

# Level-set MCMC Curve Sampling

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# Outline

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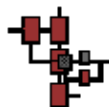
1. Overview
2. Curve evolution
3. Markov chain Monte Carlo
4. Curve sampling
5. Examples
6. Conditional simulation



# Overview

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- Curve evolution attempts to find a curve  $C$  (or curves  $C_i$ ) that best segment an image (according to some model)
- Goal is to minimize an energy functional  $E(C)$  (view as a negative log likelihood)
- Find a local minimum using gradient descent



# Sampling instead of optimization

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- Draw multiple samples from a probability distribution  $\mathbf{p}$  (e.g., uniform, Gaussian)
- Advantages:
  - Naturally handles multi-modal distributions
  - Can get out of local minima
  - Higher-order statistics (e.g., variances)
  - Conditional simulation



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# Planar curves

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- A curve is a function  $\vec{C} : [0, 1] \rightarrow \mathbb{R}^2$
- We wish to minimize an energy functional with a data fidelity term and regularization term:

$$E(\vec{C}) = D(y|\vec{C}) + \mathcal{R}(\vec{C})$$

- This results in a gradient flow:

$$\frac{d\vec{C}}{dt}(p) = \vec{F}(p)$$

- We can write any flow in terms of the normal:

$$\frac{d\vec{C}}{dt}(p) = f(p)\vec{N}(p)$$



# Level-Set Methods

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- A curve is a function (infinite dimensional)
- A natural implementation approach is to use marker points on the boundary (snakes)
  - Reinitialization issues
  - Difficulty handling topological change
- Level set methods instead evolve a surface (one dimension higher than our curve) whose zeroth level set is the curve (Sethian and Osher)



# Embedding the curve

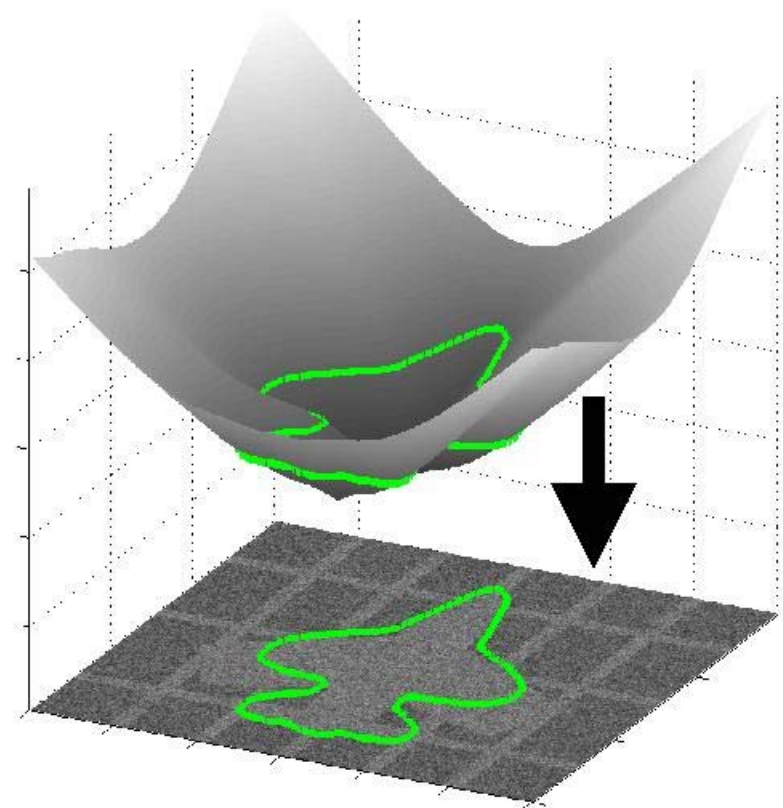
- Force level set  $\Psi$  to be zero on the curve

$$\Psi(\vec{C}(p)) = 0$$

$$\forall p \in [0, 1]$$

- Chain rule gives us

$$\begin{aligned} \frac{d\Psi}{dt} &= -\frac{d\vec{C}}{dt} \cdot \nabla\Psi \\ &= -f \|\nabla\Psi\| \end{aligned}$$







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# General MAP Model

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$$p(x|y; S) \propto p(y|x; S)p(x; S)$$

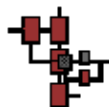
- For segmentation:
  - $x$  is a curve
  - $y$  is the observed image (can be vector)
  - $S$  is a shape model
  - Data model usually IID given the curve
- We wish to sample from  $p(x|y;S)$ , but cannot do so directly



# Markov Chain Monte Carlo

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- Class of sampling methods that iteratively generate candidates based on a previous iterate (forming a Markov chain)
- Instead of sampling from  $p(x|y;S)$ , sample from a proposal distribution  $q$  and keep samples according to an acceptance rule  $a$
- Examples include Gibbs sampling, Metropolis-Hastings



# Metropolis-Hastings

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- Metropolis-Hastings algorithm:

1. Start with  $x^0$
2. At time  $t$ , generate candidate  $\phi^t$  (given  $x^{t-1}$ )
3. Calculate Hastings ratio:

$$r^t = \frac{p(\phi^t)}{p(x^{t-1})} \cdot \frac{q(x^{t-1}|\phi^t)}{q(\phi^t|x^{t-1})}$$

4. Set  $x^t = \phi^t$  with probability  $\min(1, r^t)$ , otherwise  $x^t = x^{t-1}$
5. Go back to 2



# Asymptotic Convergence

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- We want to form a Markov chain such that its stationary distribution is  $p(x)$ :

$$p(x) = \int p(\phi) T(x|\phi) d\phi$$

- To guarantee asymptotic convergence, sufficient conditions are:

- 1) Ergodicity
- 2) Detailed balance

$$p(x^{t-1})q(\phi^t|x^{t-1})a(\phi^t|x^{t-1}) = p(\phi^t)q(x^{t-1}|\phi^t)a(x^{t-1}|\phi^t)$$



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# MCMC Curve Sampling

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- Generate perturbation on the curve:

$$\vec{C}'(s) = \vec{C}(s) + f(s)\vec{\mathcal{N}}(s)dt$$

- Sample by adding smooth random fields:

$$f \sim \mathbf{N}(0, \Sigma)$$

- $\Sigma$  controls the degree of smoothness in field
- Note for portions where  $f$  is negative, shocks may develop (so called prairie fire model)
- Implement using white noise and circular convolution



# Smoothness issues

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- While detailed balance assures asymptotic convergence, may need to wait a very long time
- In this case, smooth curves have non-zero probability under  $\mathbf{q}$ , but are very unlikely to occur
- Can view accumulation of perturbations as
$$F(s) = \sum_i f_i(s) = h \circledast \sum_i n_i$$
( $h$  is the smoothing kernel,  $n_i$  is white noise)
- Solution: make  $\mathbf{q}$  more likely to move towards high-probability regions of  $\mathbf{p}$





# Adding mean force

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- We can add deterministic elements to  $f$  (i.e., a mean to  $\mathbf{q}$ ):

$$f \sim \mathcal{N}(-\kappa + \gamma, \Sigma)$$

$$f(s) = \beta r(s) - \alpha \kappa(s) + \gamma$$

- The average behavior should then be to move towards higher-probability areas of  $\mathbf{p}$
- In the limit, setting  $f$  to be the gradient flow of the energy functional results in always accepting the perturbation



# Coverage/Detailed balance

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- It is easy to show we can go from any curve  $C_1$  to any other curve  $C_2$  (shrink to a point)
- For detailed balance, we need to compute probability of generating  $C'$  from  $C$  (and vice versa)

$$\vec{C}'(s) = \vec{C}(s) + f(s)\vec{\mathcal{N}}(s)dt$$

$$\vec{C}(s) = \vec{C}'(s) + f'(s)\vec{\mathcal{N}}'(s)dt$$

- Probability of going from  $C$  to  $C'$  is the probability of generating  $f$  (which is Gaussian) and the reverse is the probability of  $f'$  (also Gaussian)



# Approximations to $q$

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- Relationship between  $f$  and  $f'$  complicated due to the fact that the normal function changes
  - $f'$  does not always exist (given an  $f$ ). Unknown what conditions on  $f$  are necessary to guarantee existence.
  - Various levels of exactness
    - Assume  $\vec{\mathcal{N}} = \vec{\mathcal{N}}'$  (then  $f' = -f$ )
    - Infinitesimal approximation (ignore tangential)
$$f'(s) = \langle -f(s)\vec{\mathcal{N}}(s), \vec{\mathcal{N}}'(s) \rangle$$
    - Trace along  $\vec{\mathcal{N}}$  (technical issues)
  - Unknown how approximations affect convergence
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# SAD Target

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- We define symmetric area difference (SAD) as:

$$d_{\text{SAD}}(\Psi_1, \Psi_2) = \int_{\Omega} (\mathcal{H}(-\Psi_1(x)) - \mathcal{H}(-\Psi_2(x)))^2 dx$$

- Use a Boltzmann distribution:

$$p(\vec{C}|\vec{C}_0) = \frac{1}{Z} \exp(-d_{\text{SAD}}(\vec{C}, \vec{C}_0)/T) p(\vec{C})$$

