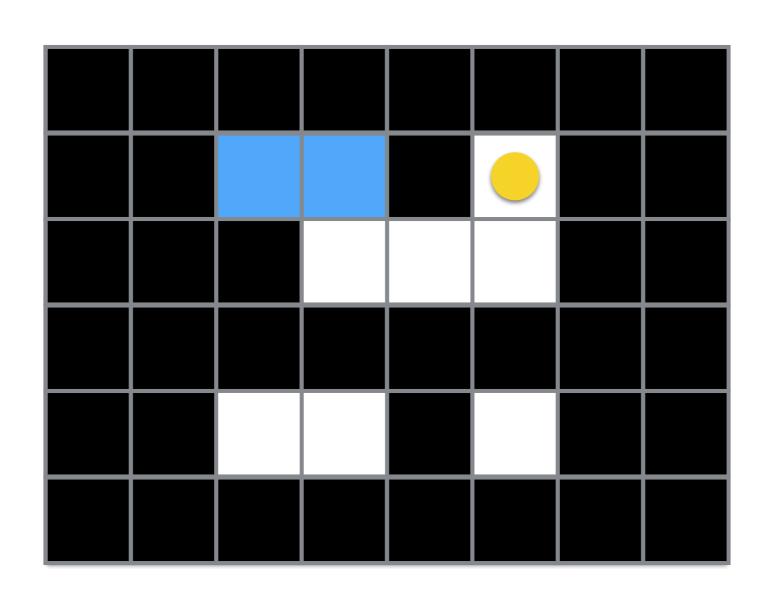
Then, we move right, and we repeat the process.

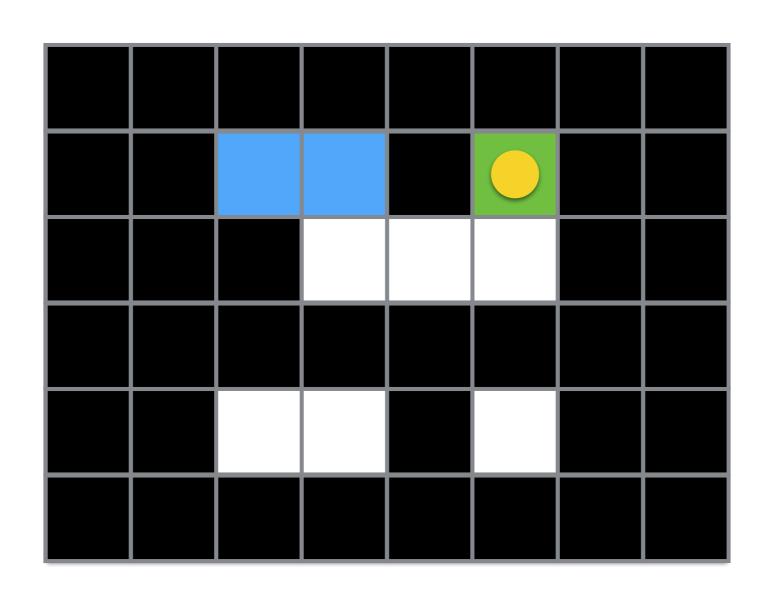


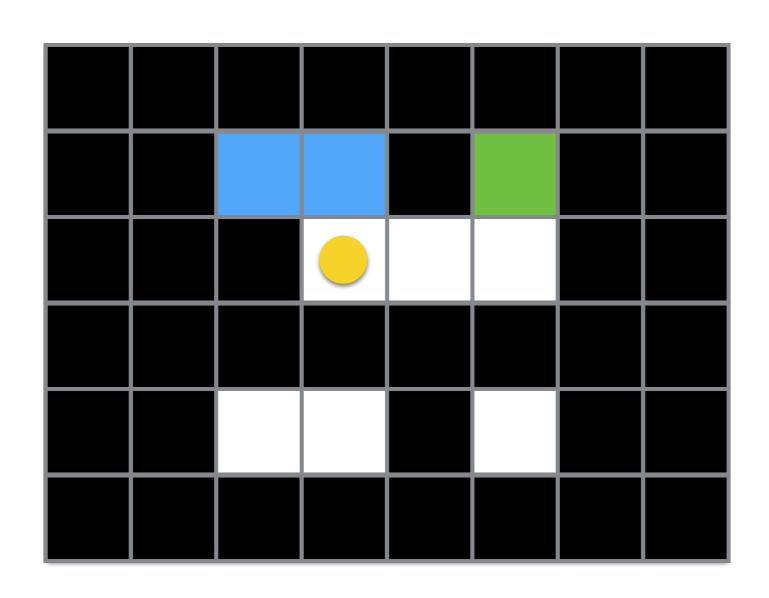
In this case, the left neighbor has a label, so we reuse it.

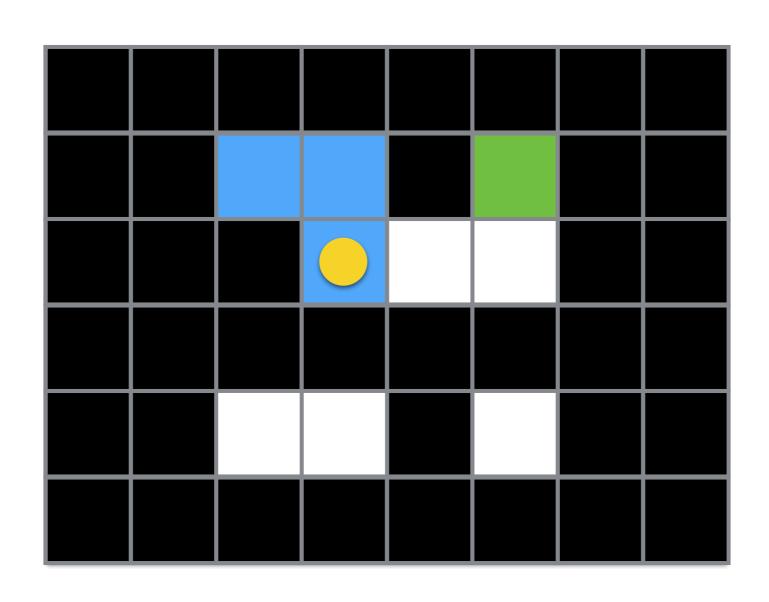


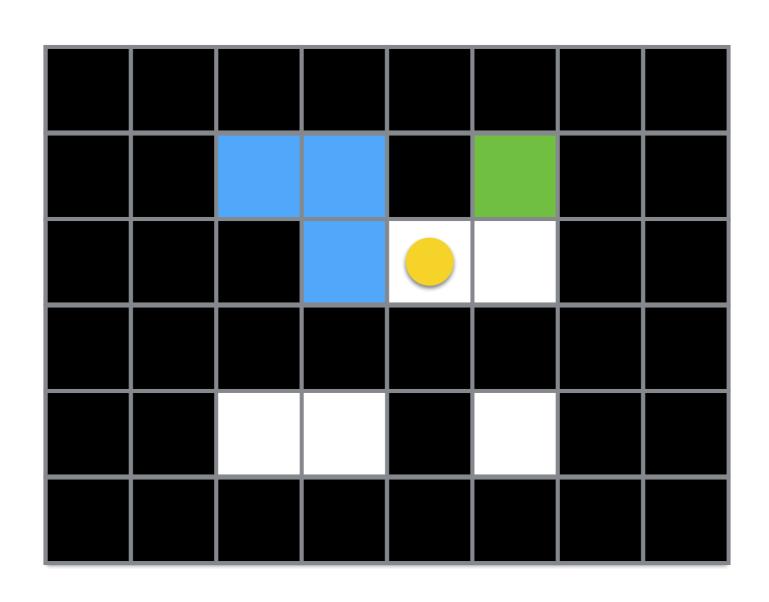
In this case, the left neighbor has a label, so we reuse it.

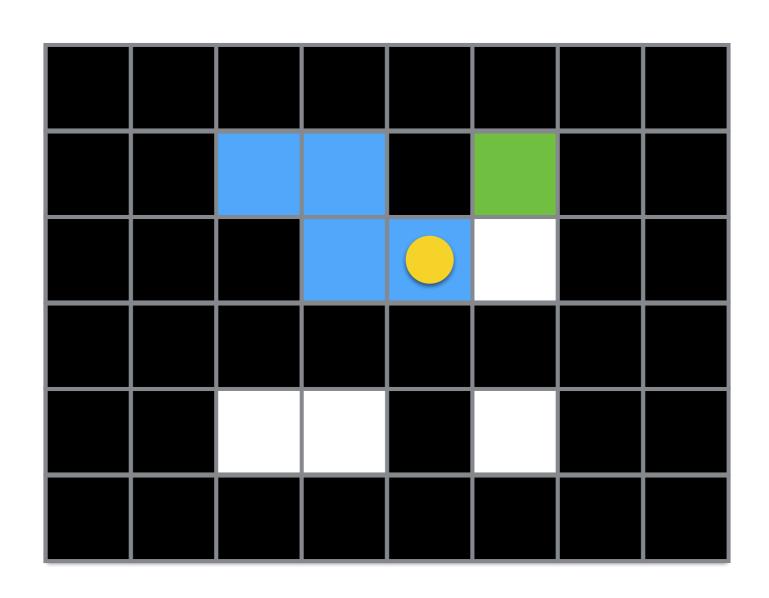






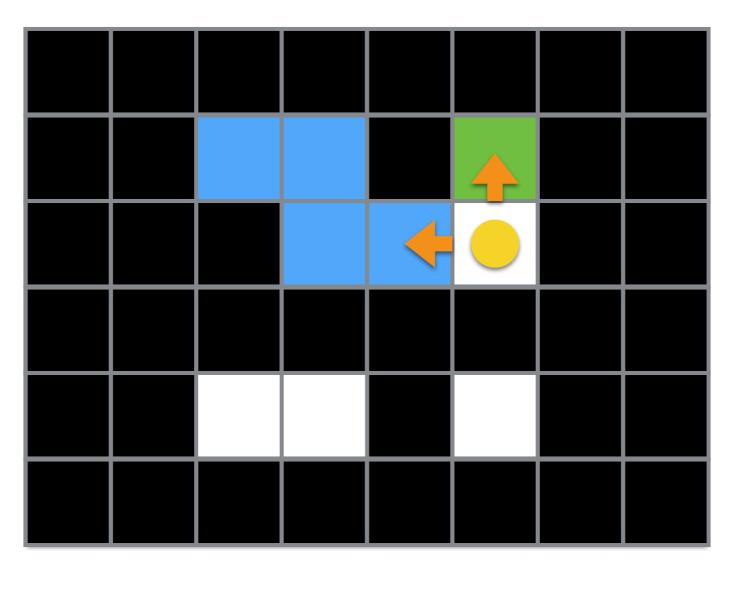




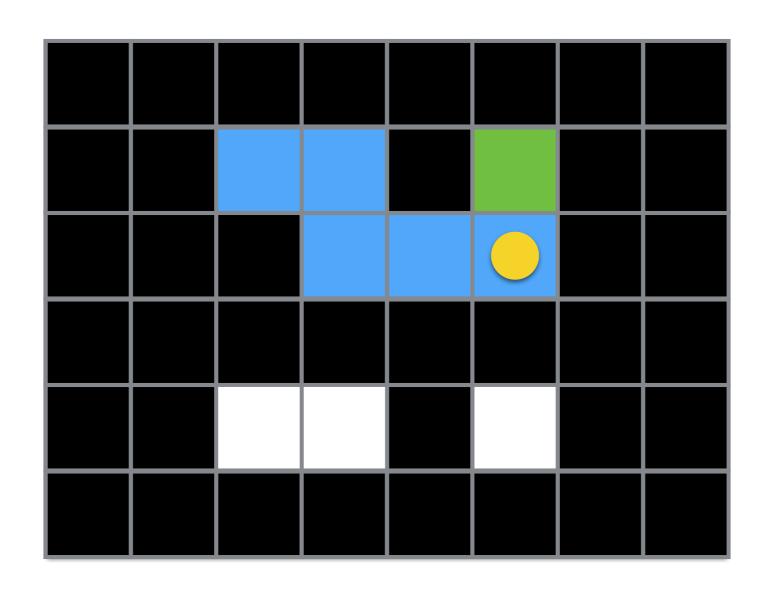


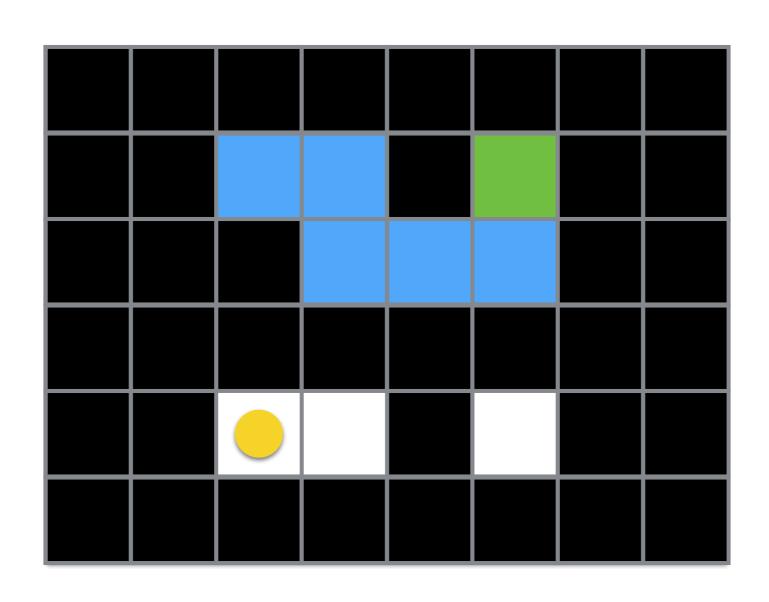
1

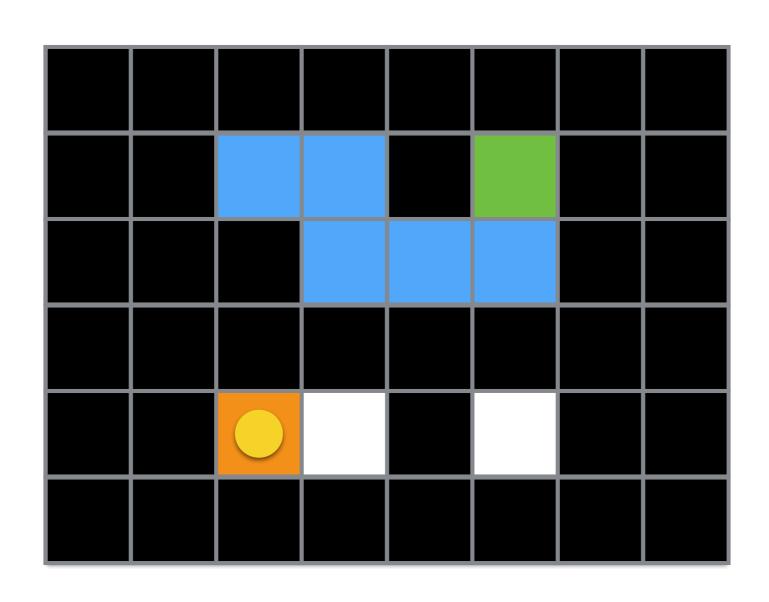
2

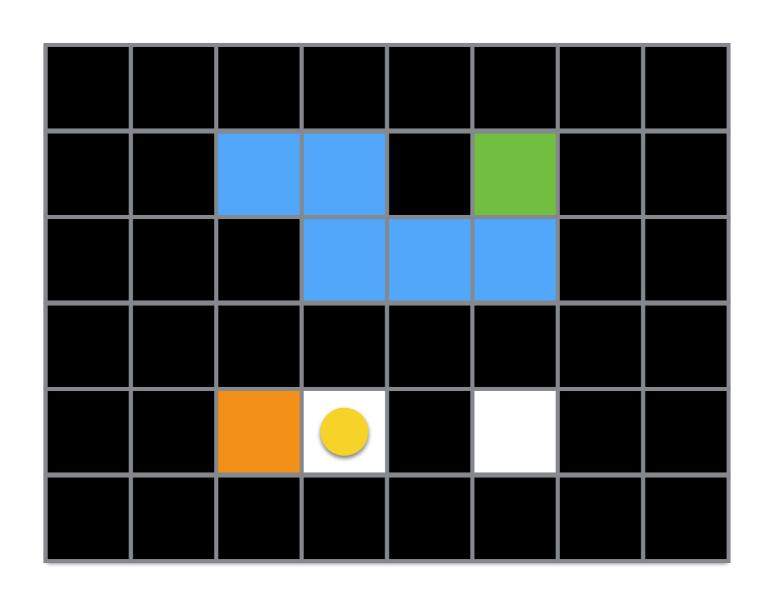


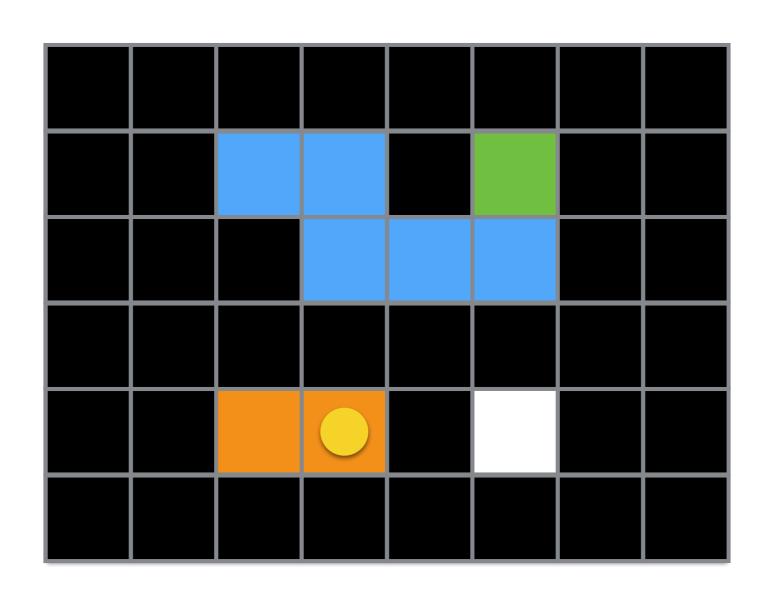
In this case, we choose the lowest label, and we store that 1 is equivalent to 2