- T is a parameter we can use to control how likely we are to keep less likely samples
- We will keep a sample with T log(2) additional errors with probability 1/2
- Single mode distribution



# Target Shape

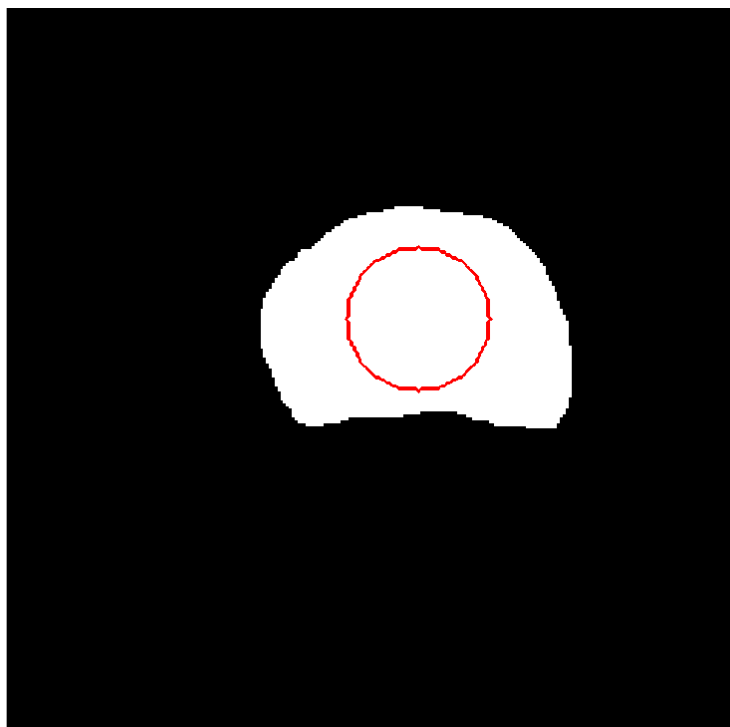
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# Initialization

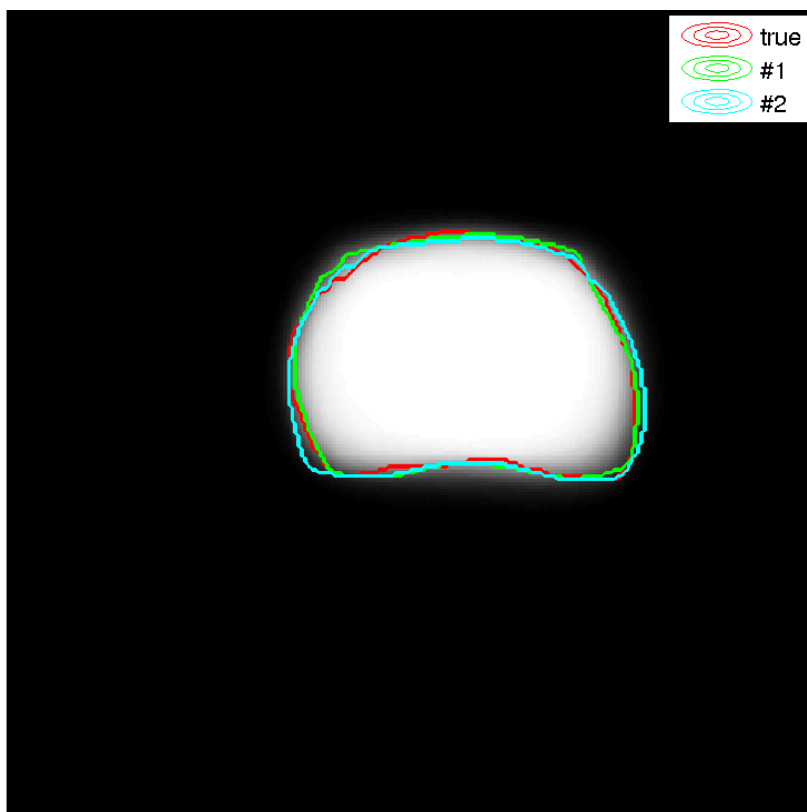
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# Most likely samples

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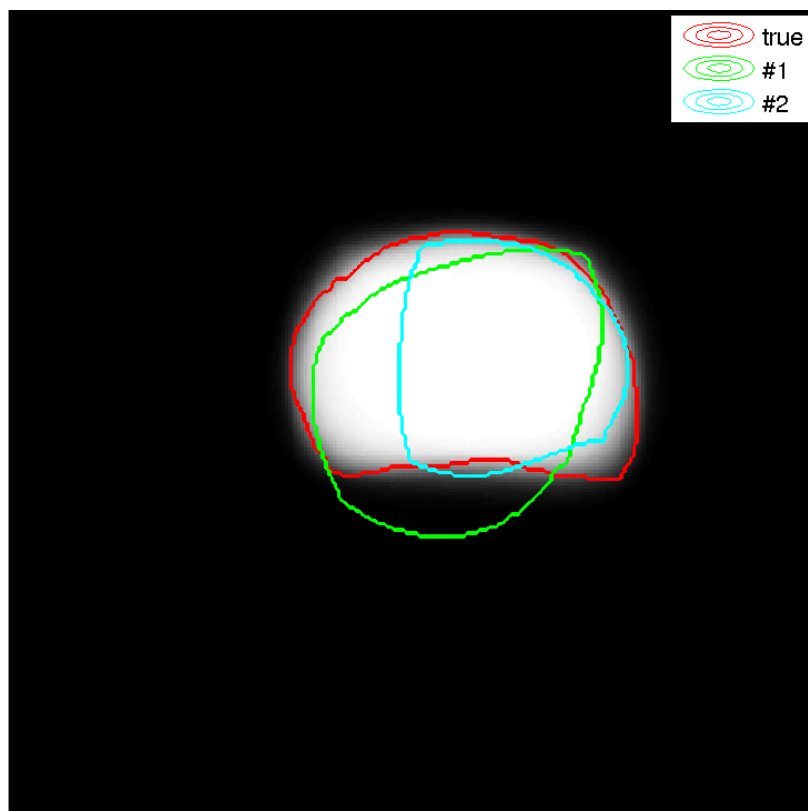






# Least likely samples

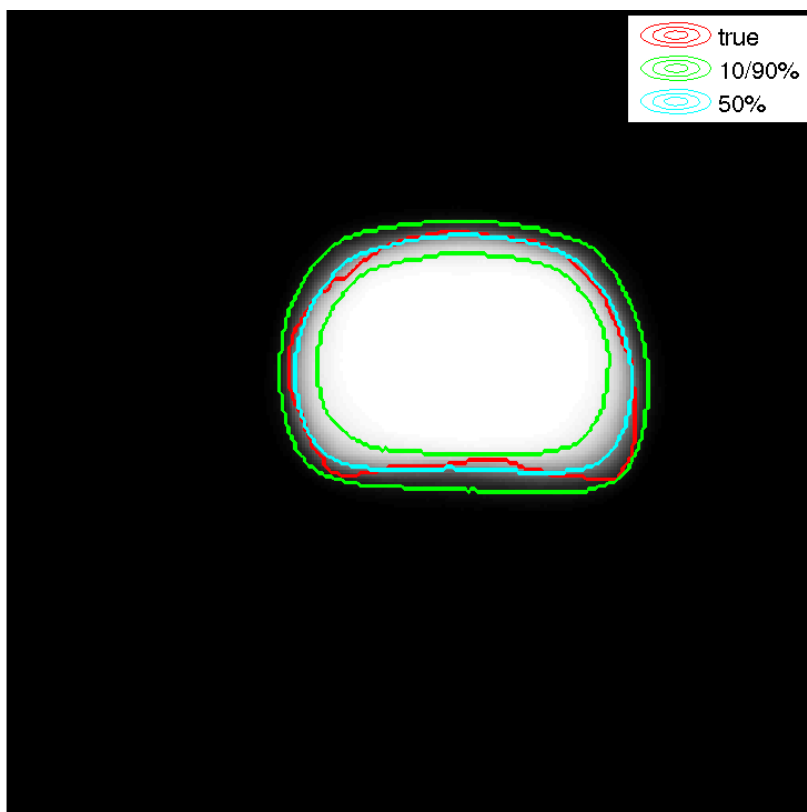
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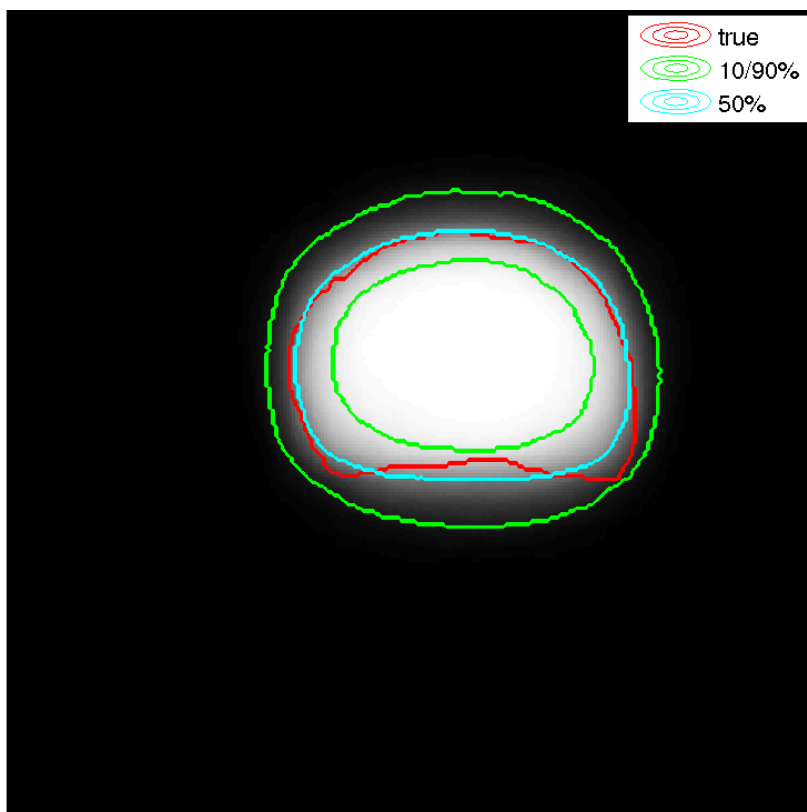
# “Confidence intervals”

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# Doubling the temperature





# Synthetic noisy image

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- Piecewise-constant observation model:

$$y(x) = \mu(x) + n(x)$$

- Chan-Vese energy functional:

$$E(\vec{C}) = \iint_{R_0} (y - \mu_0)^2 dx + \iint_{R_1} (y - \mu_1)^2 dx + \alpha \oint_{\vec{C}} ds$$

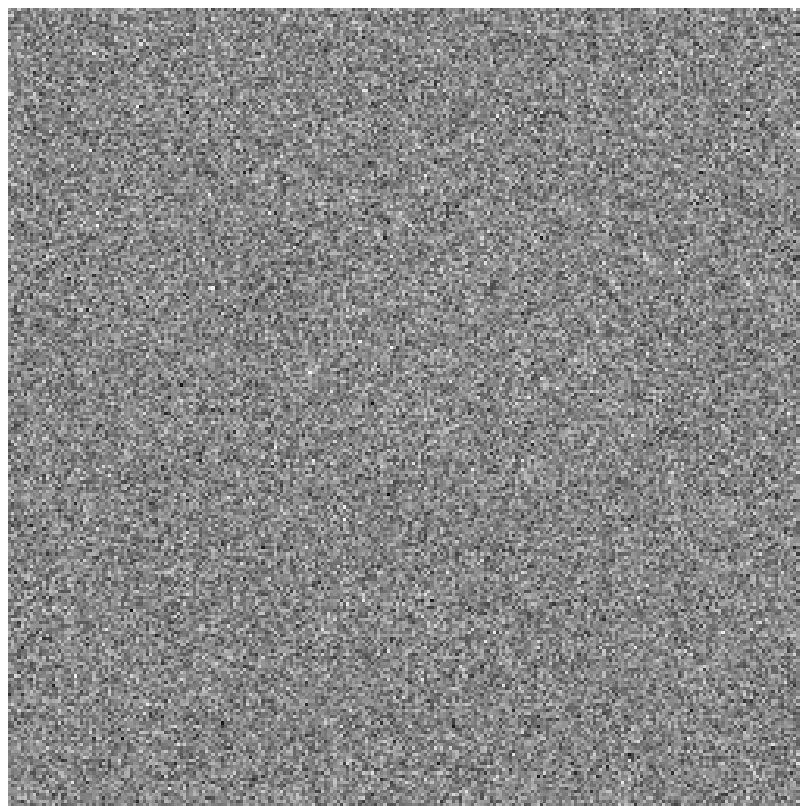
- Probability distribution ( $T=2\sigma^2$ ):

$$p(\vec{C}) = \frac{1}{Z} \exp(-E(\vec{C})/T)$$



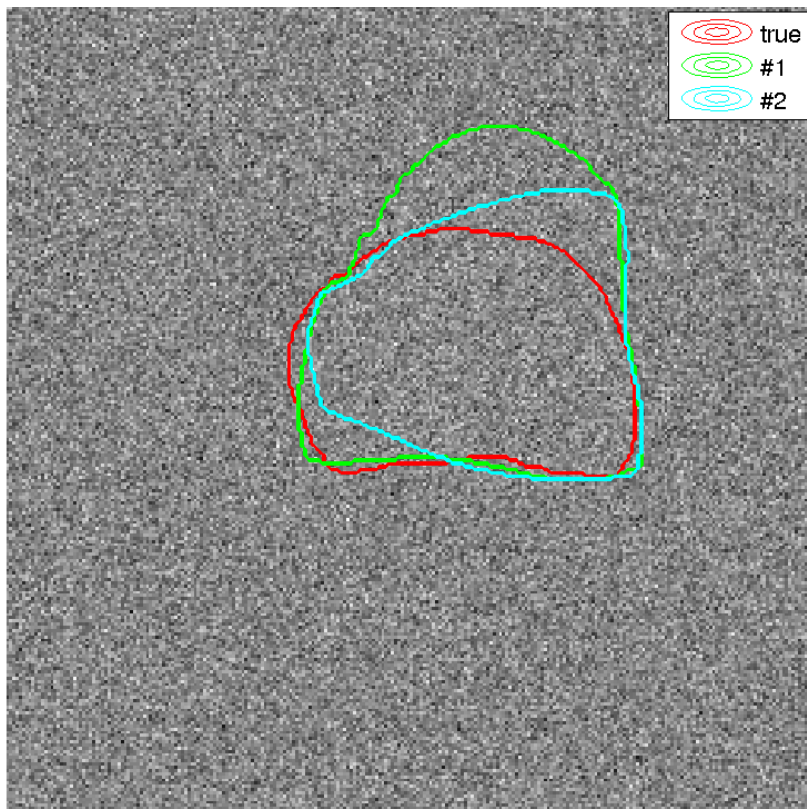
# Prostate in a Haystack

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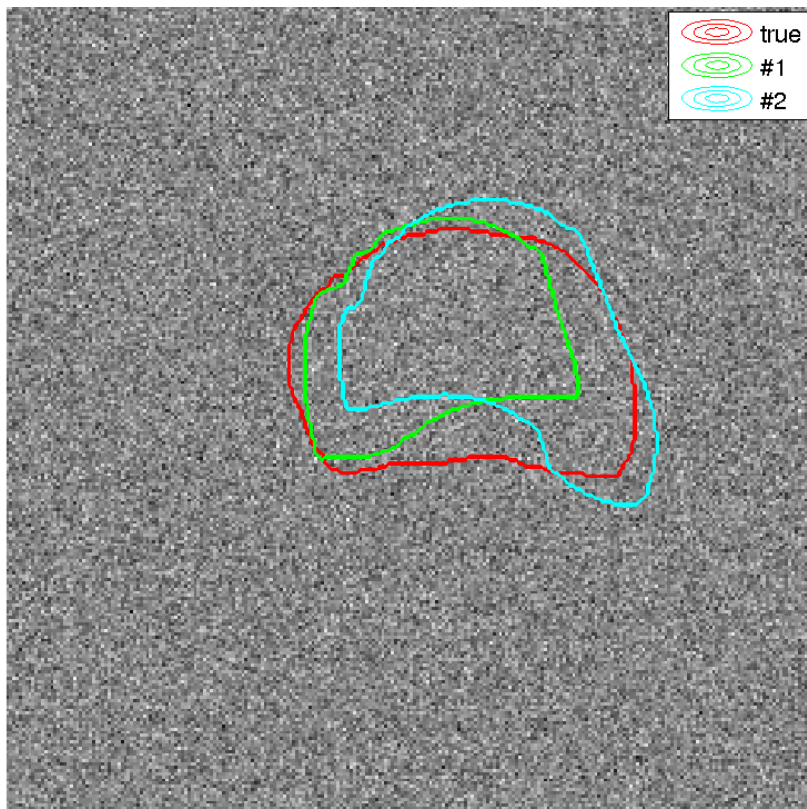