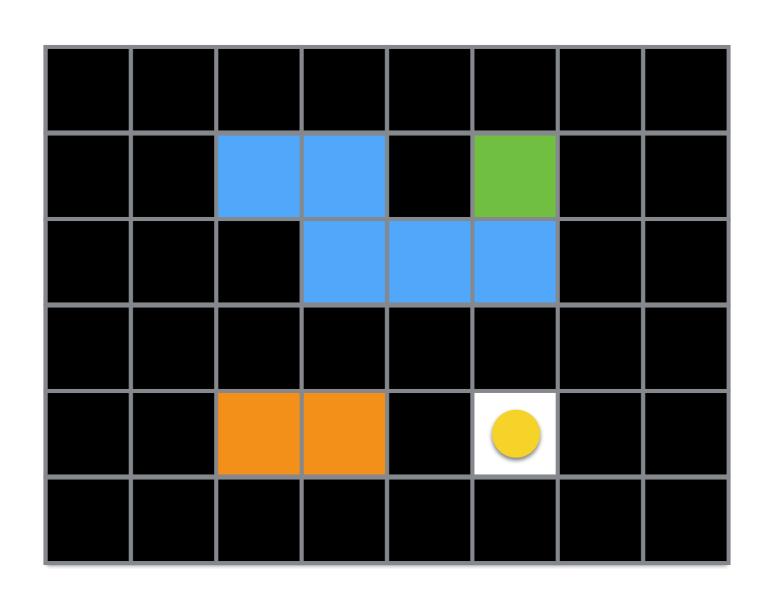


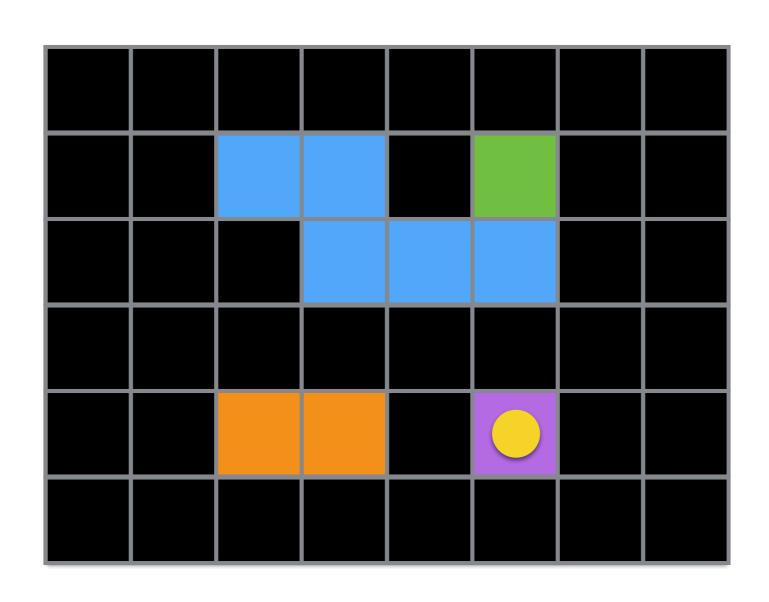




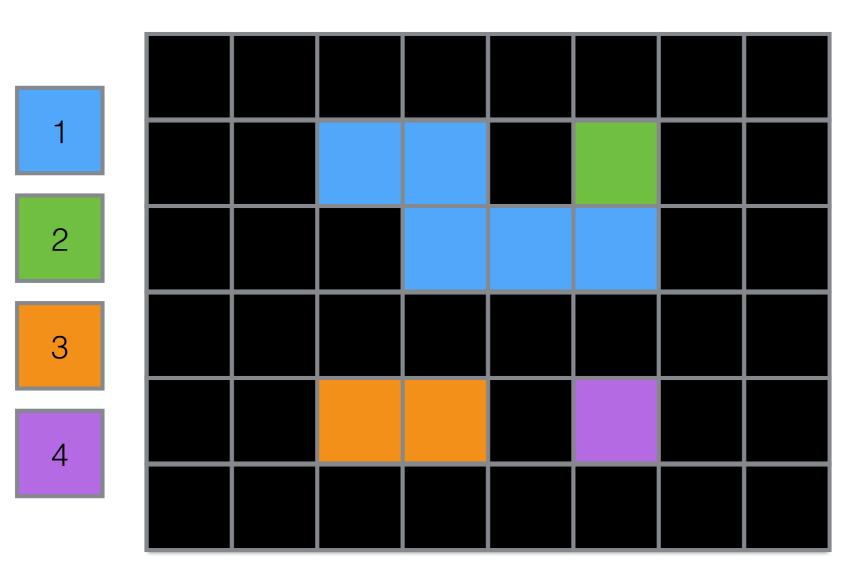




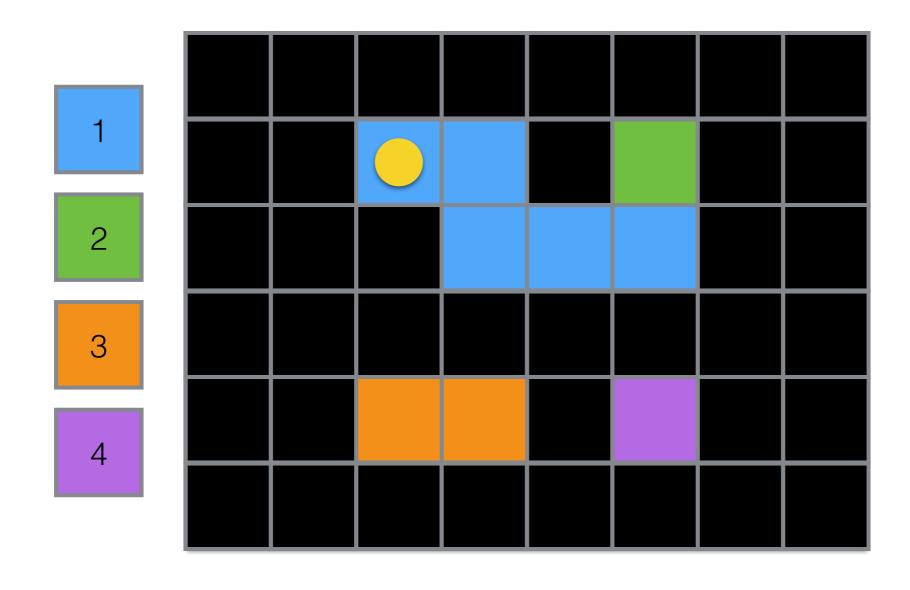


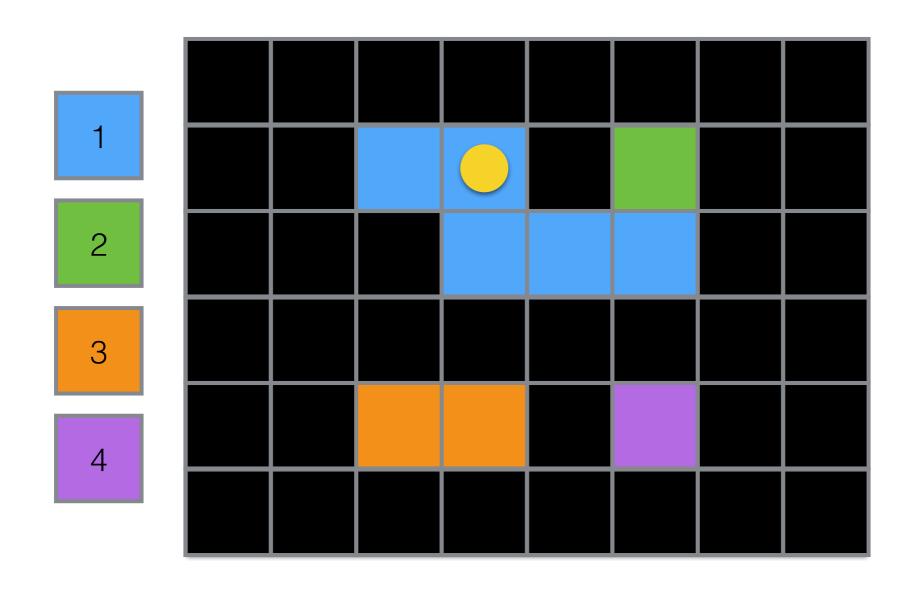


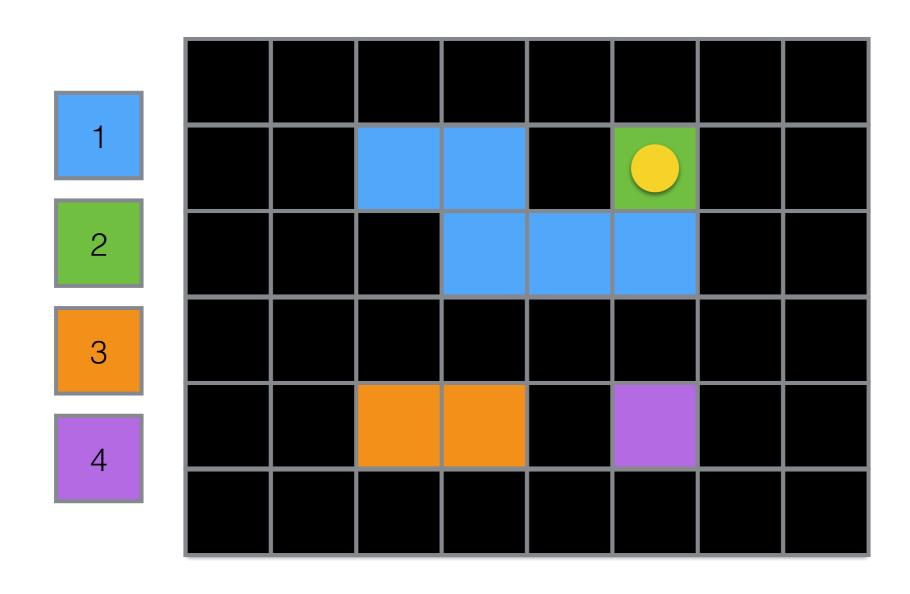
Second Pass

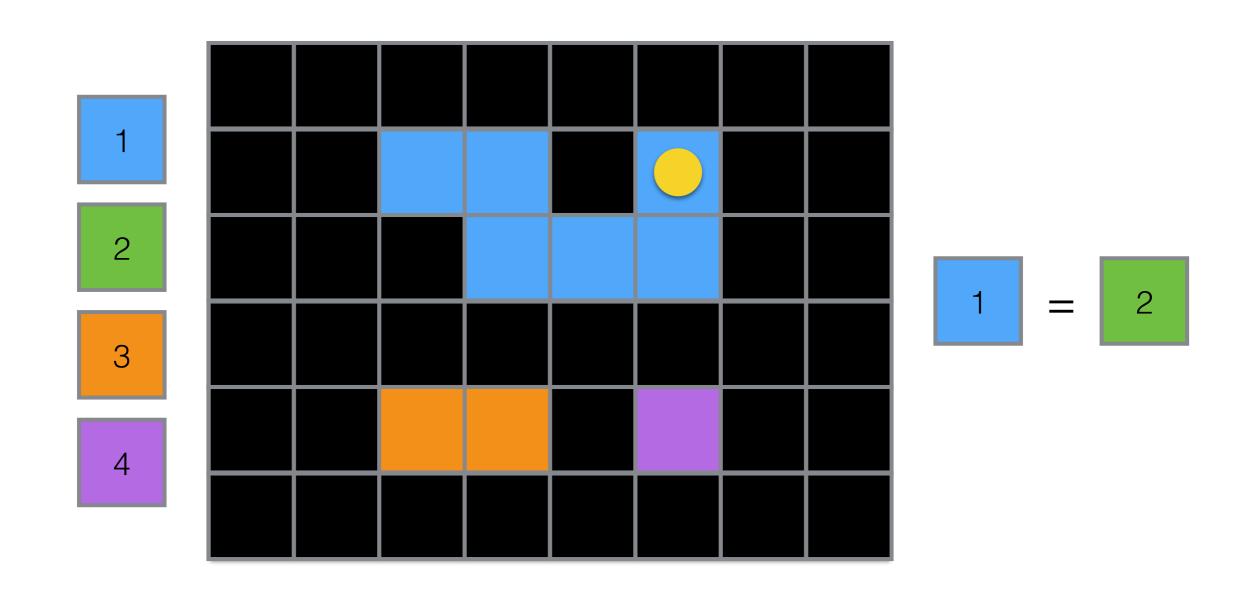


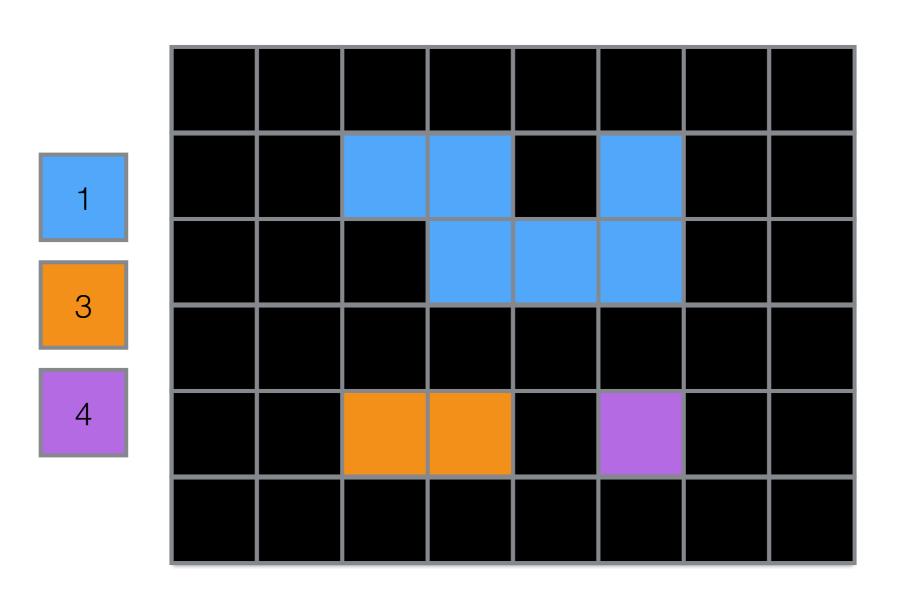
We go through all pixels. For each pixel we set the value of lowest equivalent.











Thresholding: Connected Components Example



Thresholding

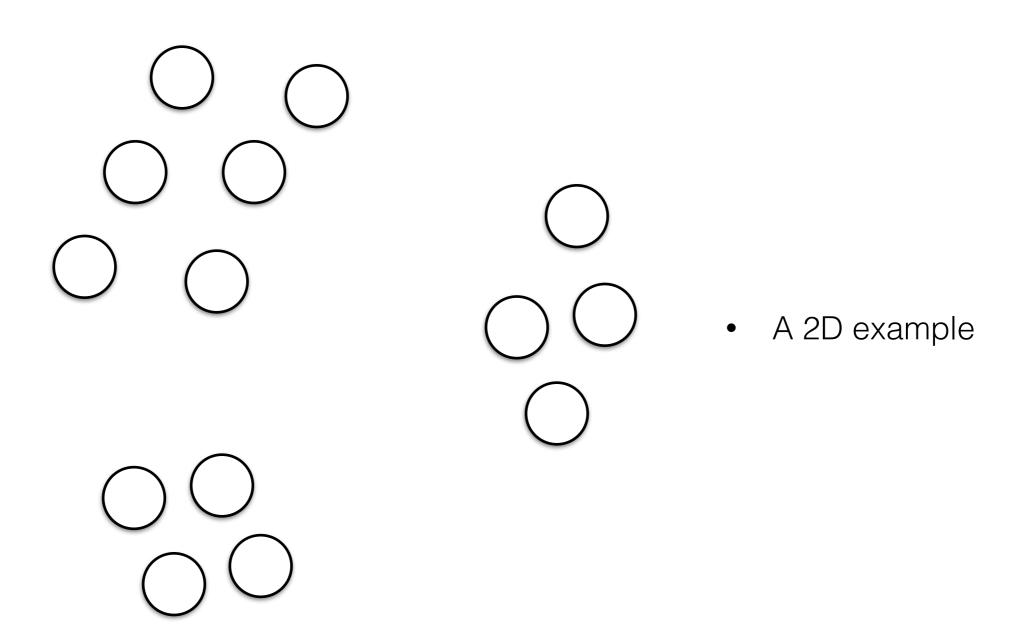
- It works if each object has a unique intensity value/ color; this is a very limiting constraint!
 - However, it could be used as a starting point for other algorithms.
- The user needs to set the threshold!
 - The I_t value for each class may be inferred by analyzing the histogram of the input image.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

k-Means

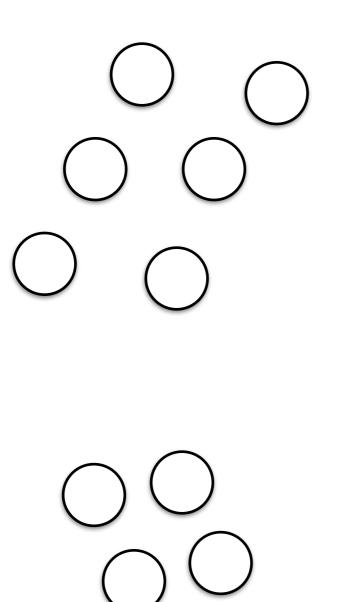
k-Means

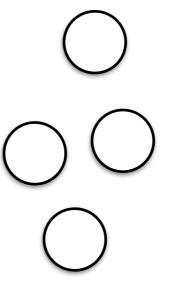
- k-means is a clustering algorithm for clustering n-D vectors/points in an n-D space:
 - A pixel with position (x, y) and intensity l is a 3D vector: $\langle x, y, l \rangle$
 - A voxel with position (x, y, z) and intensity l is a 4D vector: $\langle x, y, z, l \rangle$
- Let's assume we have k objects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

k-Means: How it Works



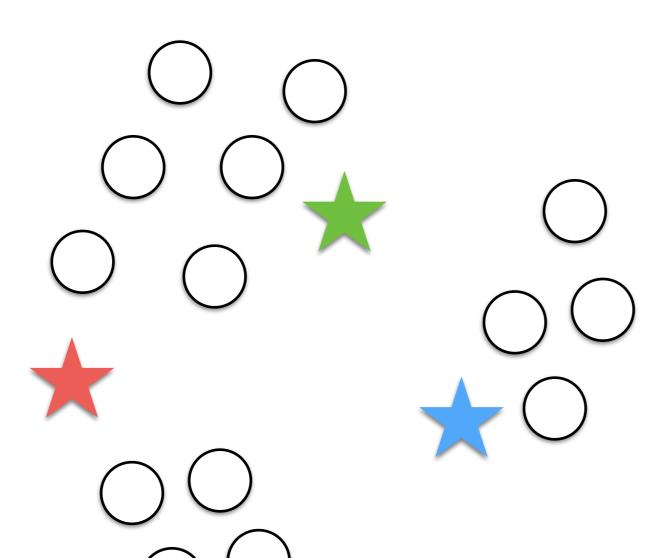
k-Means: Initialization



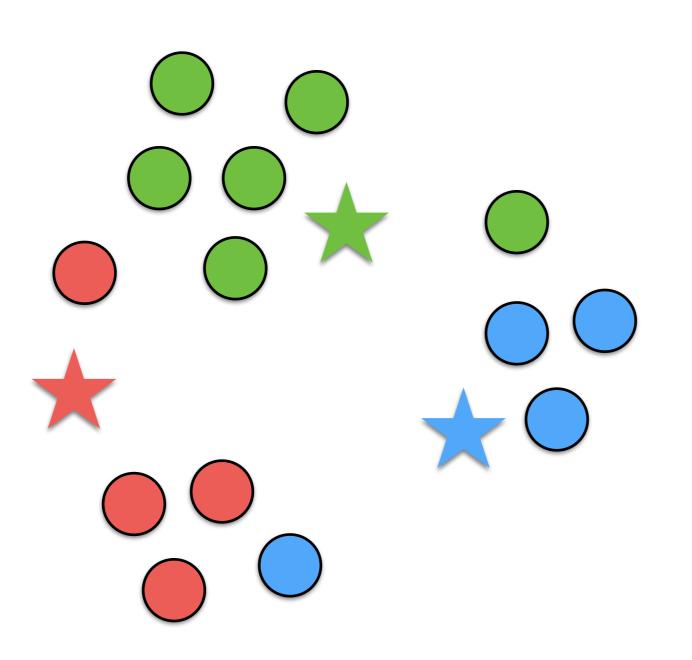


- Let's assume k=3
- We make a random guess on the k-centroids; i.e., the stars.