# Most likely samples



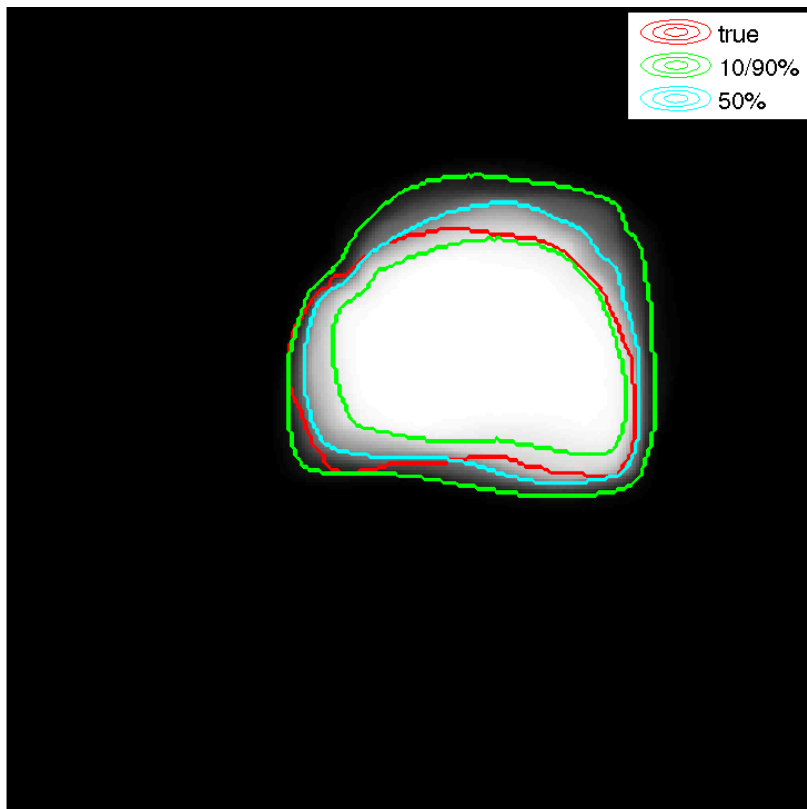


# Least likely samples





# Confidence intervals







## When “best” is not best

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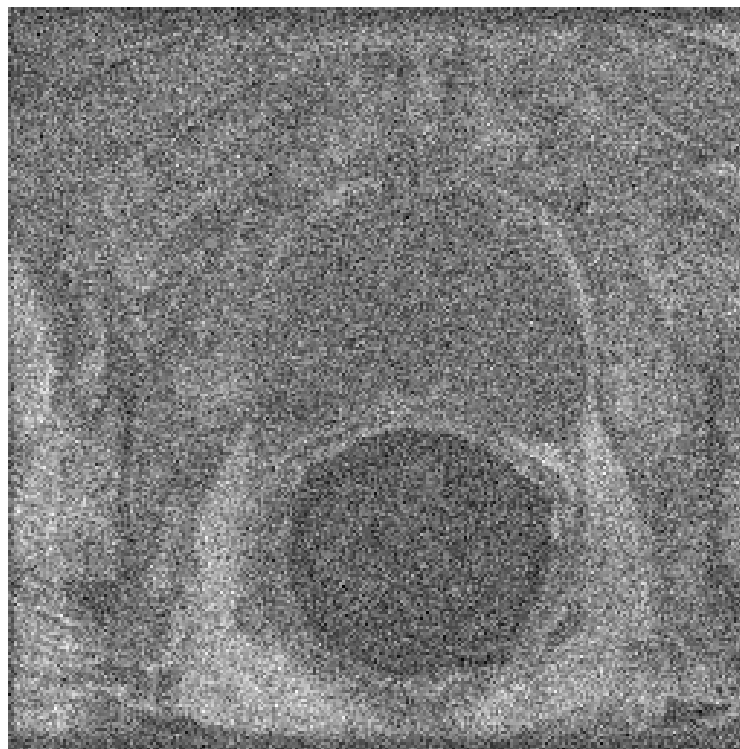
- In this example, the most likely samples under the model are not the most accurate according to the underlying truth
- 10%/90% “confidence” bands do a good job of enclosing the true answer
- Histogram image tells us more uncertainty in upper-right corner
- “Median” curve is quite close to the true curve
- Optimization would result in subpar results



# Bias-corrected prostate

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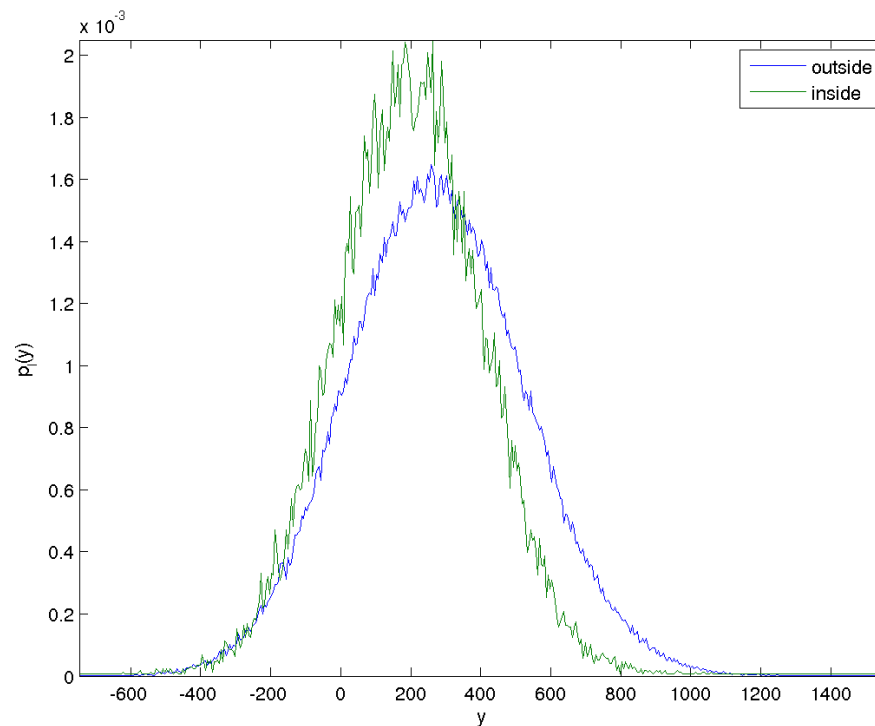
- “Expert” segmentation, add noise (simulate body coil image)





# Learn probability densities

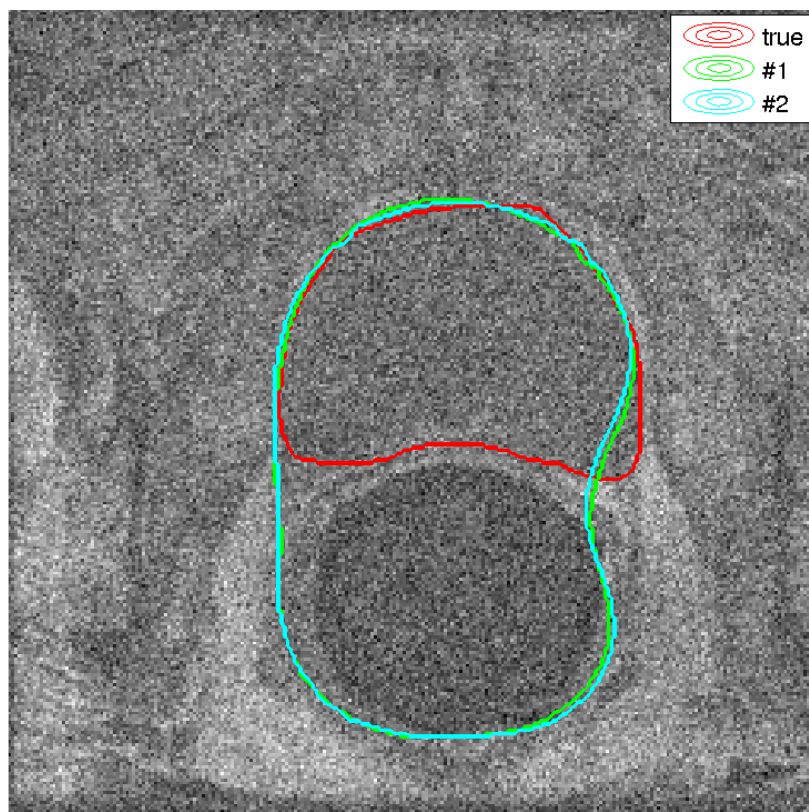
- Use histograms
- Learn pdf inside  $p(y|1)$  and pdf outside  $p(y|0)$  and assume iid given curve:



$$E(\vec{C}) = - \int_{\Omega} \log p(y(x) | \mathcal{H}(-\Psi(x))) dx + \alpha \oint_{\vec{C}} ds$$

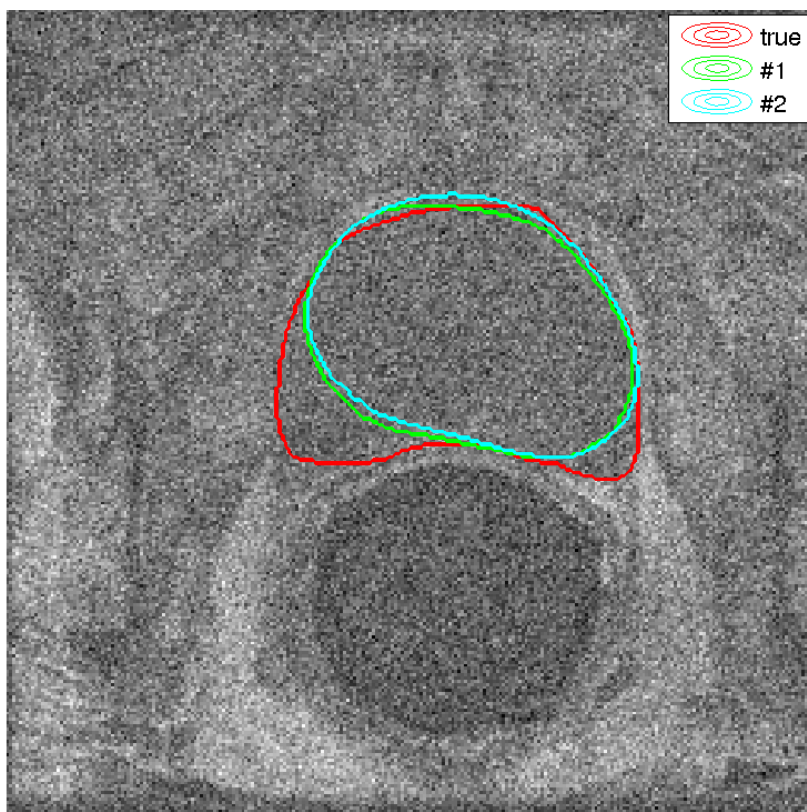


# Most likely samples





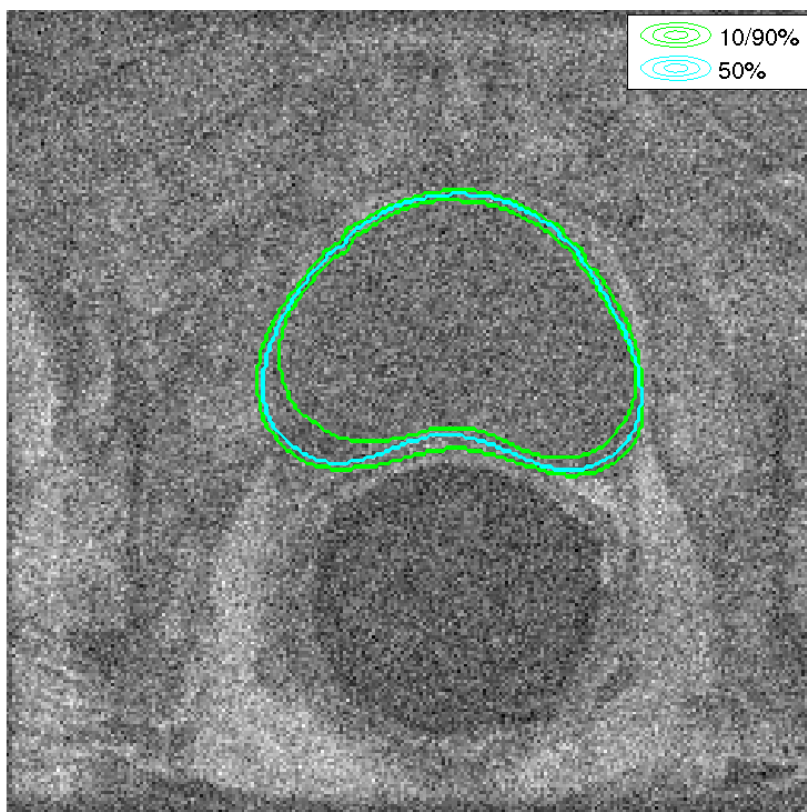
# Least likely samples





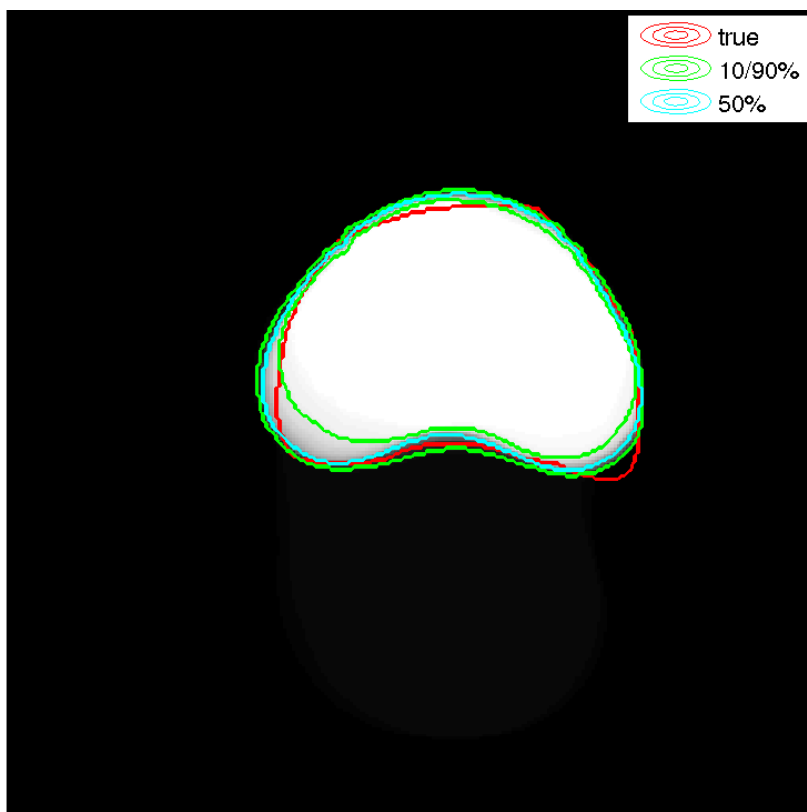
# Confidence intervals

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# Confidence intervals (with histogram)





# Multimodality and convergence

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- Natural multimodal distribution
- Burn-in time is not long enough (otherwise more samples would have clustered near the more-likely mode)
- When starting near one mode, need a lot of time to traverse valley between modes
- Clustering could help with presenting results





# Gravity inversion

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- Supplement to standard seismic data to segment bottom salt using an array of surface gravimeters ( $\sim 10^{-15}$  N accuracy)
- Subtract base effects (geoid, centrifugal force, etc.) to leave salt effects:

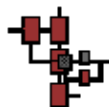
$$\vec{g}(i) = G \int_{\Omega} \frac{\rho(x; \vec{C}) \hat{r}_i(x)}{\|\vec{r}_i(x)\|^2} dx$$

- Assume constant density inside and outside:

$$\rho(x; \vec{C}) = \Delta\rho \mathcal{H}(-\Psi(x))$$

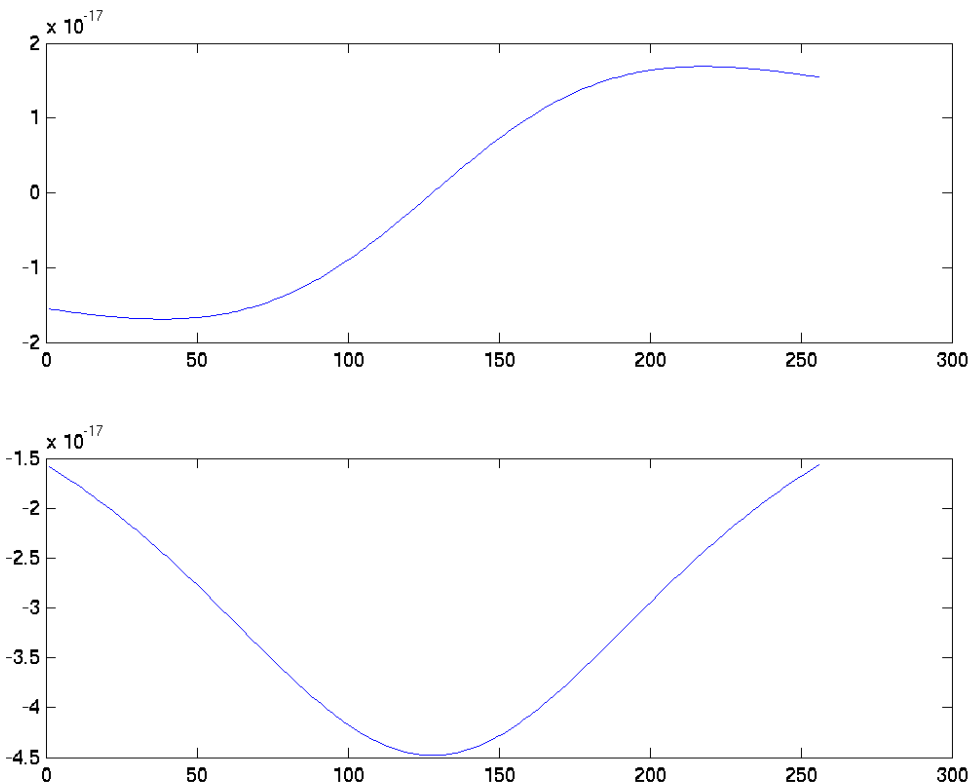
- Model energy as L2 estimation error (probability as Boltzmann distribution):

$$E(\vec{C}) = \sum_{i=1}^{N_{\text{array}}} \|\vec{g}_{\text{obs}}(i) - \vec{g}(i; \vec{C})\|^2 + \alpha \int_{\vec{C}} ds$$