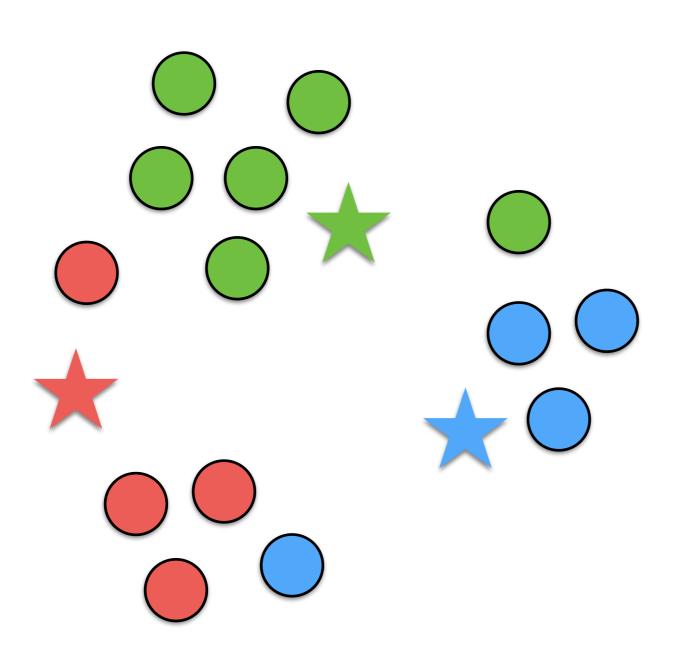
k-Means: Initialization



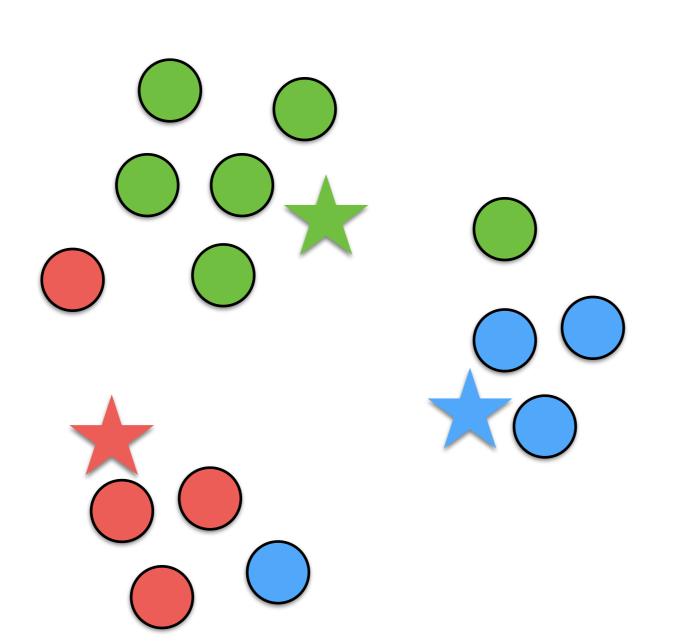
- Let's assume k=3
- We make a random guess on the k-centroids; i.e., the stars.



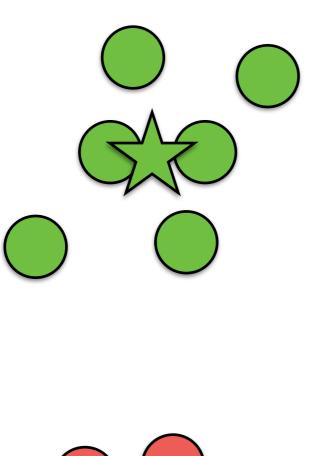
 We now assign a sample to a cluster if the distance (L1, L2, etc.), between a centroid is the minimum.

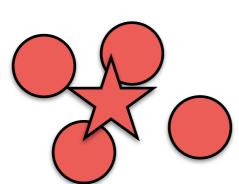


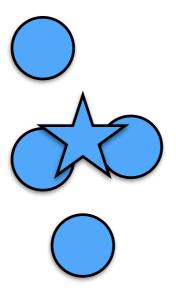
 We re-compute the centroid as the mean of samples of a cluster.



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

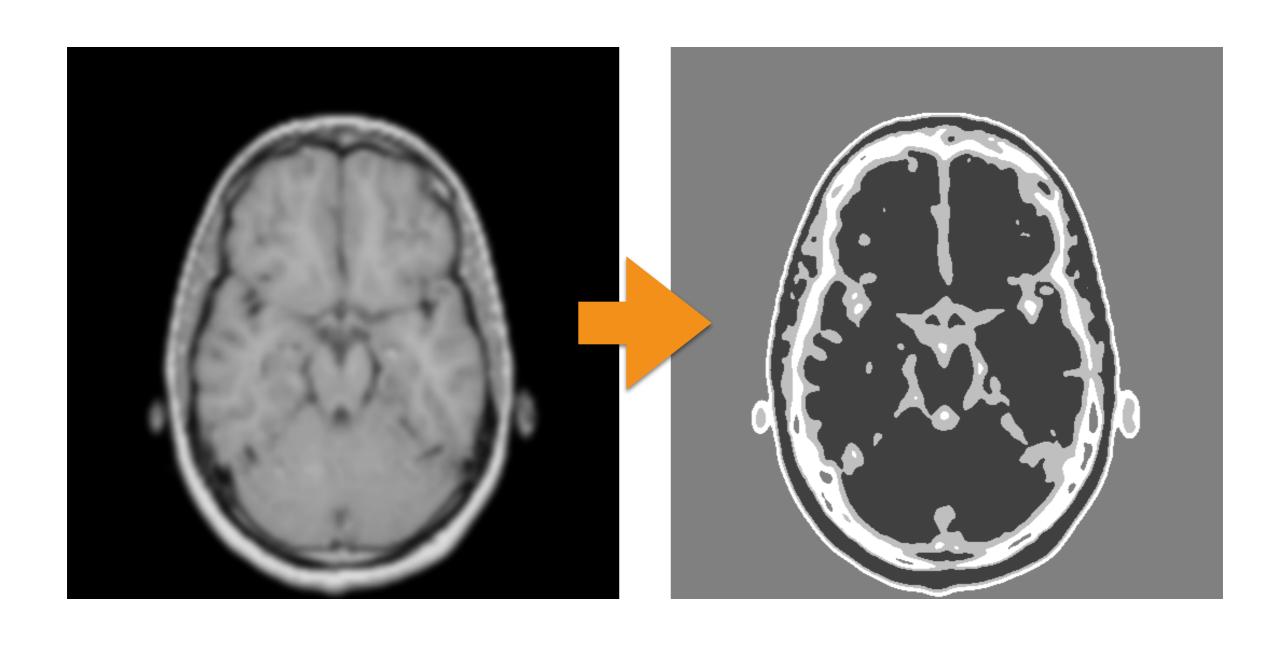




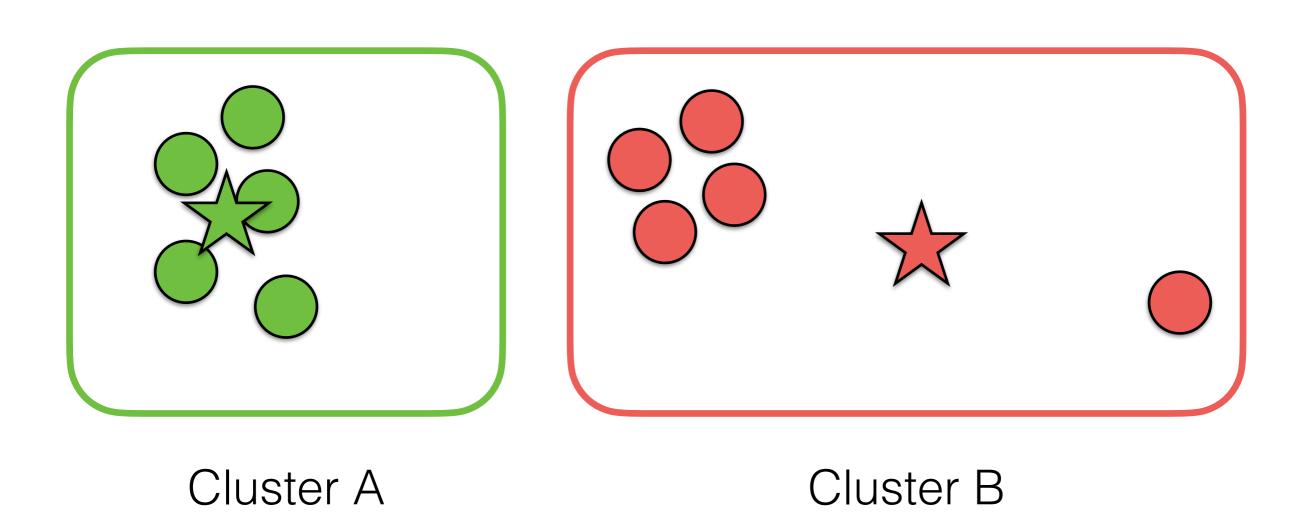


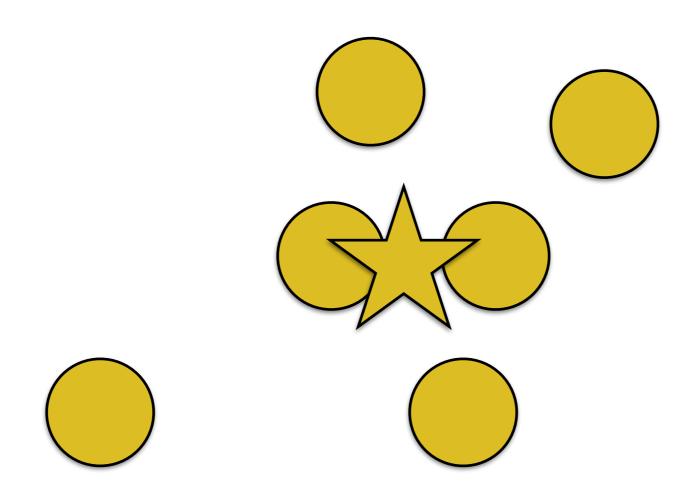
 We repeat the process until convergence (no more changes) or after m iterations.

k-Means Example

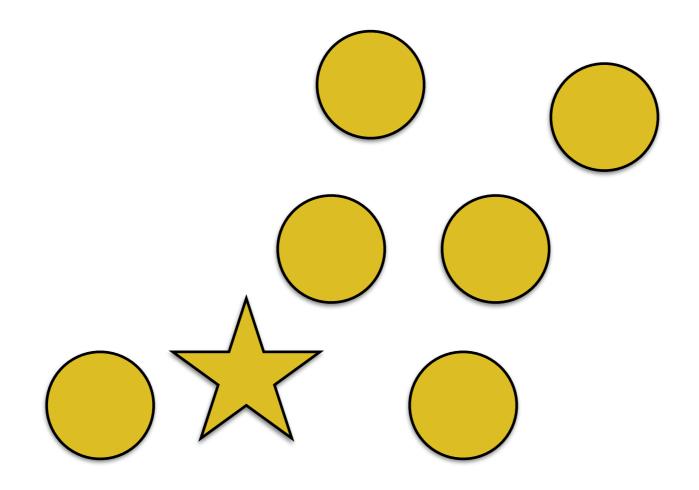


k-Means: Outliers

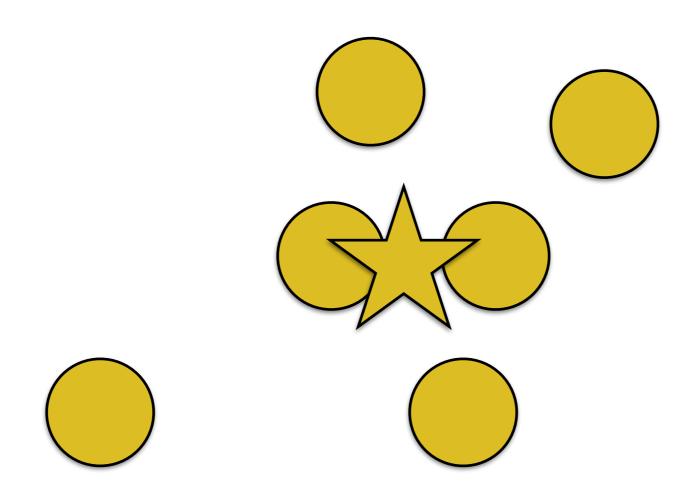




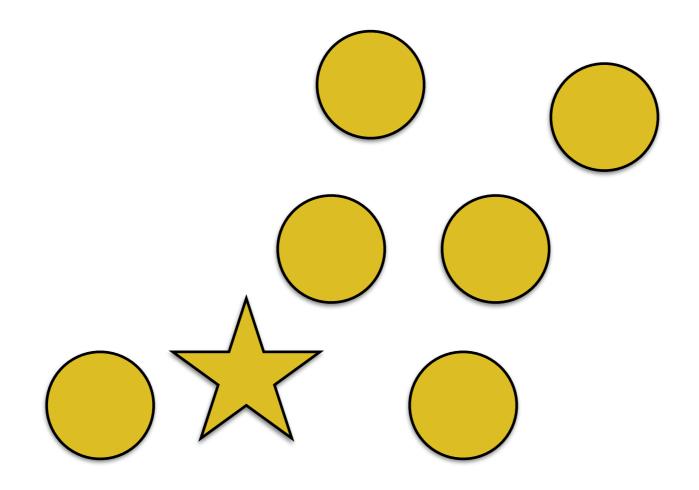
Even Iteration



Odd Iteration



Even Iteration



Odd Iteration

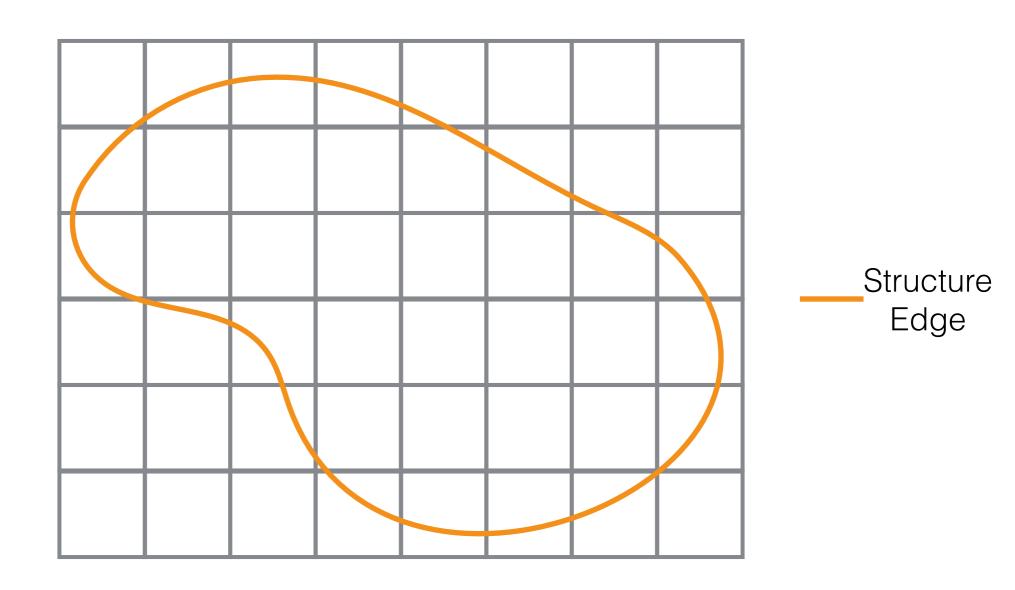
k-Means: Advantages

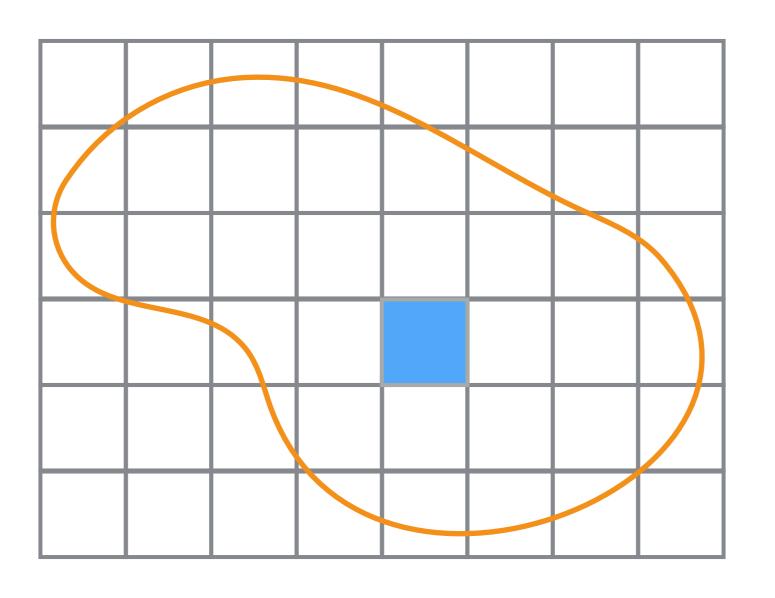
- The method is fully automatic
- This works for 2D and 3D volumes
- This can "understand" neighbors in an implicit way

k-Means: Disadvantages

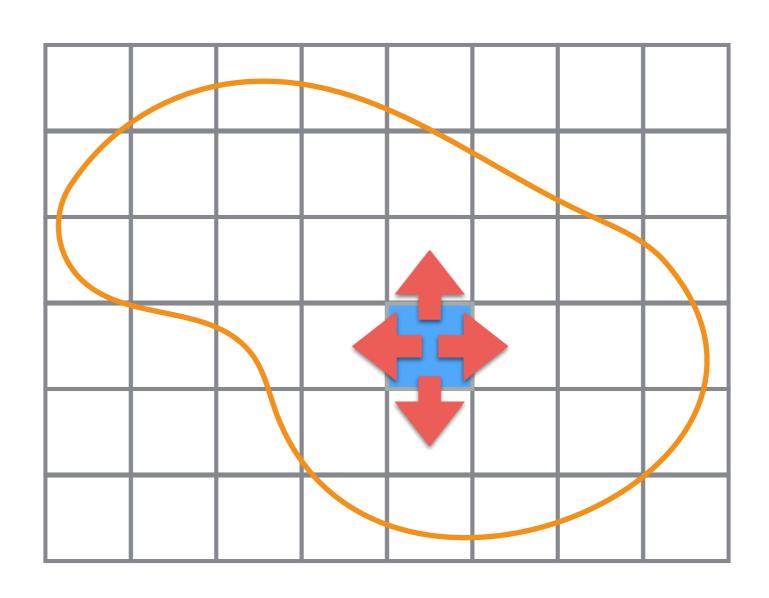
- We need to know how many objects (including the background) are in the image:
 - We may run k-means multiple times until a certain criterion is met (e.g., reaching the 80% of percentage of explained variance)
- Outliers:
 - better initialization (sampling)
- The method may not converge
 - we need to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

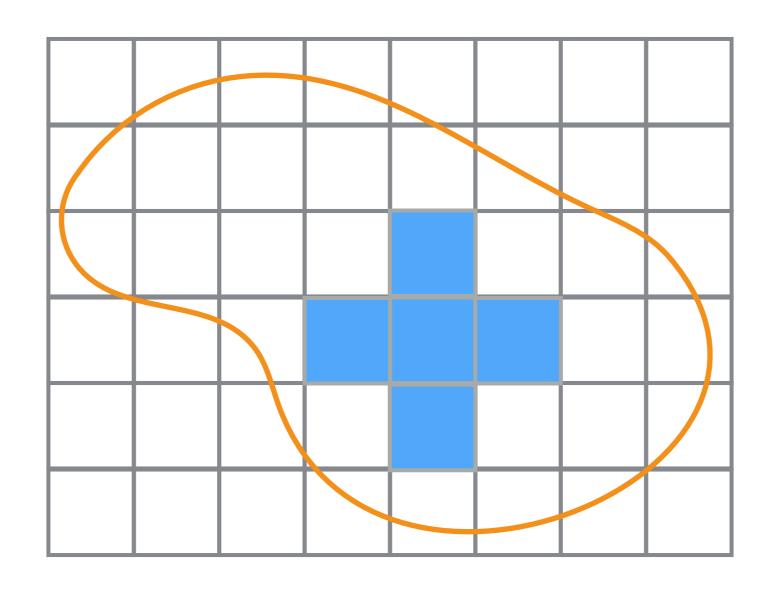
- This algorithms expands a painted initial mask until it reaches strong edges
- Therefore, we need to compute edges first!