# A strange segmentation problem

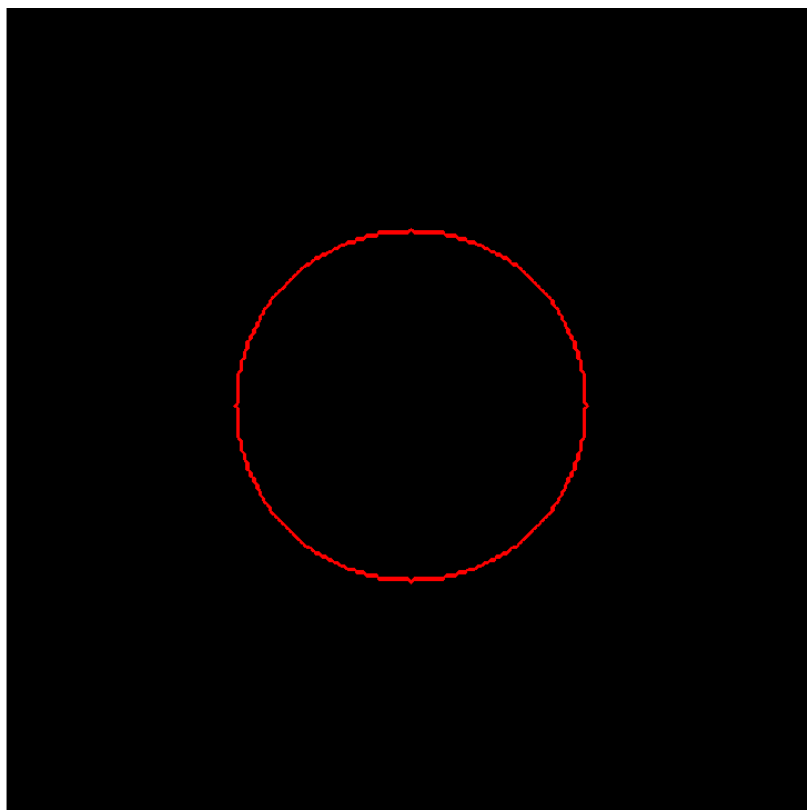
X- and z-  
components of  
gravity profile  
for synthetic salt  
body





# Circle salt

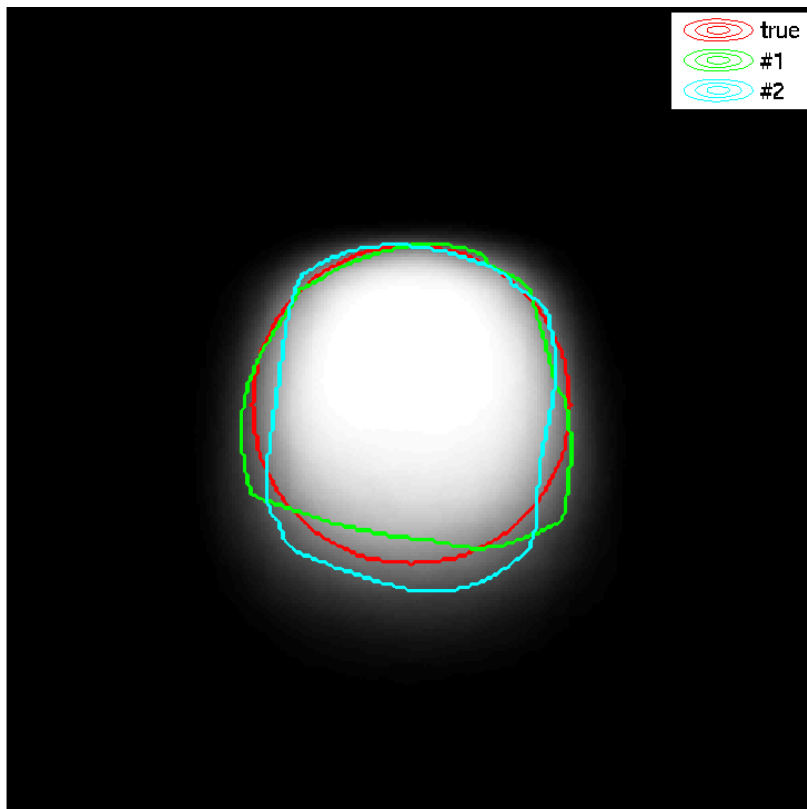
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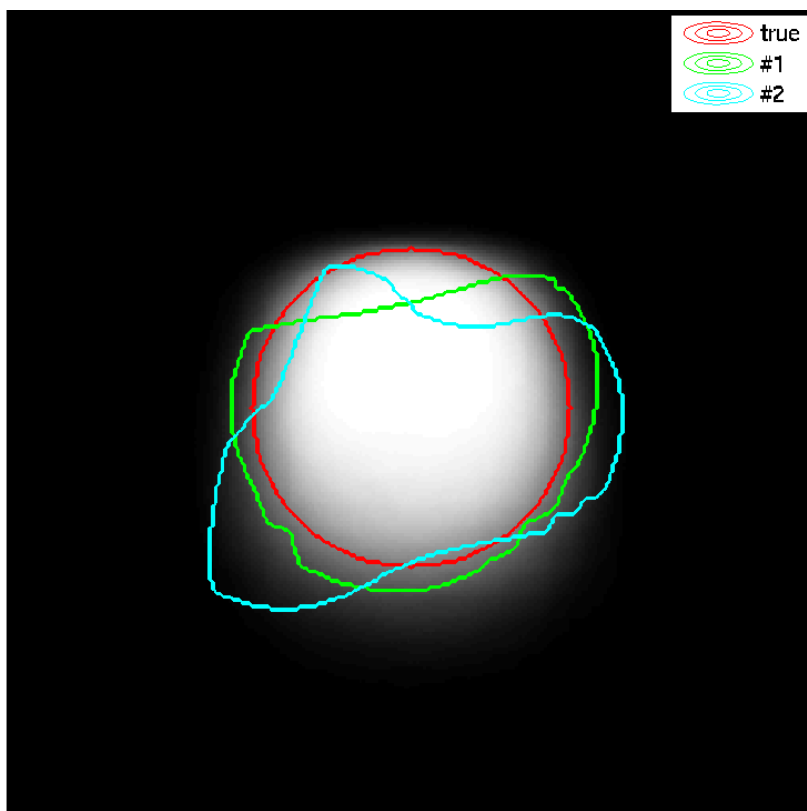
# Most likely samples

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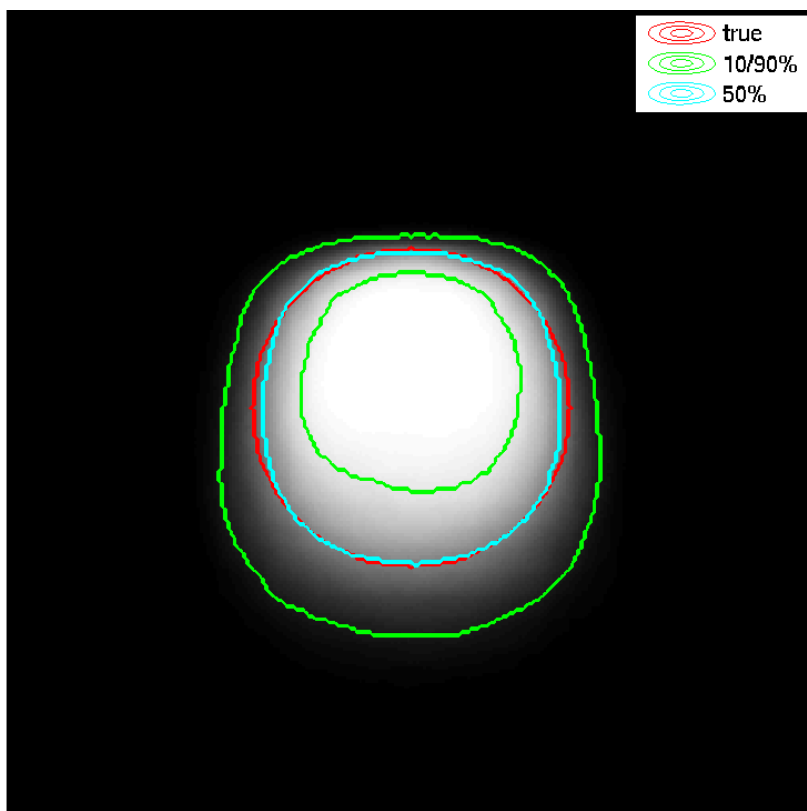
# Least likely samples