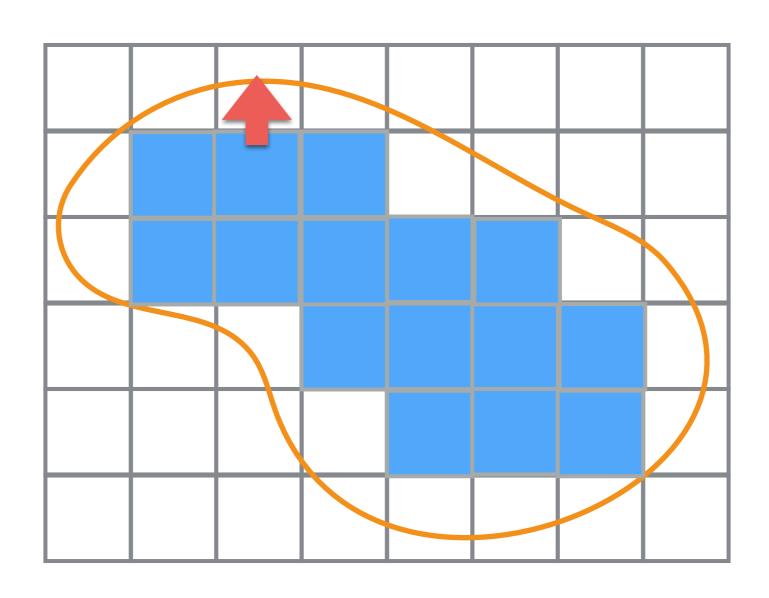


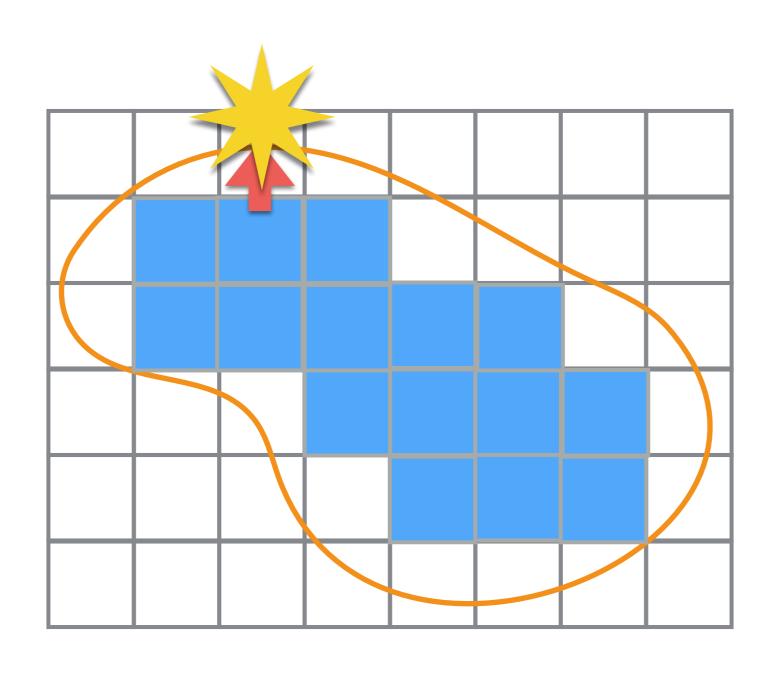
Seed





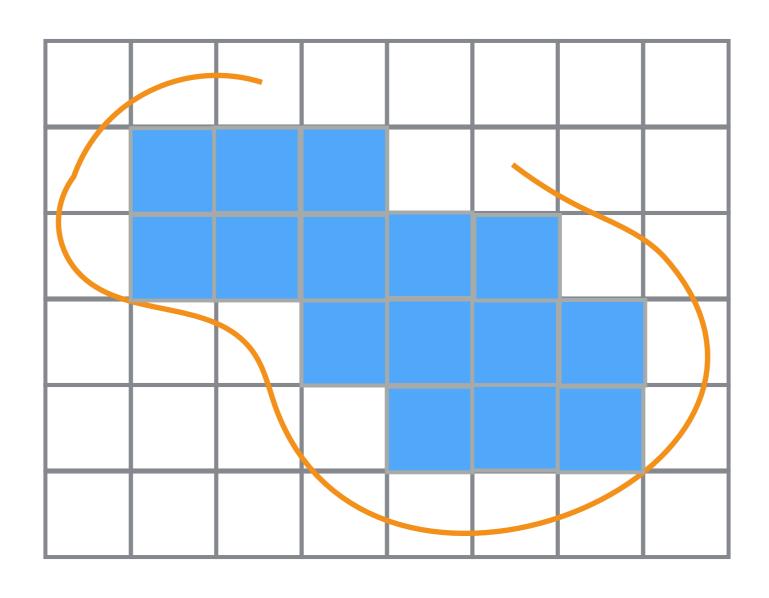
after a while...



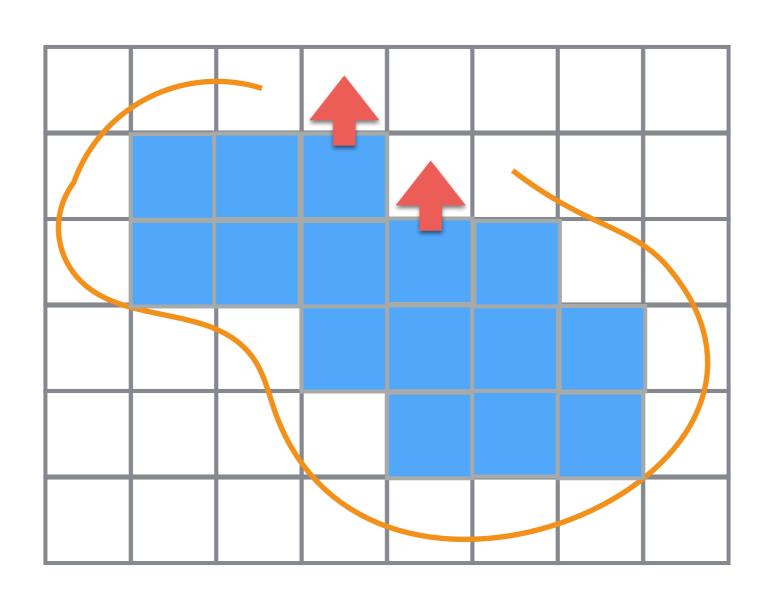


- It is straightforward to extend to 3D!
- This algorithm depends on:
 - The threshold of edge detection
- It may be slow:
 - From an initial seed, the growing region needs to reach the farthest edge pixel/voxel.
 - Computational complexity is a function of the area/ volume of the object we want to segment.

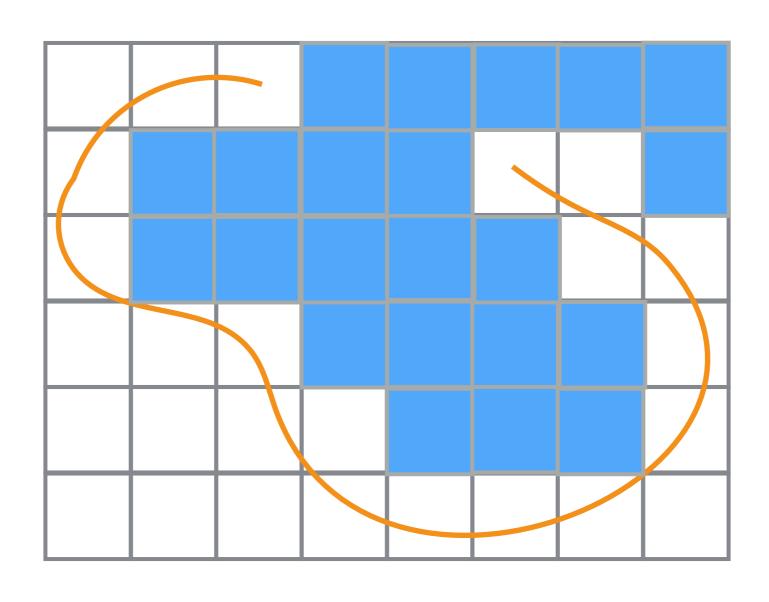
Region Growing: Epic Fail



Region Growing: Epic Fail



Region Growing: Epic Fail



Active Contour Model aka Snakes

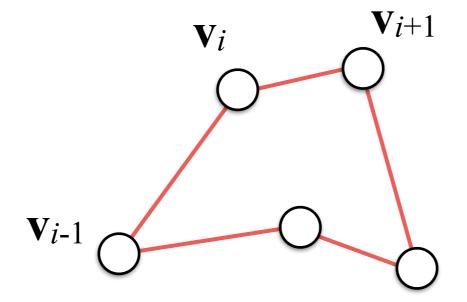
A snake is a parametric curve:

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

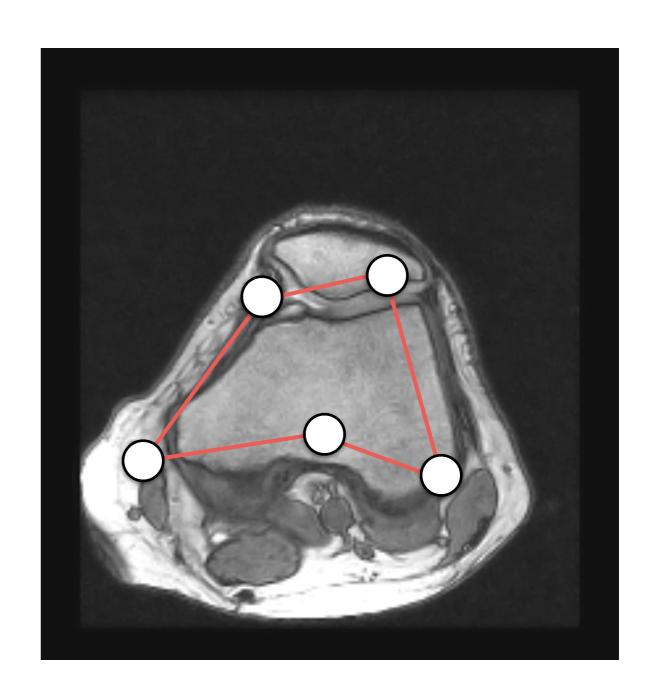
 Typically, it is a spline (original paper), but for sake of simplicity let's assume a piecewise linear curve.

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

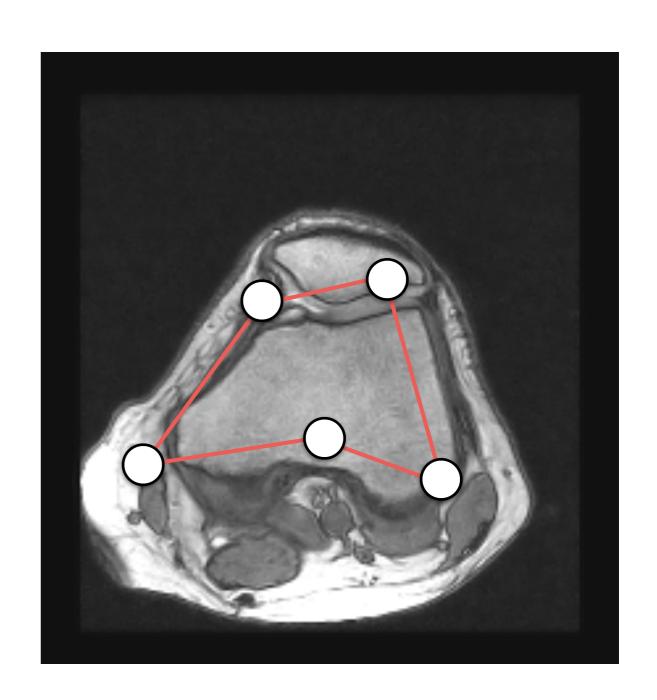
$$C = \{v_i | i \in [1,n]\}$$
 $v_i = [x_i, y_i]$



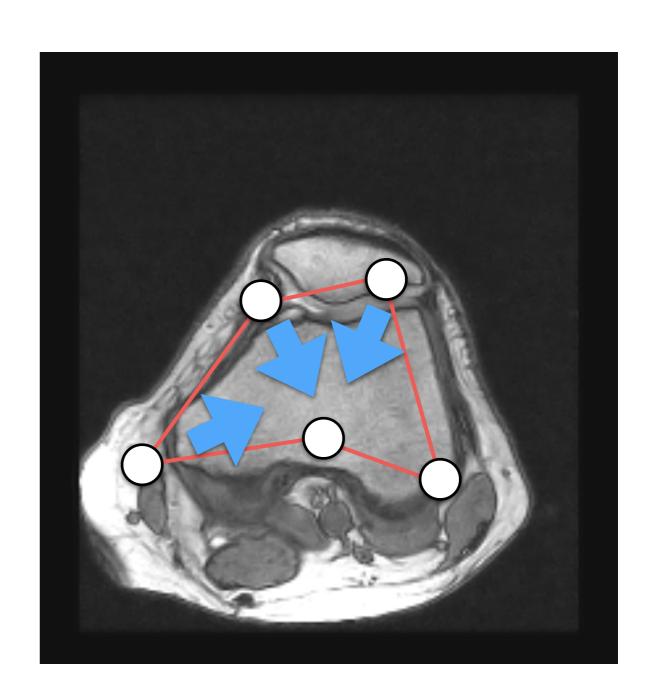
 A first step, we draw a snake close to the boundary of the object we want to segment.



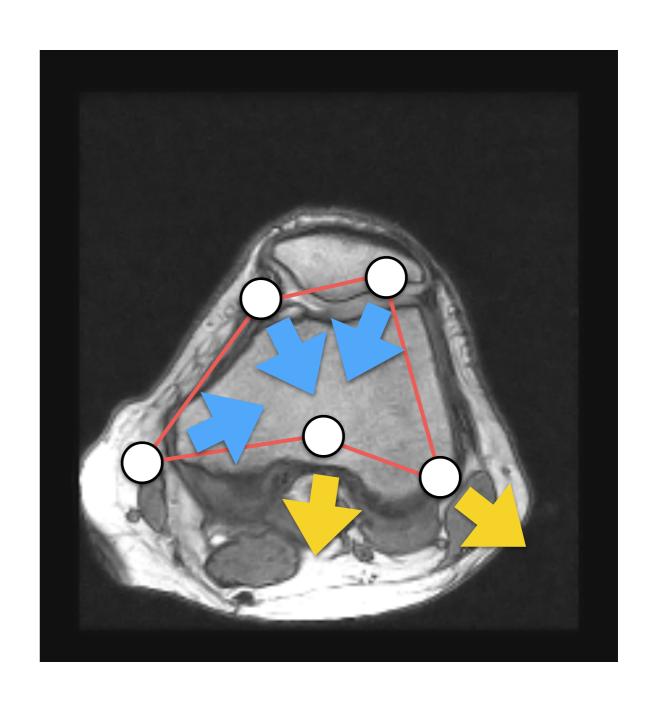
 Then, we *deform* its control points in order to move them towards the object's boundary.



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



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



- 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.

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

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

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

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

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

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

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

Elasticity

$$E_{\text{internal}} = \int_{0}^{1} E_{\text{internal}}(\mathbf{v}(t)) dt$$

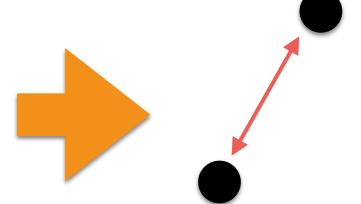
 This energy represents the internal energy of the cure due to bending. It is defined per point as

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

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

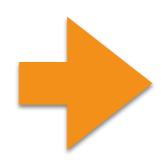
The first term is an elastic energy:

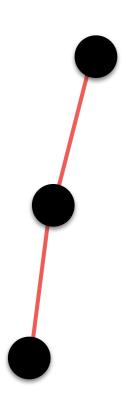
$$\frac{d\mathbf{v}(t)}{dt} \approx \mathbf{v}_{i+1} - \mathbf{v}_i$$



The second term is a bending energy:

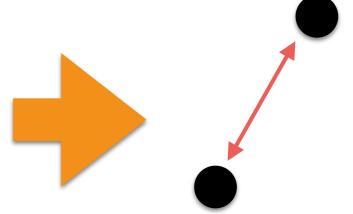
$$\frac{d^2\mathbf{v}(t)}{d^2t} \approx \mathbf{v}_{i+1} - 2\mathbf{v}_i + \mathbf{v}_{i-1}$$





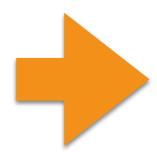
The first term is an elastic energy:

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



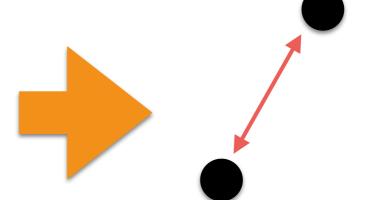
The second term is a bending energy:

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



• The first term is an elastic energy:

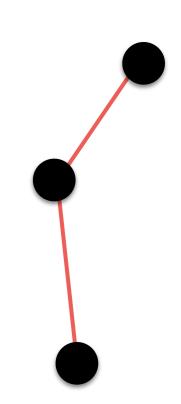
$$\frac{d\mathbf{v}(t)}{dt} \approx \mathbf{v}_{i+1} - \mathbf{v}_i$$



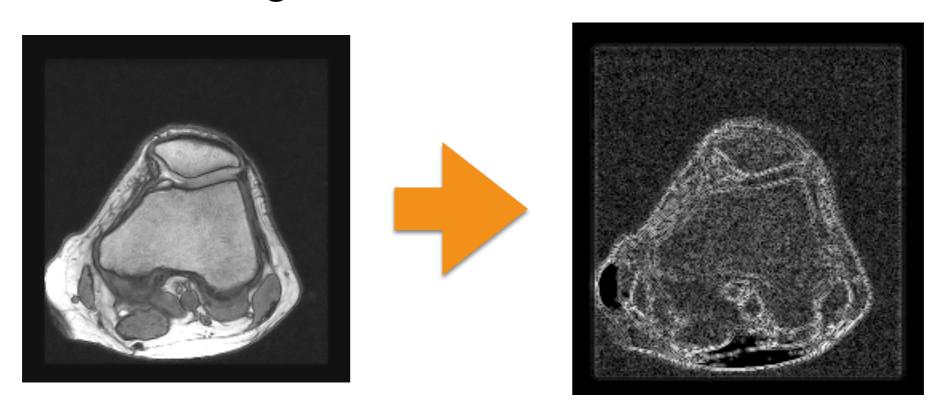
The second term is a bending energy:

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





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

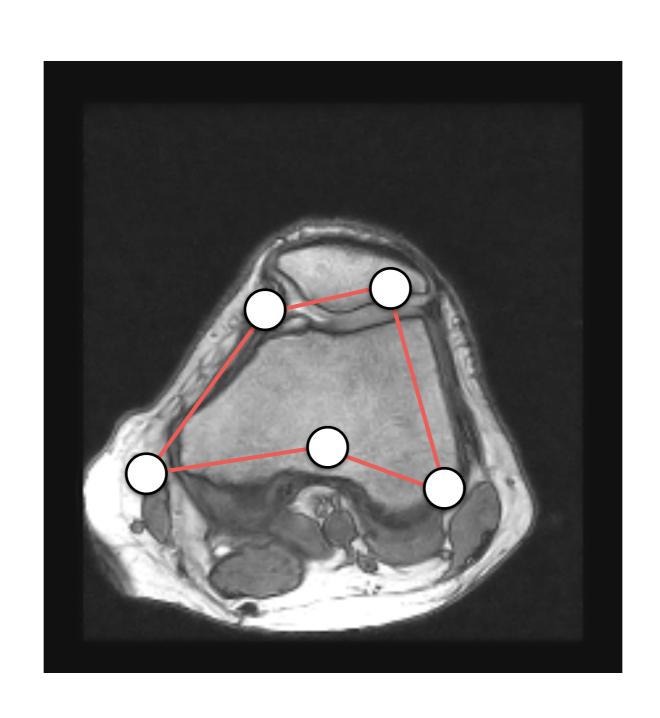


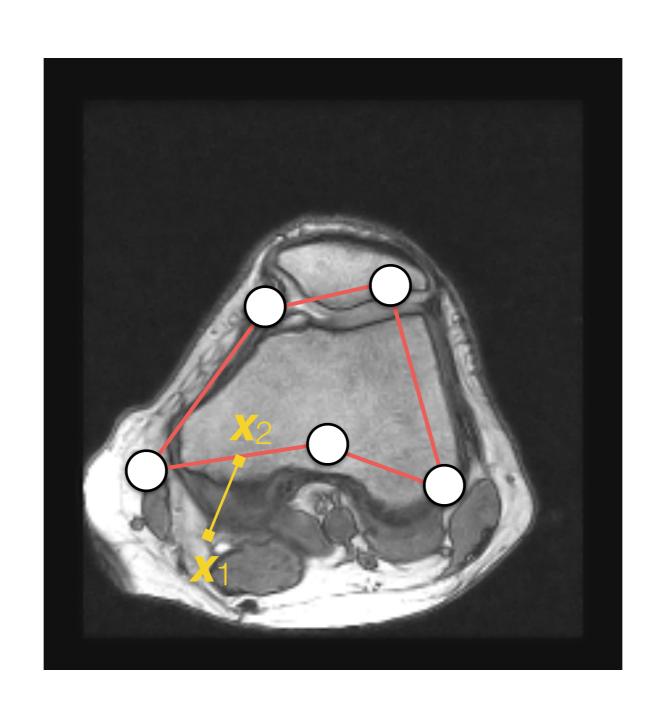
It is defined per point as

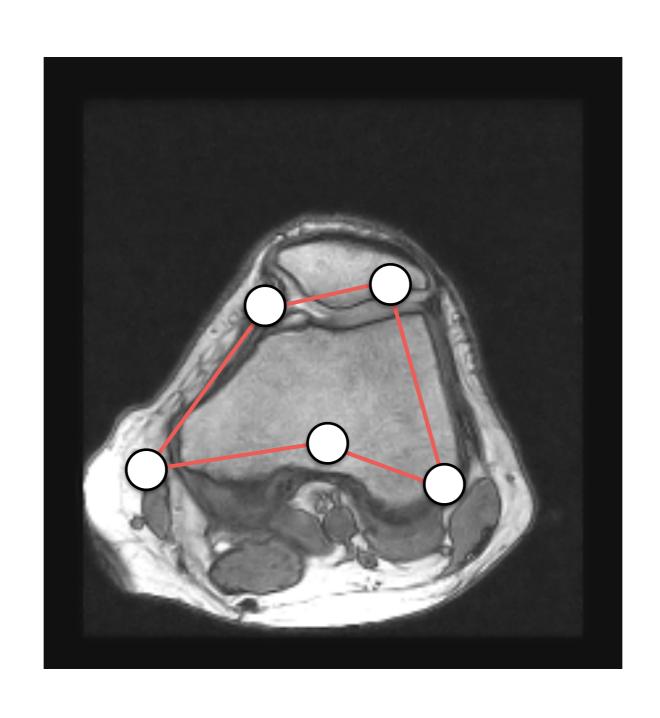
$$E_{\text{external}}(\mathbf{v}(t)) = -\|\nabla I(\mathbf{v}(t))\|^2$$

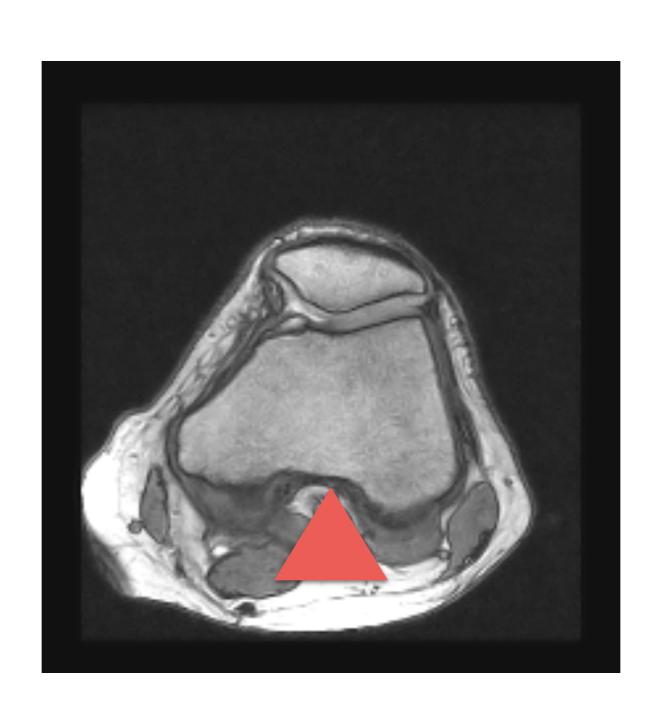
$$E_{\text{external}} = \int_0^1 E_{\text{external}}(\mathbf{v}(t))dt$$

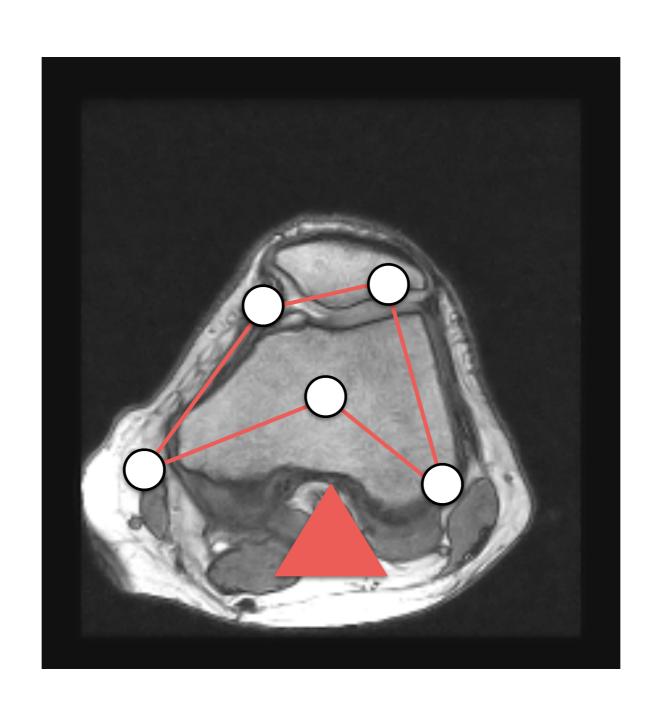
- This energy is meant for interactive systems.
- The user interactively monitors the minimization, and she/he can push/pull vertices using the mouse cursor's position:
 - Repulsion forces or "vulcano": $\frac{1}{r^2}$
 - Spring forces: $-k\|\mathbf{x}_1 \mathbf{x}_2\|^2$









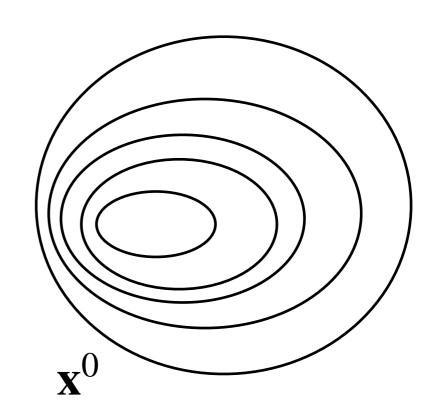


How do we solve E?

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

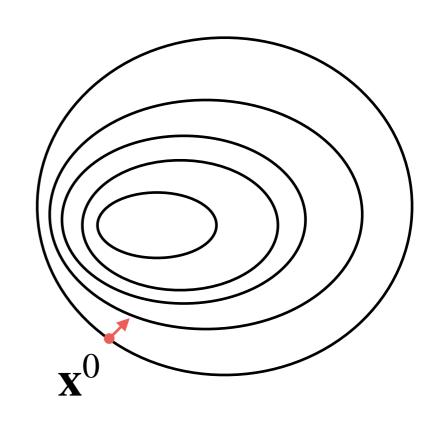
$$\mathbf{x}_{j}^{i+1} = \mathbf{x}_{j}^{i} - \alpha \frac{\partial}{\partial \mathbf{x}_{j}} f(\mathbf{x}^{i})$$

- We need to start with a g
- It will find a local minimum!
- f has to be differentiable.
- \mathbf{x}^0 is a "good" guess.



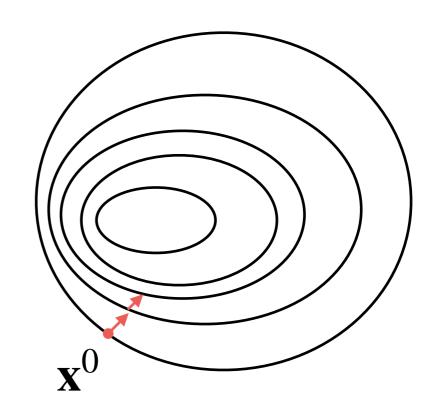
$$\mathbf{x}_{j}^{i+1} = \mathbf{x}_{j}^{i} - \alpha \frac{\partial}{\partial \mathbf{x}_{j}} f(\mathbf{x}^{i})$$

- We need to start with a g
- It will find a local minimum!
- f has to be differentiable.
- \mathbf{x}^0 is a "good" guess.



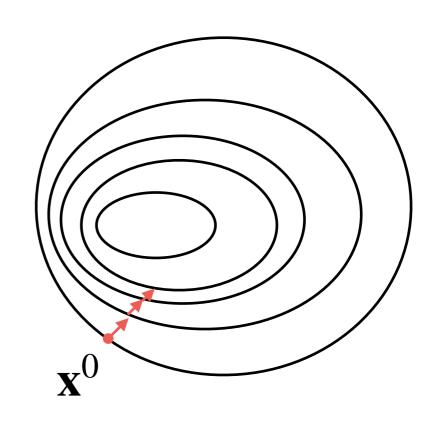
$$\mathbf{x}_{j}^{i+1} = \mathbf{x}_{j}^{i} - \alpha \frac{\partial}{\partial \mathbf{x}_{j}} f(\mathbf{x}^{i})$$

- We need to start with a g
- It will find a local minimum!
- f has to be differentiable.
- \mathbf{x}^0 is a "good" guess.



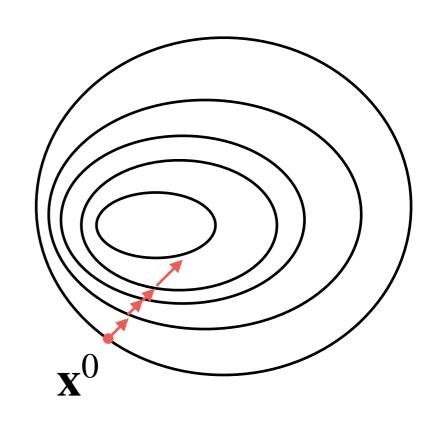
$$\mathbf{x}_{j}^{i+1} = \mathbf{x}_{j}^{i} - \alpha \frac{\partial}{\partial \mathbf{x}_{j}} f(\mathbf{x}^{i})$$

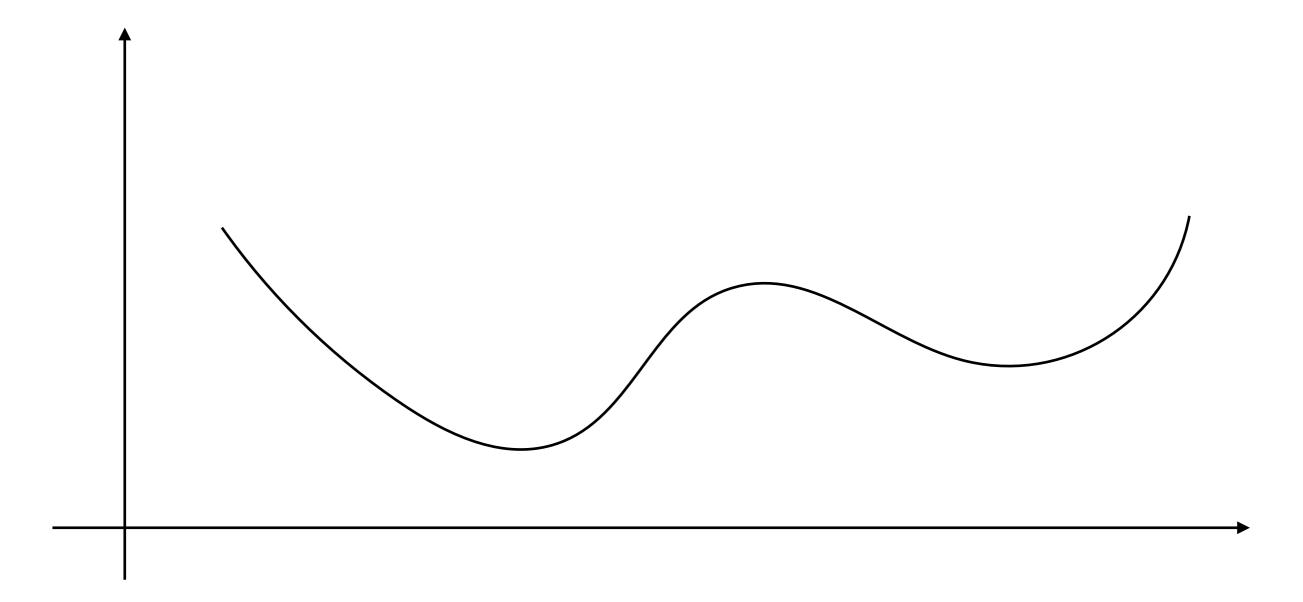
- We need to start with a g
- It will find a local minimum!
- f has to be differentiable.
- \mathbf{x}^0 is a "good" guess.

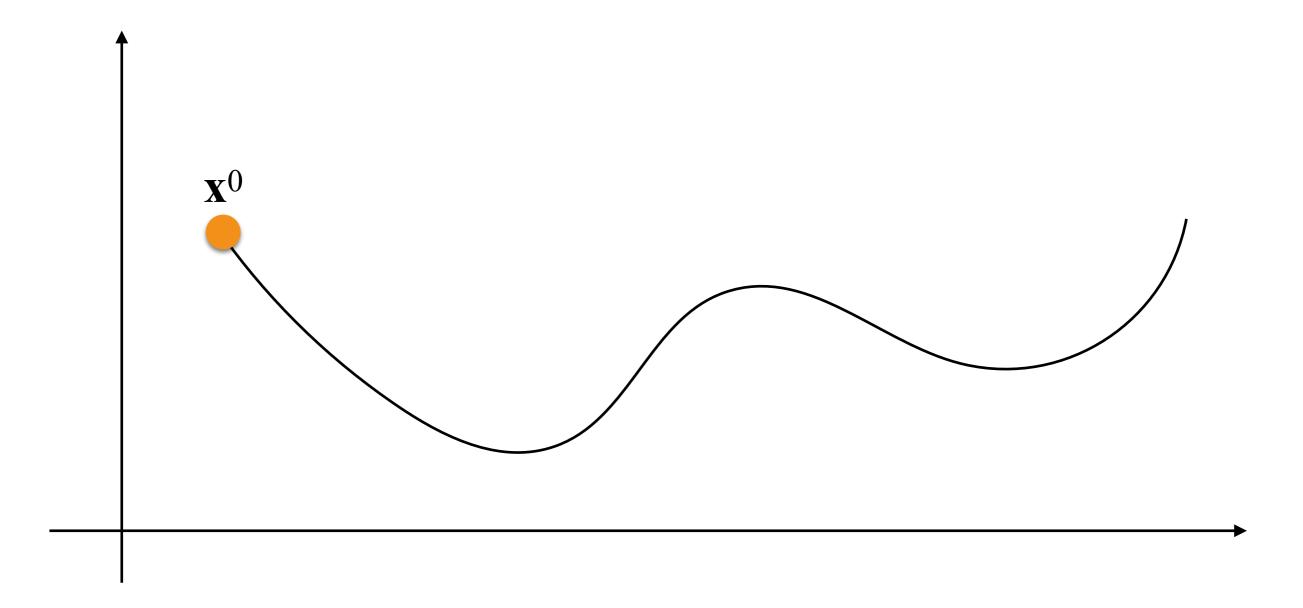


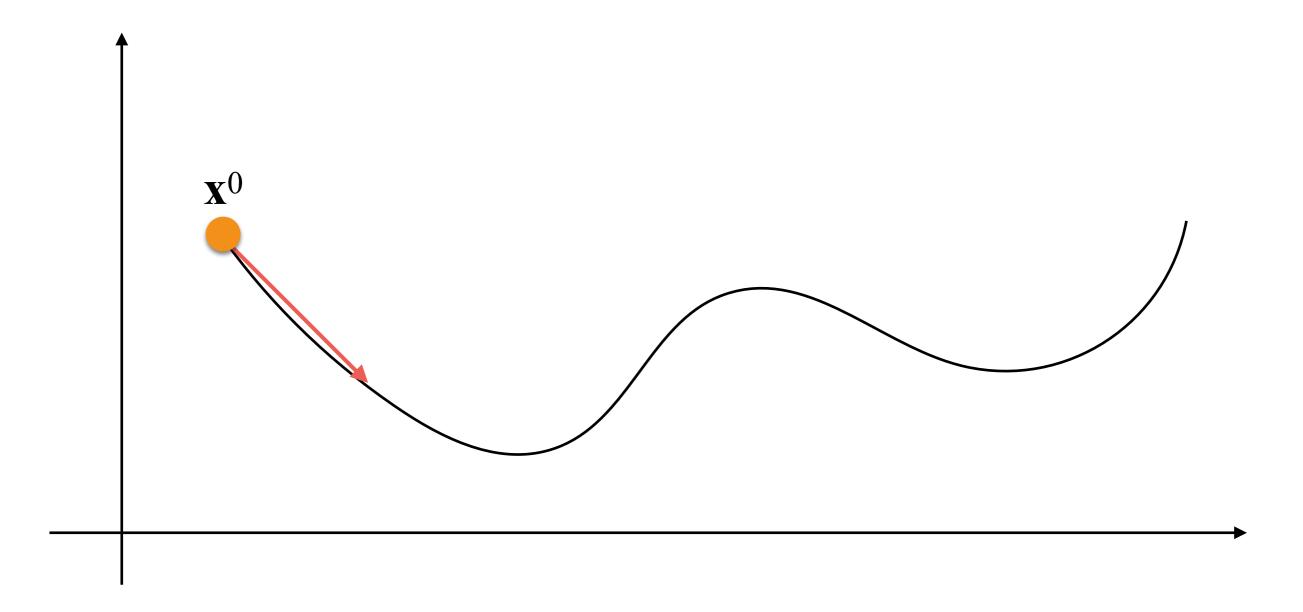
$$\mathbf{x}_{j}^{i+1} = \mathbf{x}_{j}^{i} - \alpha \frac{\partial}{\partial \mathbf{x}_{j}} f(\mathbf{x}^{i})$$