# Confidence intervals

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# Notable features

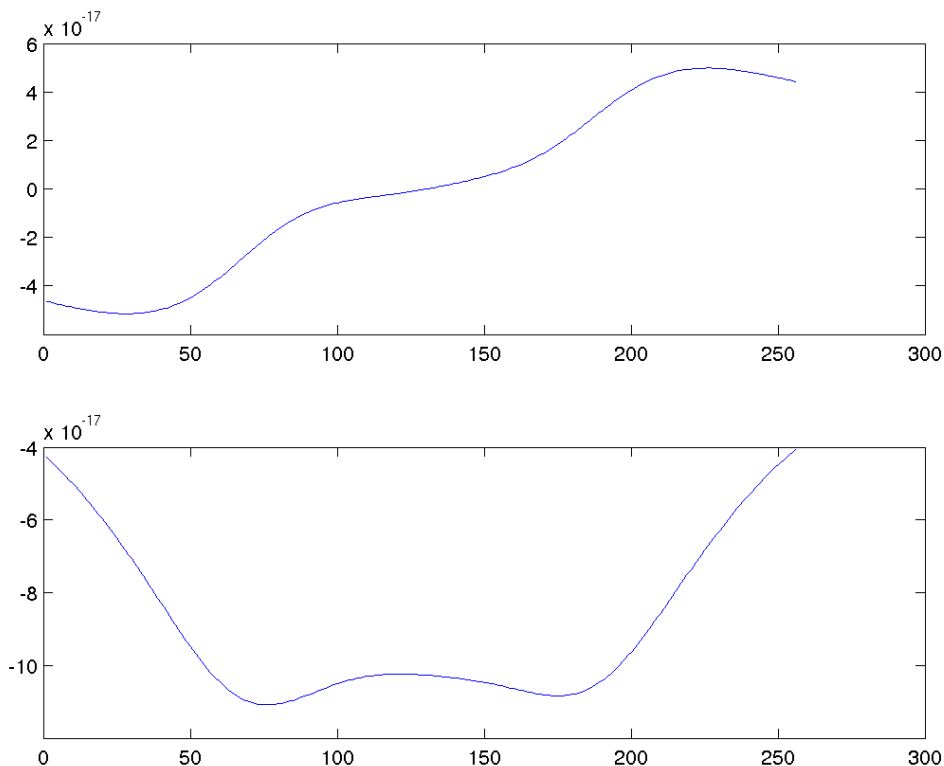
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- Measurement points  $\ll$  image pixels, but we can do a reasonable job
- Much higher uncertainty at the bottom than the top (weaker measurements)
- Less uncertainty in middle than on sides
- Median of histogram not necessarily related to median of distribution

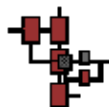


# Mystery example

- Same x- and z-components of gravity
- Synthetic image with unknown truth

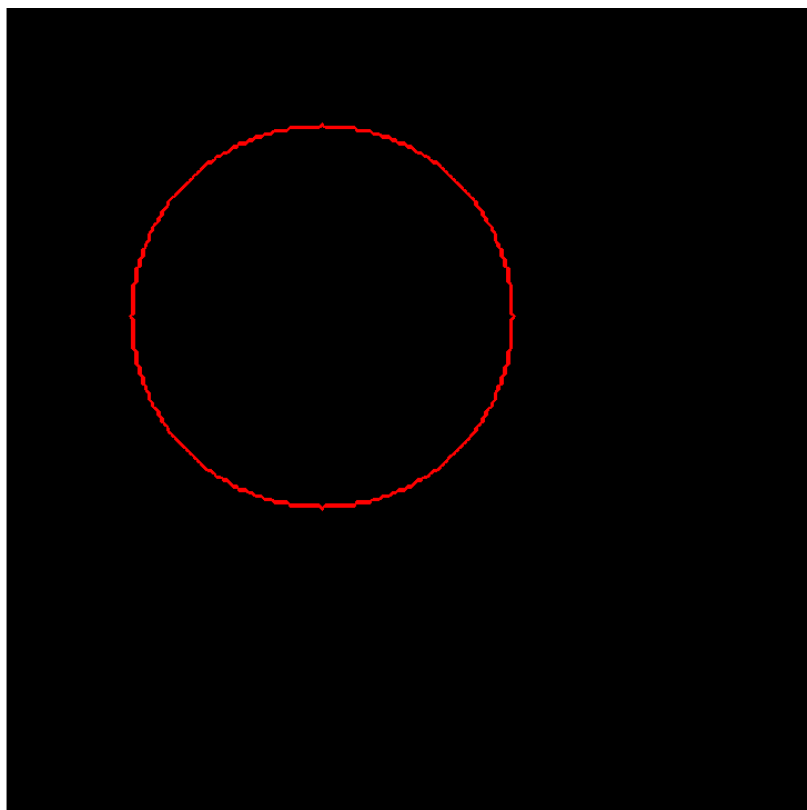






# Initial state

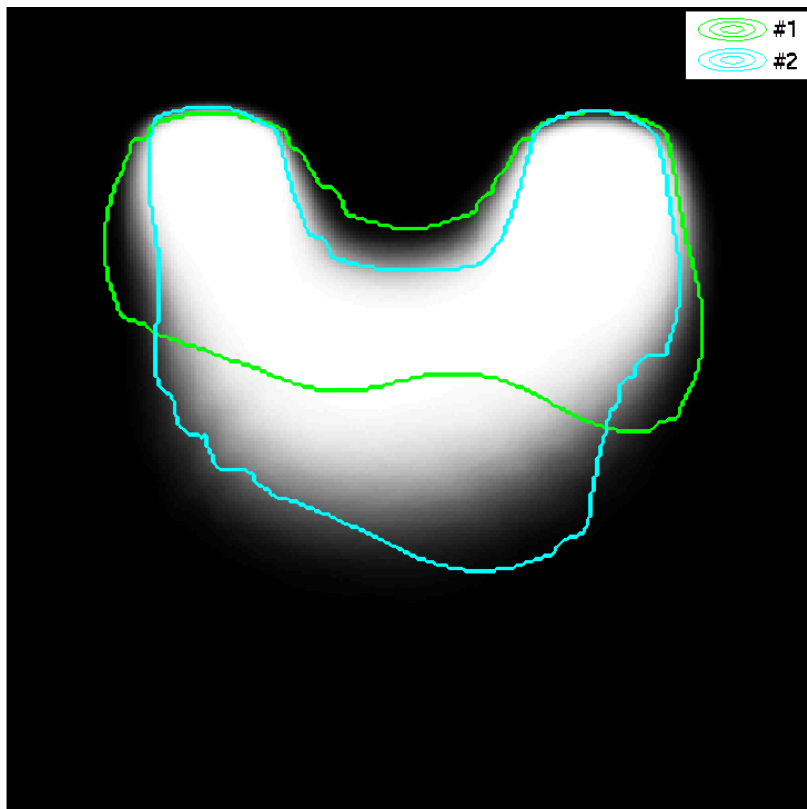
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# Most likely samples