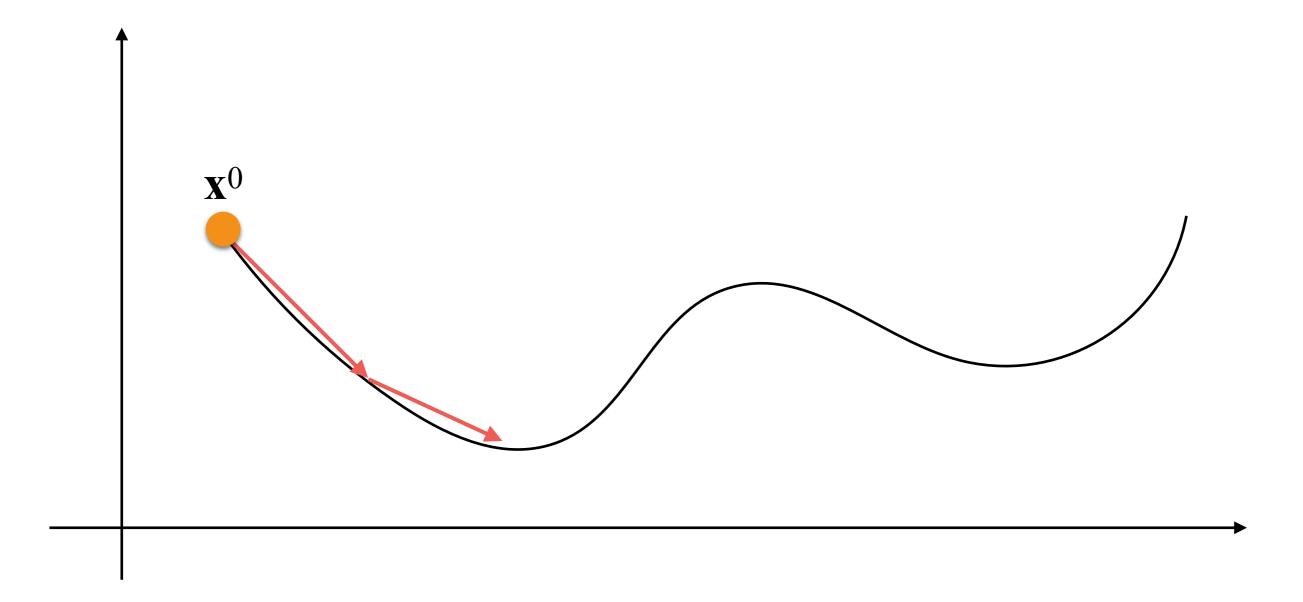
- We need to start with a g
- It will find a local minimum!
- f has to be differentiable.
- \mathbf{x}^0 is a "good" guess.

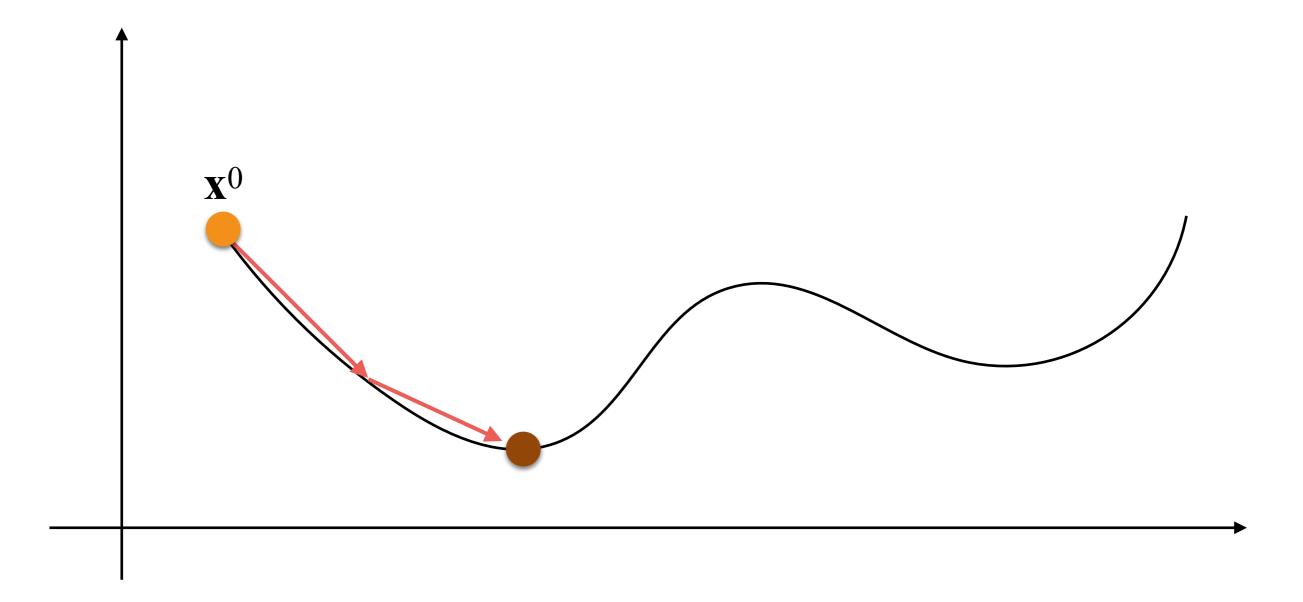


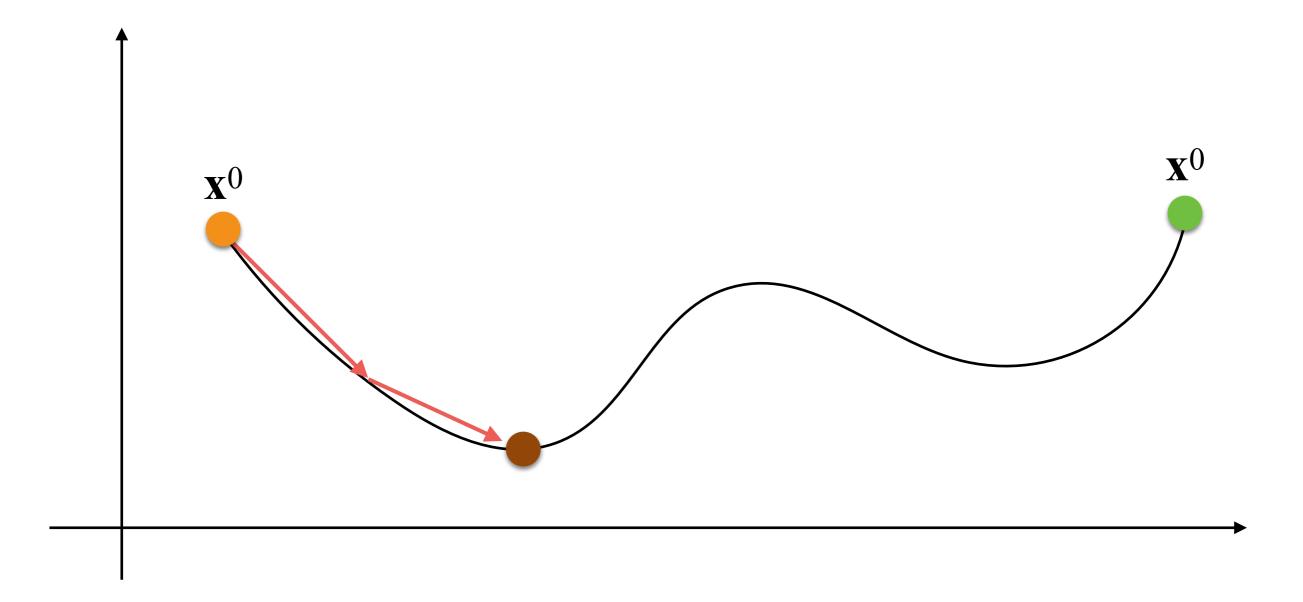


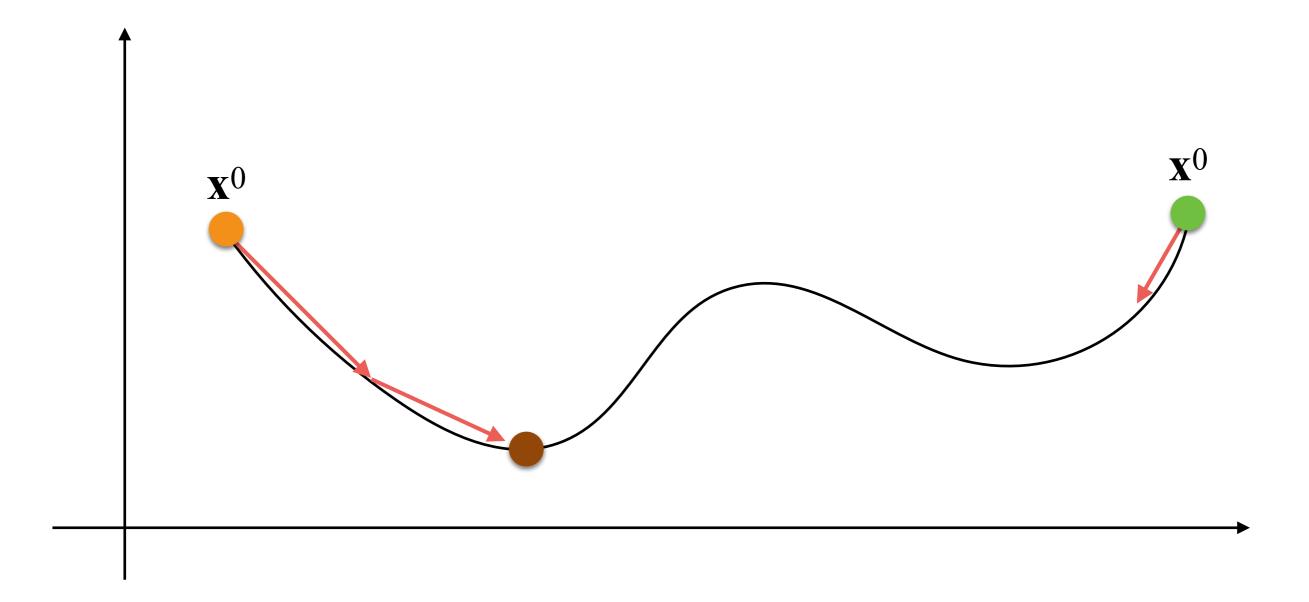


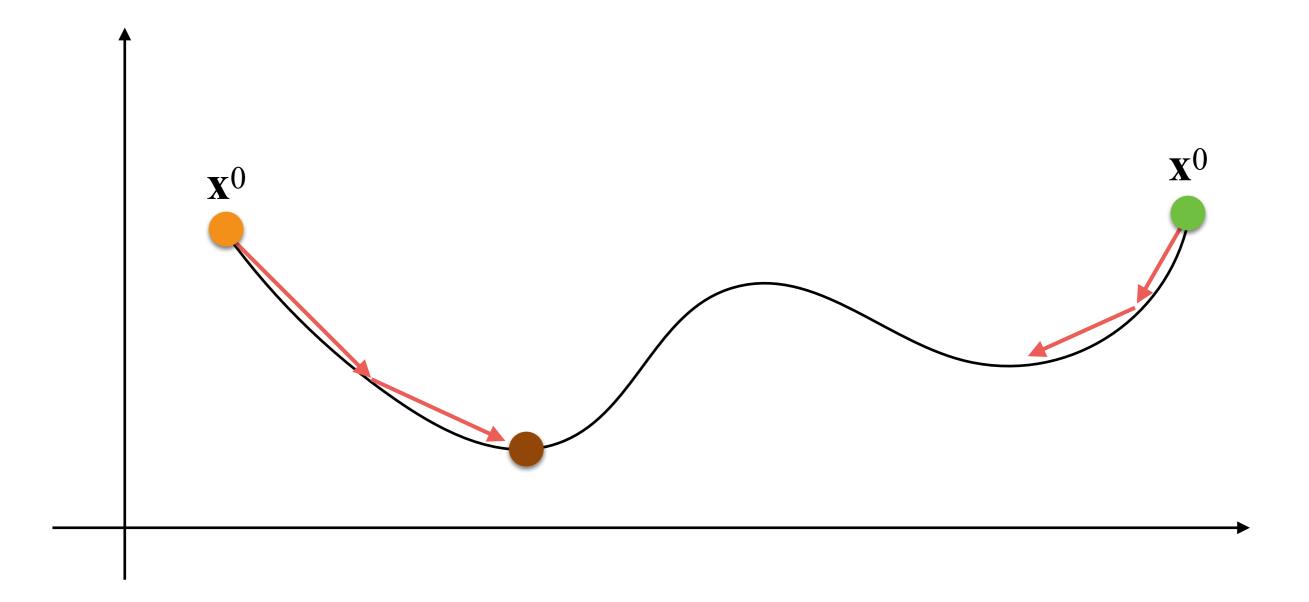


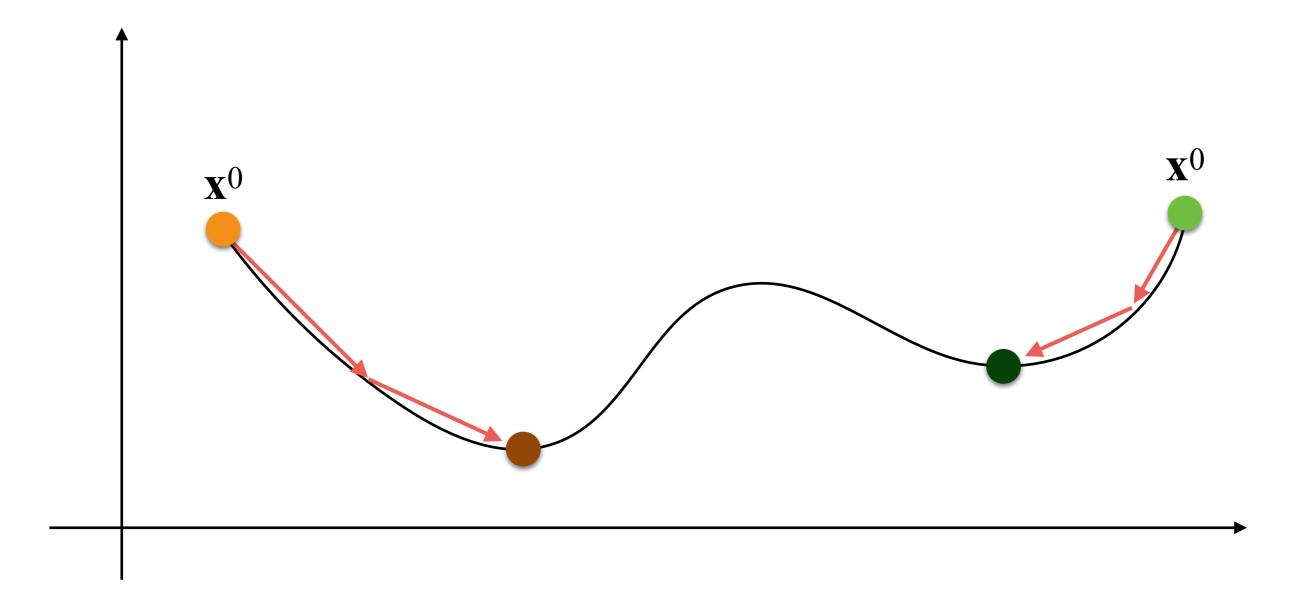








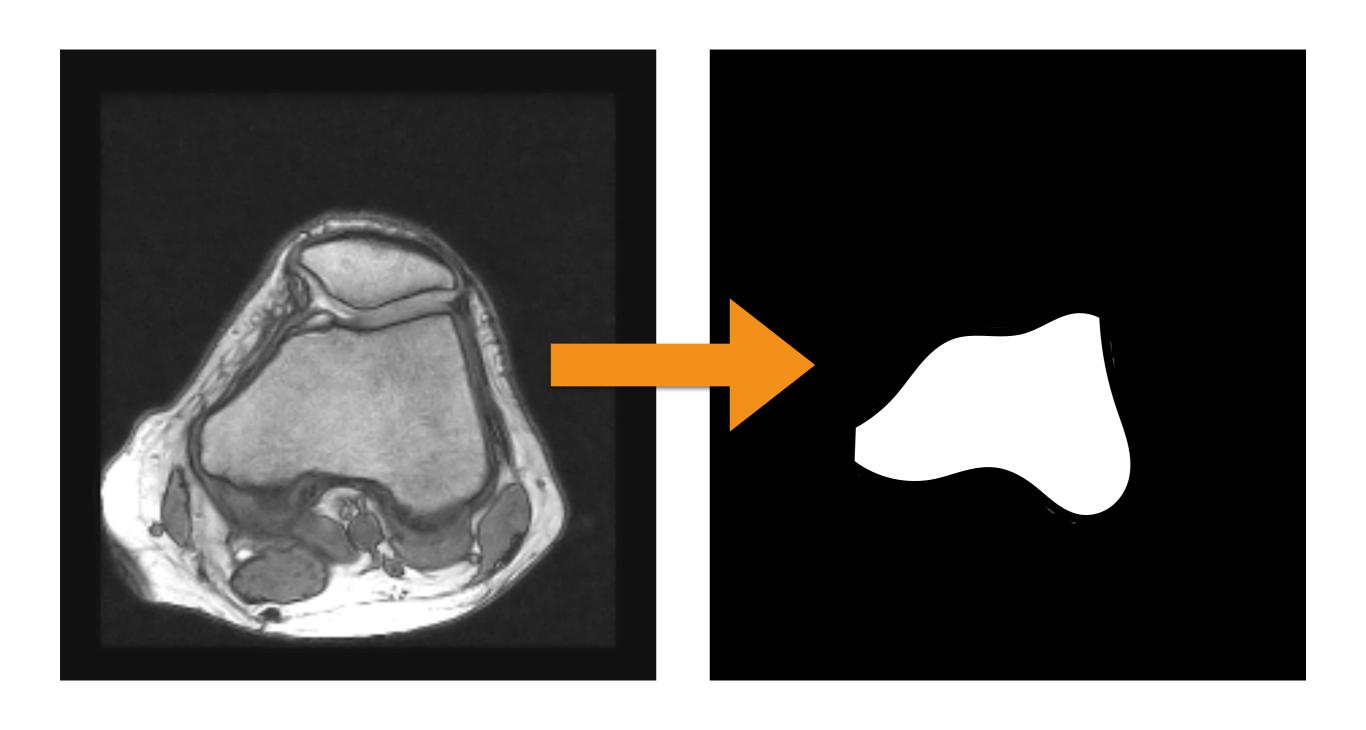




Snakes: Gradient Descent

- 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!

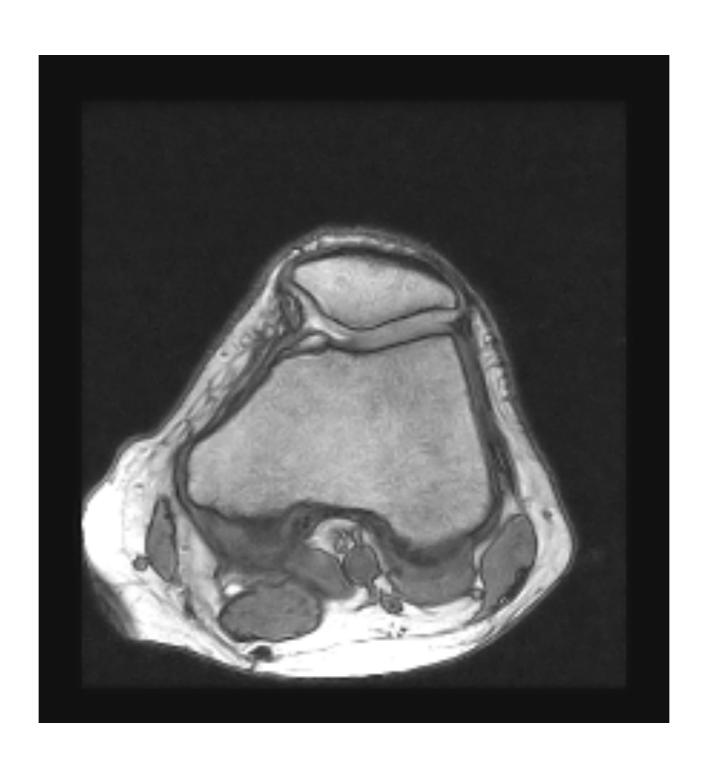
Snakes An Example



Snakes

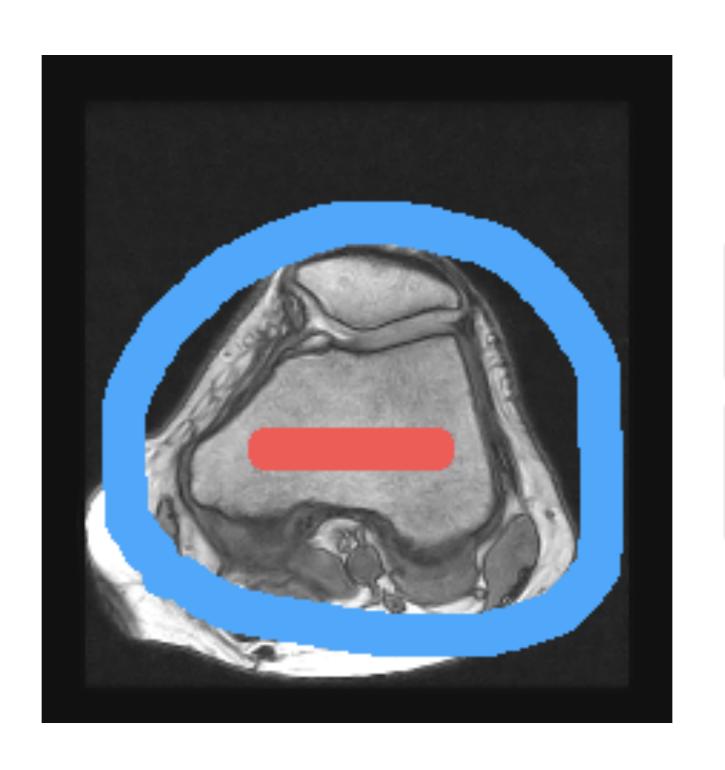
- Extension to the 3D case:
 - Instead of a curve we have a parametric surface; e.g., we can start using a sphere.
- Disadvantages:
 - We may have an over-smooth boundaries when using splines
 - How many *n* control points?
 - Not trivial to avoid self-intersection!

- Stroke-based 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.



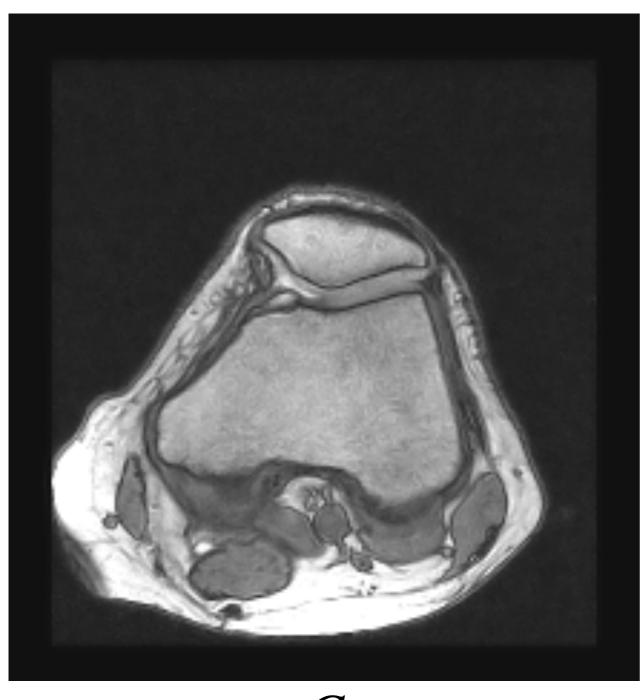
+1

-1



+1

_





(

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:

$$< l_i; \theta_i; C_i >$$

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

$$< l_i = 0; \theta_i = 0; C_i = I(x_i, y_i) > \forall_i s(x_i, y_i) = 0$$

Initialization for pixels covered by a stroke (s):

$$< l_i = s(x_i, y_i); \theta_i = 1; C_i = I(x_i, y_i) > \forall_i s(x_i, y_i) \neq 0$$

Stroke-Based: Grow-Cut

• For each pixel, we have:

Strength
Label
$$\langle l_i; \theta_i; C_i \rangle$$
 Intensity

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

$$< l_i = 0; \theta_i = 0; C_i = I(x_i, y_i) > \forall_i s(x_i, y_i) = 0$$

Initialization for pixels covered by a stroke (s):

$$< l_i = s(x_i, y_i); \theta_i = 1; C_i = I(x_i, y_i) > \forall_i s(x_i, y_i) \neq 0$$

Stroke-Based: A Single Grow-Cut Pass

- For each pixel I in the image:
 - We copy the previous status:

$$< l_i^{t+1}, \theta_i^{t+1}, I_i^{t+1} > = < l_i^t, \theta_i^t, I_i^t >$$

- For each neighbor j of i:
 - if $g(\|C_i^t C_j^t\|_2)\theta_j^t > \theta_i^t$ then

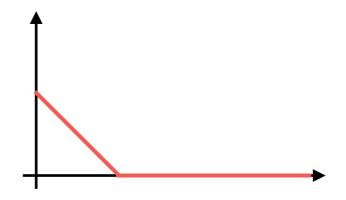
$$l_i^{t+1} = l_j^t$$

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

Stroke-Based: A Single Grow-Cut Pass

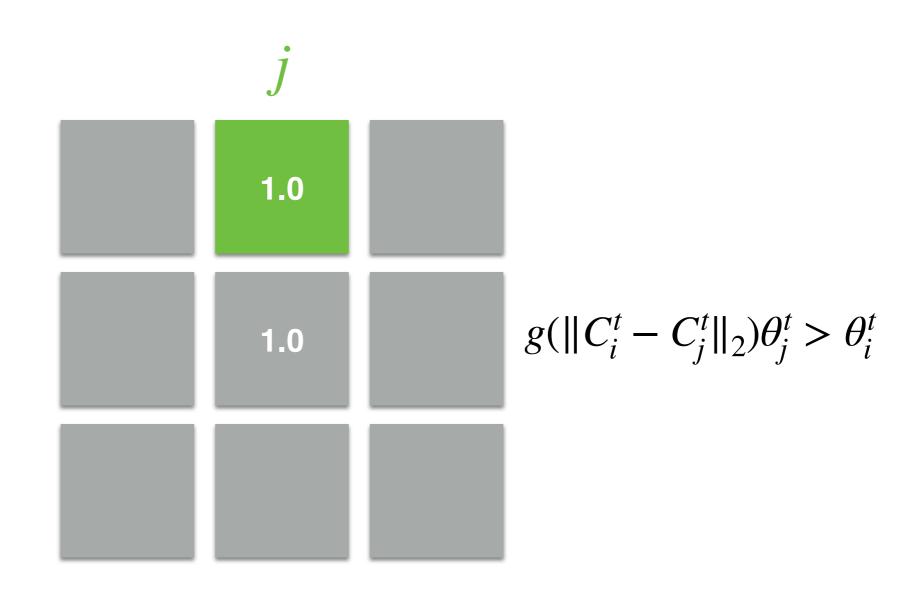
• Note that *g* is a decreasing function. For example:

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

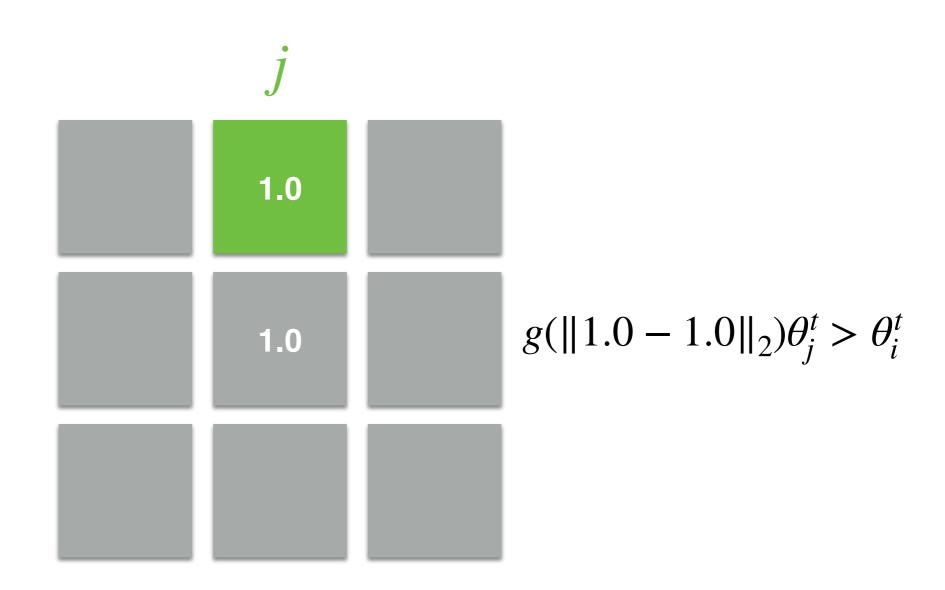


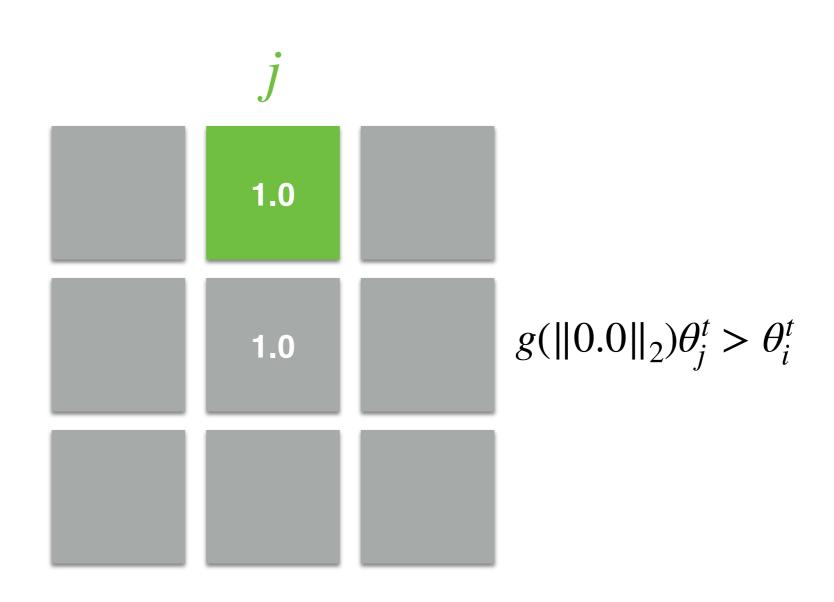
- This means that if the two pixels, which we compare, are close in intensity/color values they should have the same label *l*.
- They should also share the same label the neighbors have a higher strength!

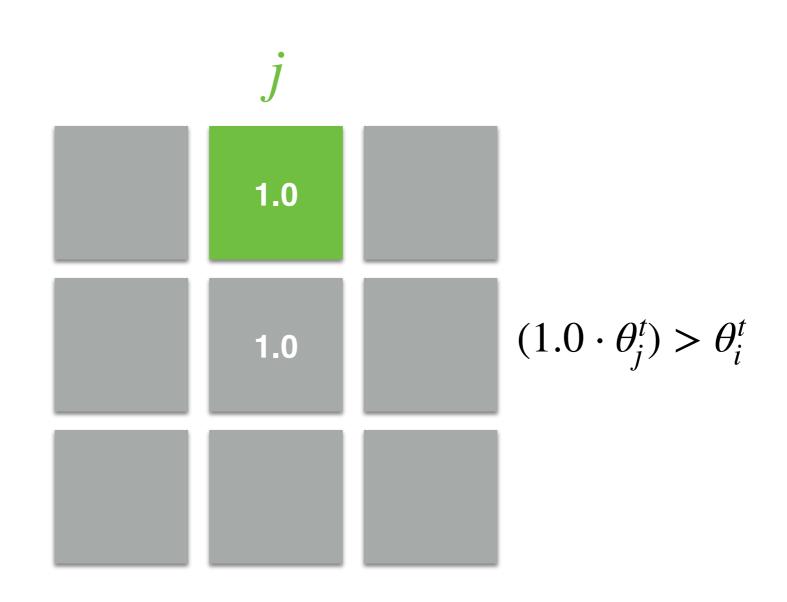
Stroke-Based: Example 1 - Switching Labels

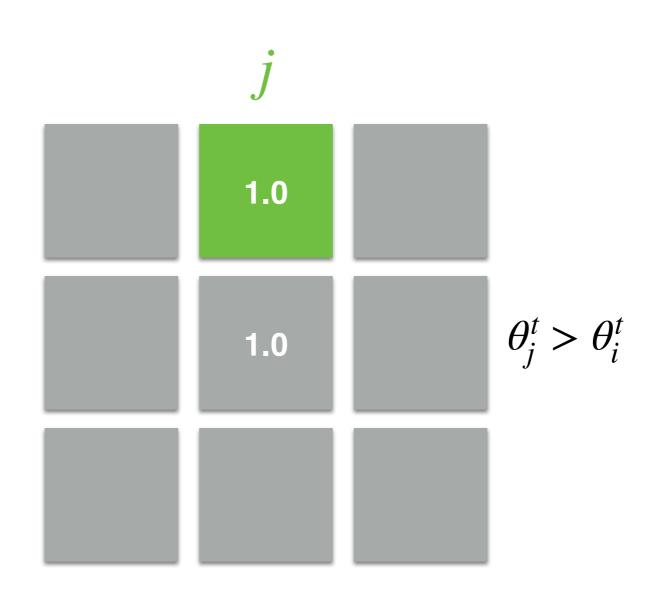


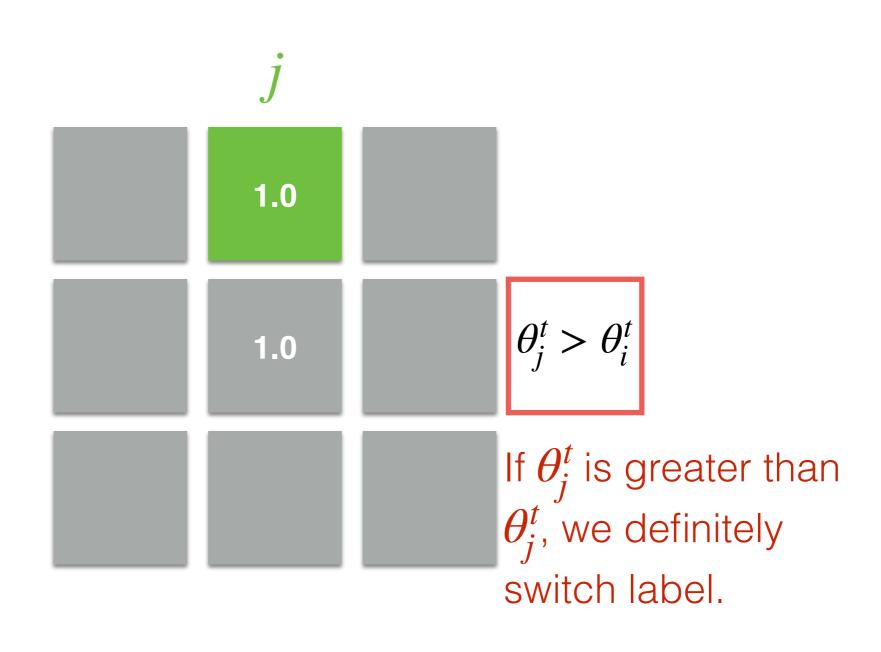
Stroke-Based: Example 1 - Switching Labels

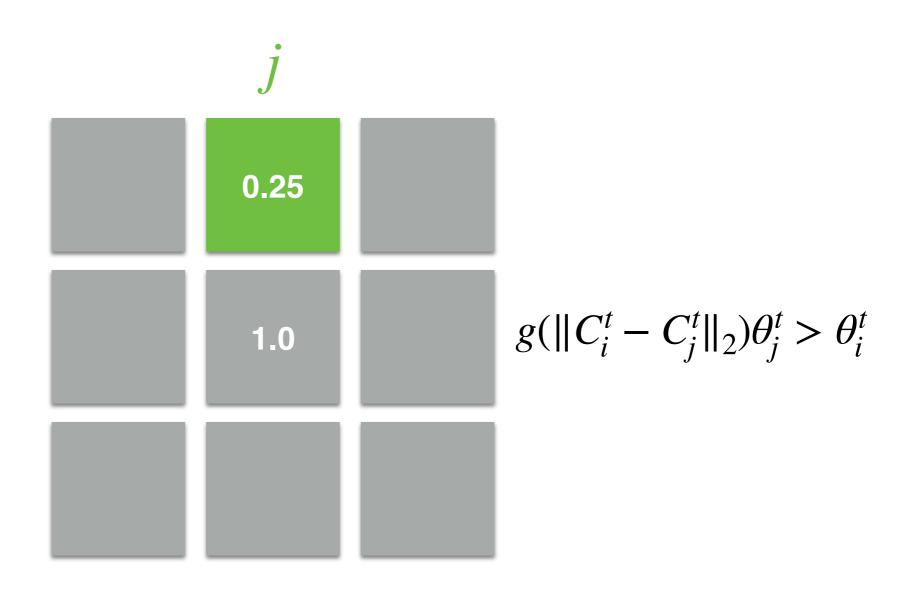


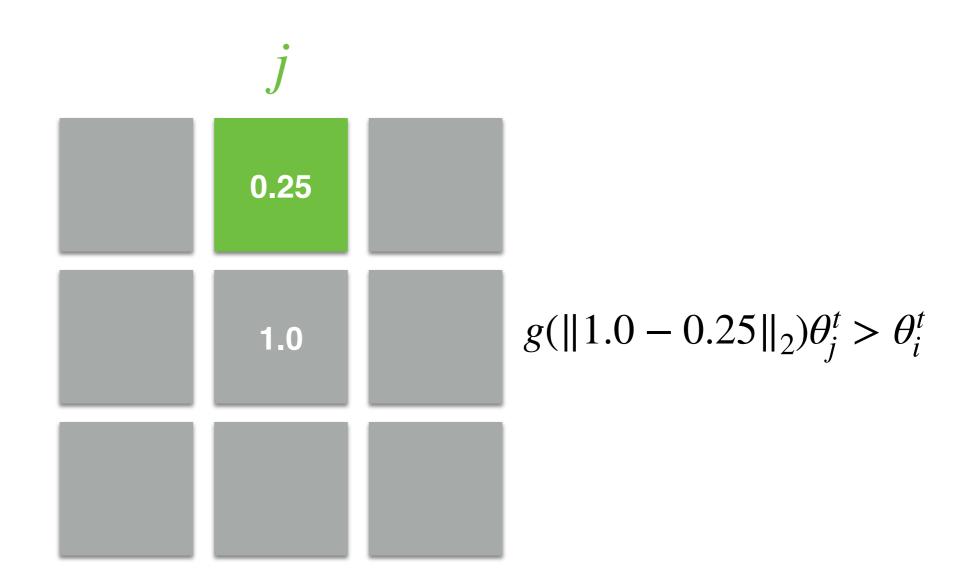


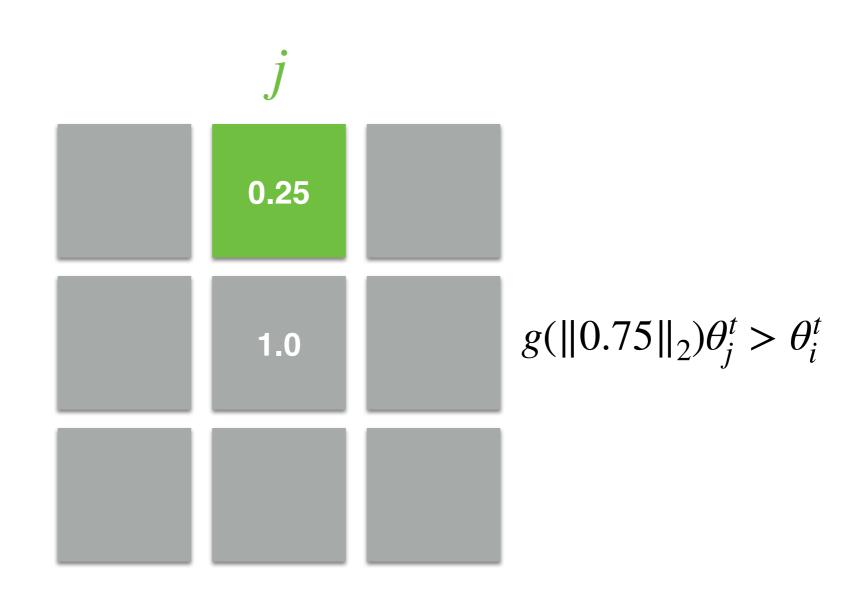


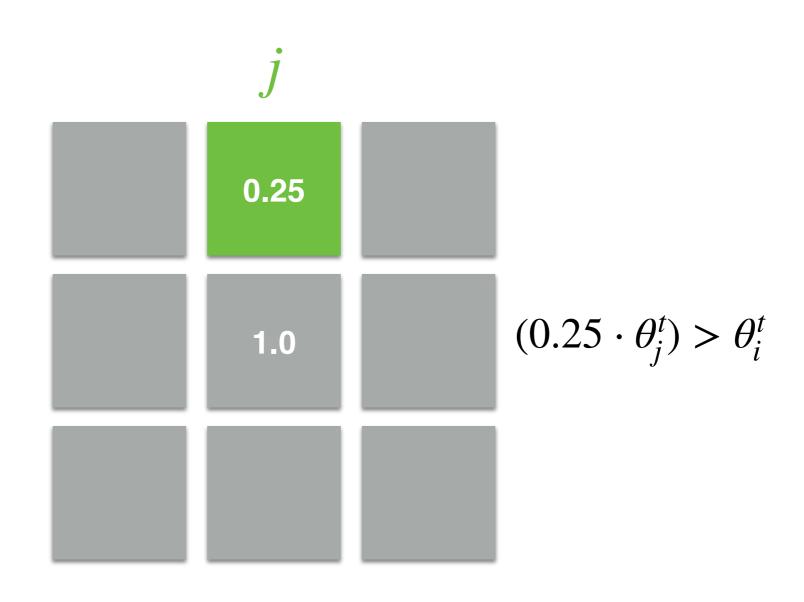


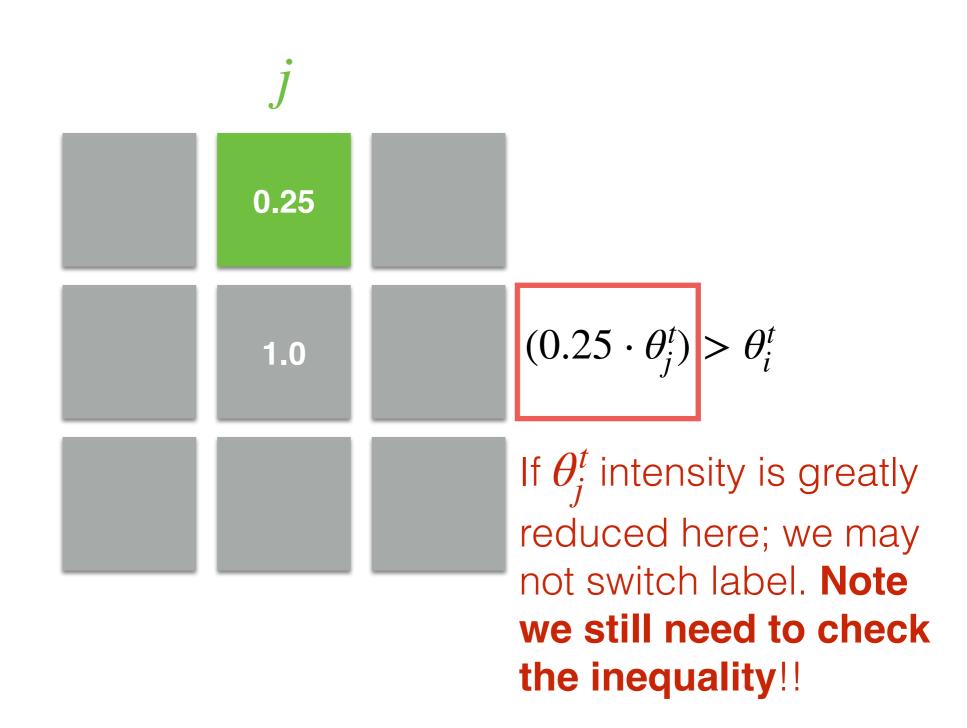








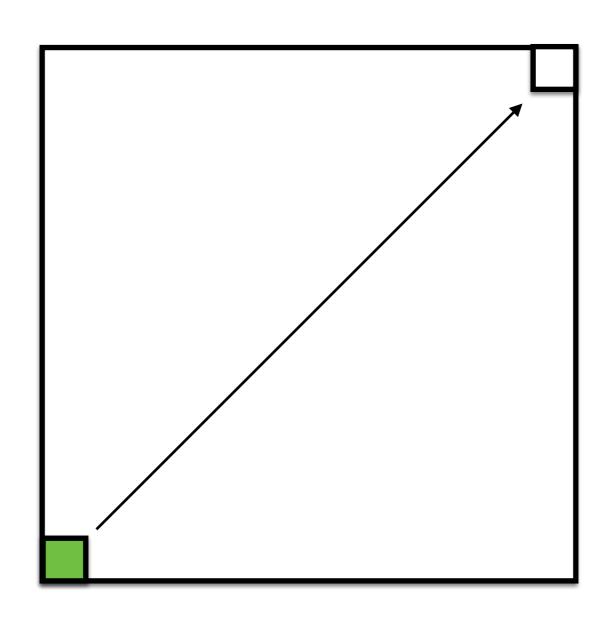


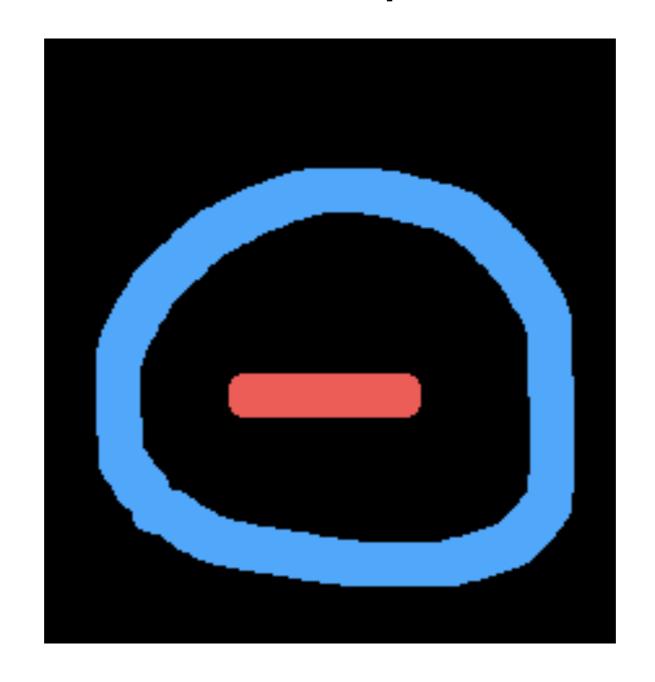


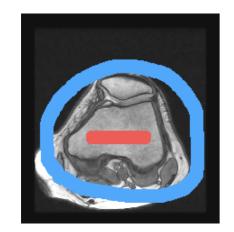
Stroke-Based: Grow-Cut

- 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

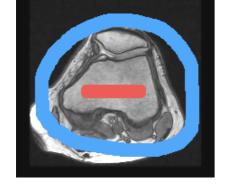




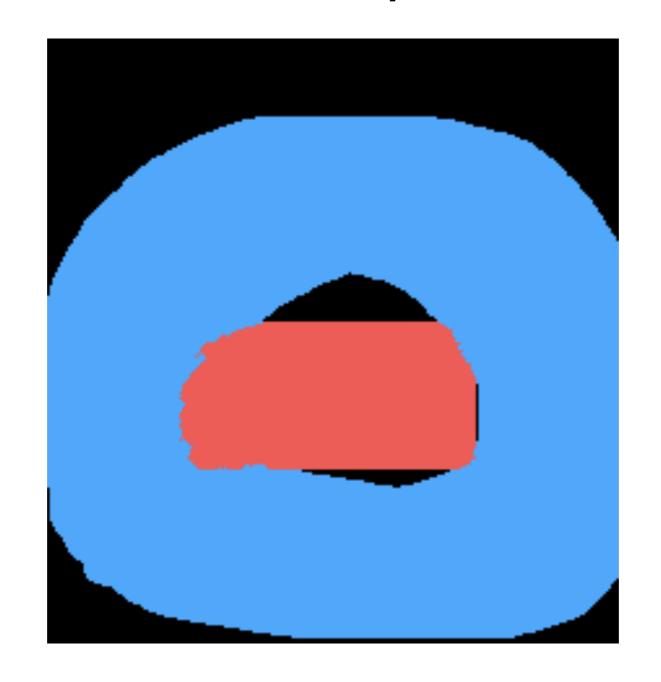


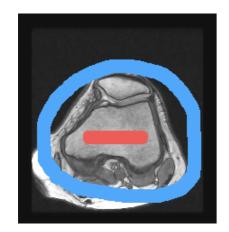
Iteration = 0



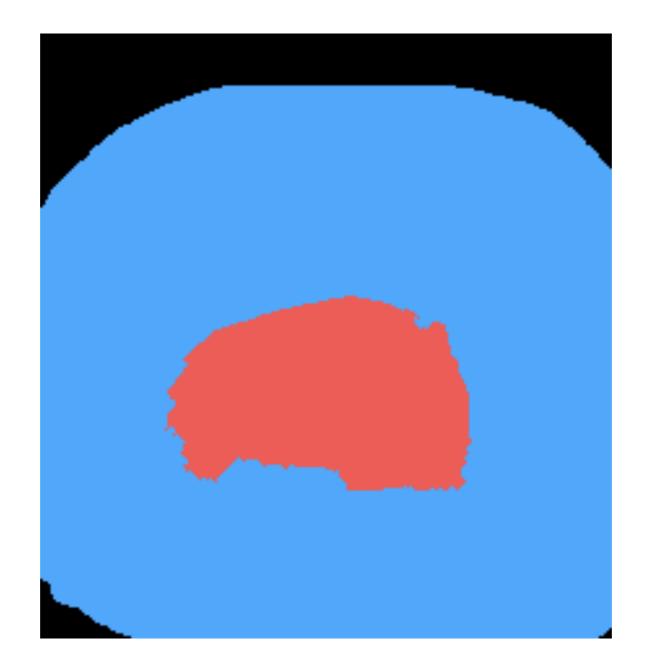


Iteration = 10

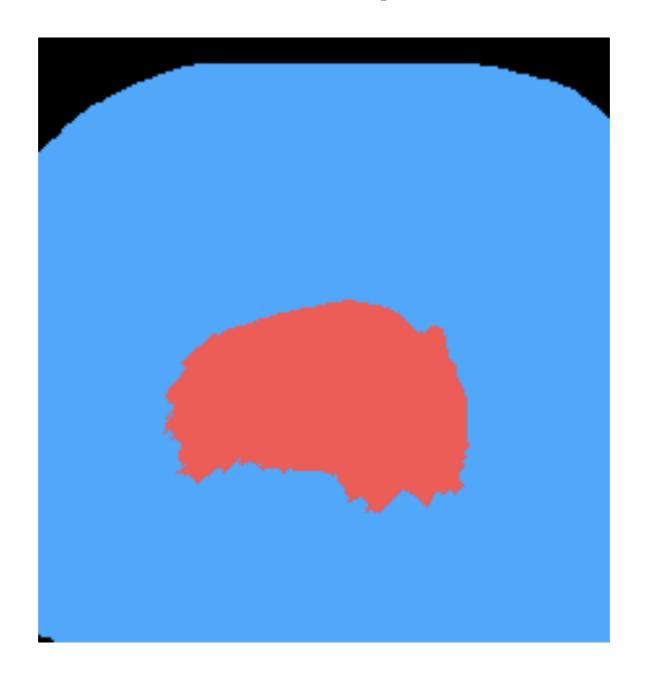


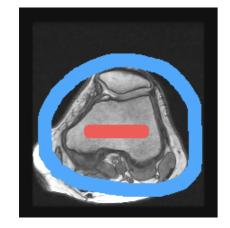


Iteration = 20

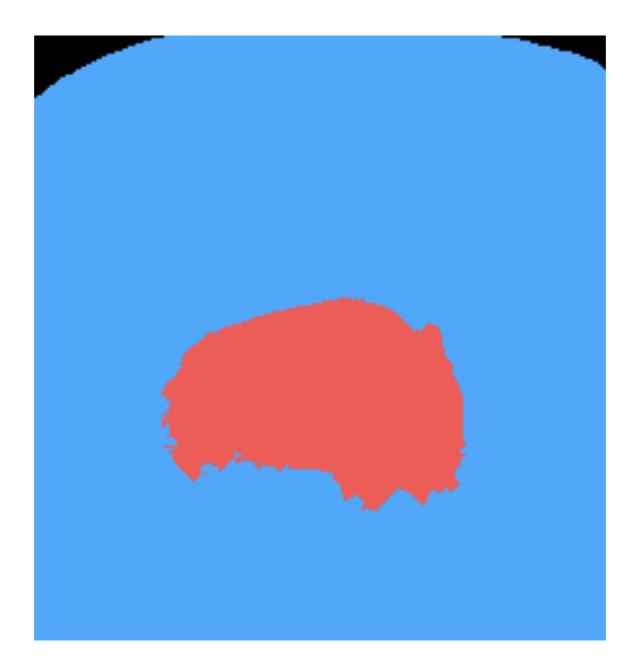


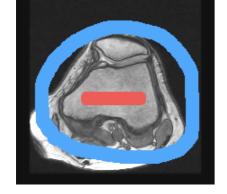






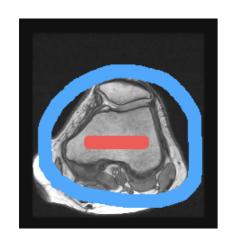
Iteration = 40



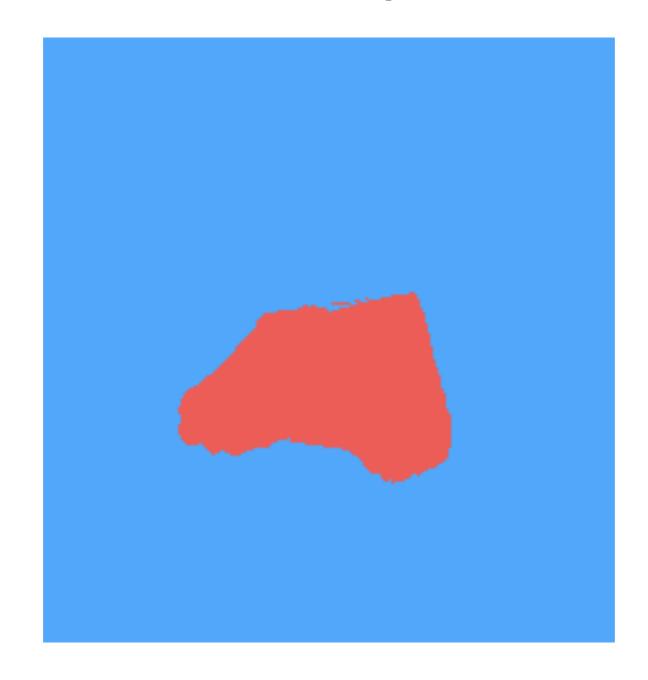


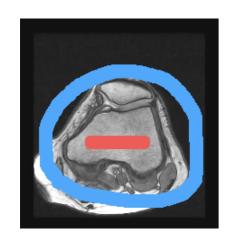
Iteration = 50



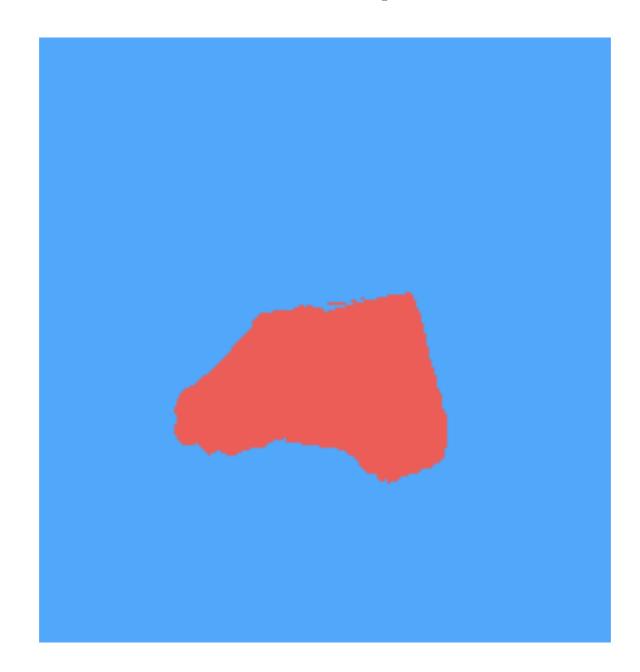


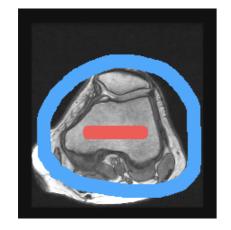
Iteration = 100





Iteration = 200





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!