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# Least likely samples

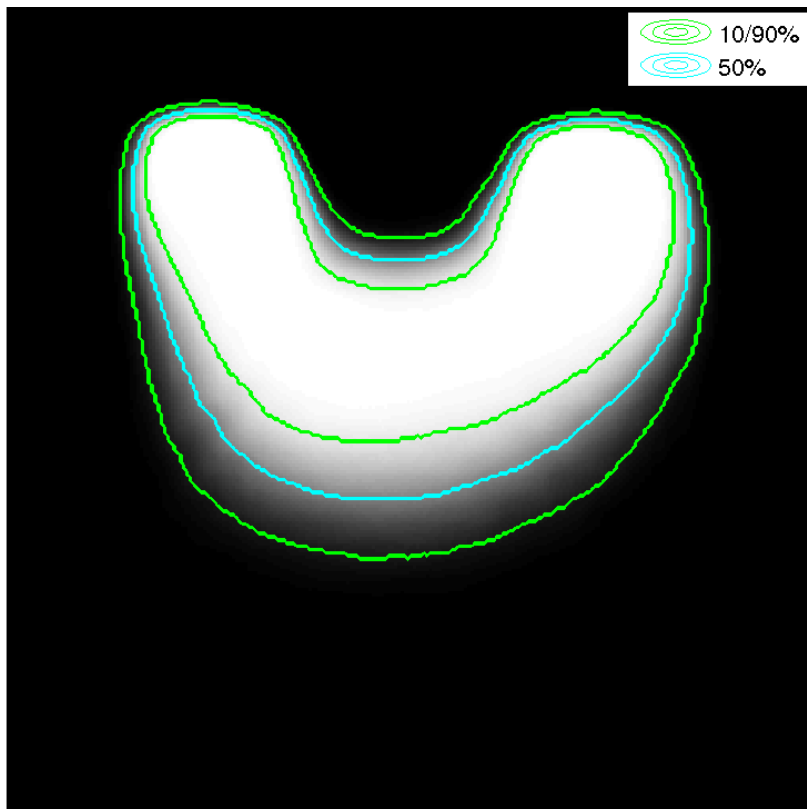
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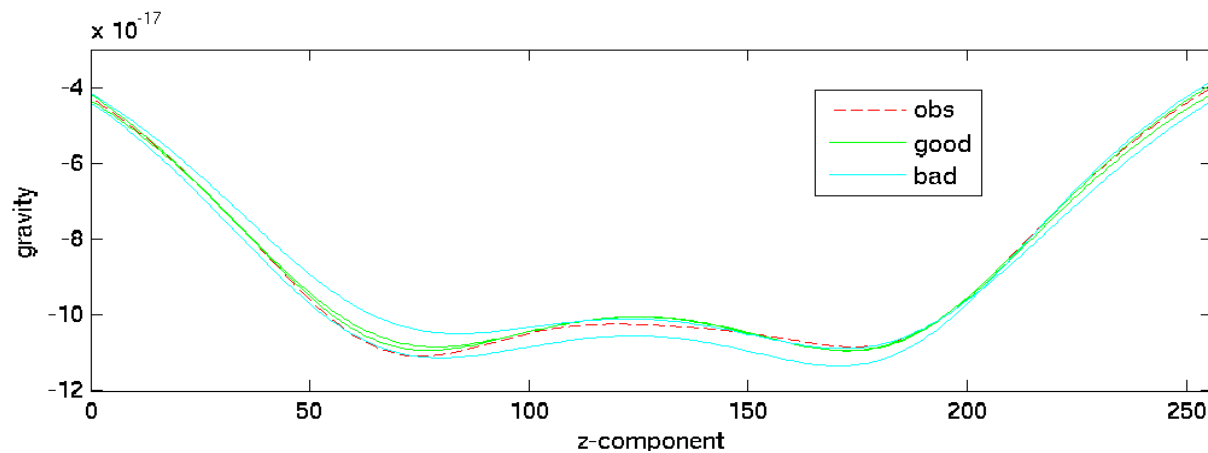
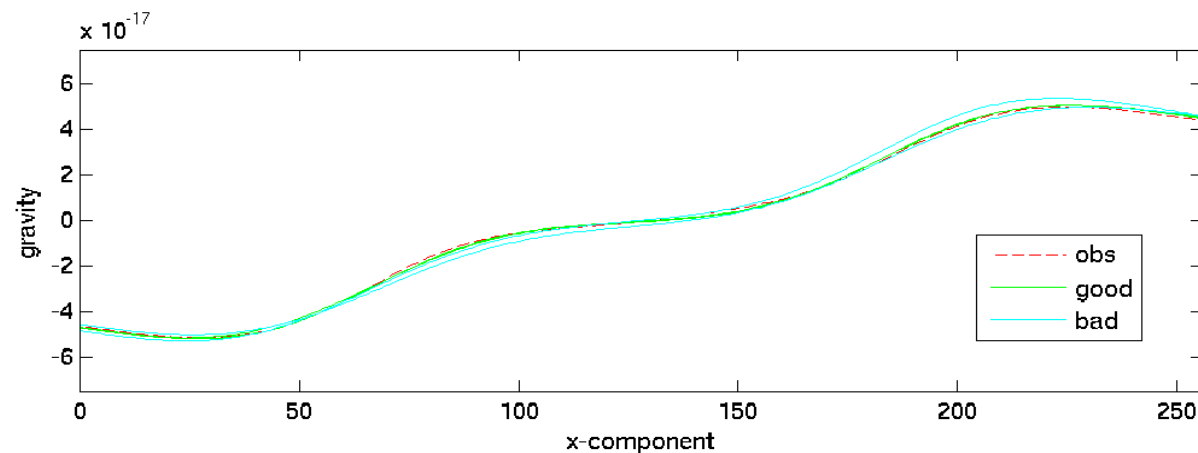
# Confidence intervals

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# Gravity error





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# User Information

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- In many problems, the model admits many reasonable solutions
- Currently user input largely limited to initialization
- We can use user information to reduce the number of reasonable solutions
  - Regions of inclusion or exclusion
  - Partial segmentations
- Can help with both convergence speed and accuracy
- Interactive segmentation



# Conditional simulation

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- With conditional simulation, we are given the values on a subset of the variables
- We then wish to generate sample paths that fill in the remainder of the variables (e.g., simulating Brownian motion)





# Simulating curves

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- Say we are given  $C_s$ , a subset of  $C$  (with some uncertainty associated with it)
- We wish to sample the unknown part of the curve  $C_u$

- One way to view is as sampling from:

$$p(y|\vec{C})p(\vec{C}) = p(y|\vec{C})p(\vec{C}_u|\vec{C}_s)p(\vec{C}_s)$$

- Difficulty is being able to evaluate the middle term as theoretically need to integrate  $p(C)$



# Simplifying Cases

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Under special cases, evaluation of  $p(\vec{C}_u | \vec{C}_s)$  is tractable:

1. When  $C$  is low-dimensional (can do analytical integration or Monte-Carlo integration)
2. When  $C_s$  is assumed to be exact
3. When  $p(C)$  has special form (e.g., independent)
4. When willing to approximate



# Chan-Vese in 3D

- Energy functional with surface area regularization:

$$E(\vec{C}) = \frac{1}{2\sigma_1^2} \iiint_{R_1} (y - \mu_1)^2 d\mathbf{x} + \frac{1}{2\sigma_2^2} \iiint_{R_2} (y - \mu_2)^2 d\mathbf{x} \\ + \alpha \iint_{\vec{C}} dA$$

- With a slice-based model, we can write the regularization term as:

$$\iint_{\vec{C}} dA = \sum_{i=1}^{n-1} \iint_{\vec{c}_i \oplus \vec{c}_{i+1}} dA$$

where  $\vec{c}_i \oplus \vec{c}_{i+1}$  is the surface between  $c_i$  and  $c_{i+1}$



# Zero-order hold approximation

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- Approximate volume as piecewise-constant “cylinders”:

$$\vec{C}(s, z) = \vec{c}_i(s), \quad \forall |z - i\Delta z| < \frac{\Delta z}{2}$$

- Then we see that the surface areas are:

$$\iint_{\vec{c}_i \oplus \vec{c}_{i+1}} dA = \frac{\Delta z}{2} \oint_{\vec{c}_i} ds + \frac{\Delta z}{2} \oint_{\vec{c}_{i+1}} ds + \iint_{R_{i,i+1}^{\text{diff}}} d\mathbf{x}$$

- We see terms related to the curve length and the difference between neighboring slices
- Upper bound to correct surface area



# Overall regularization term

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- Adding everything together results in:

$$\oint \vec{c} dA = \underbrace{\Delta z \sum_{i=1}^n \oint_{\vec{c}_i} ds}_{\text{self potentials}} + \underbrace{\sum_{i=1}^{n-1} \int \int_{R_{i,i+1}^{\text{diff}}} dx}_{\text{edge potentials}}$$



## 2.5D Approach

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- In 3D world, natural (or built-in) partition of volumes into slices
- Assume Markov relationship among slices
- Then have local potentials (e.g., PCA) and edge potentials (coupling between slices)
- Naturally lends itself to local Metropolis-Hastings approach (iterating over the slices)

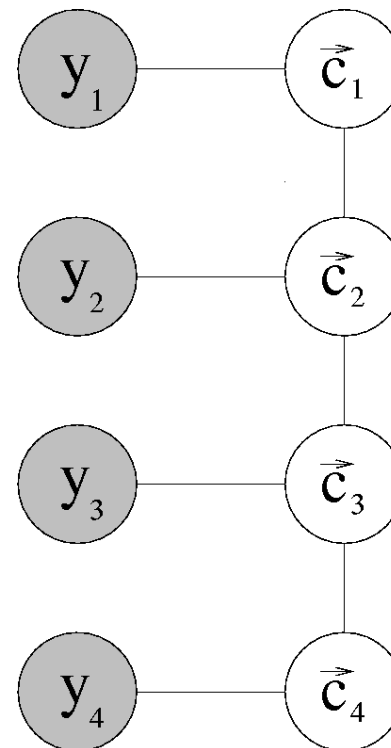


## 2.5D Model

- We can model this as a simple chain structure with pairwise interactions
- This admits the following factorization:

$$p(Y|\vec{C}) = \prod_{i=1}^n p(y_i|\vec{c}_i)$$

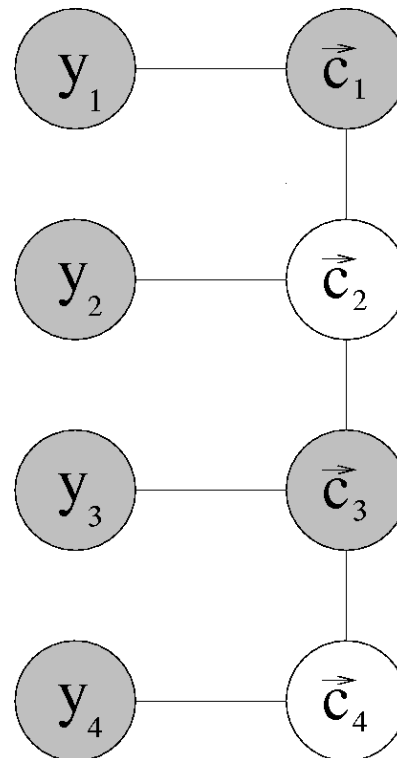
$$p(\vec{C}) = \prod_{i=1}^n \psi_i(\vec{c}_i) \prod_{i=1}^{n-1} \psi_{i,i+1}(\vec{c}_i, \vec{c}_{i+1})$$





# Partial segmentations

- Assume that we are given segmentations of every other slice
- We now want to sample surfaces conditioned on the fact that certain slices are fixed
- Markovianity tells us that  $c_2$  and  $c_4$  are independent conditioned on  $c_3$







## Log probability for $c_2$

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- We can then construct the probability for  $c_2$  conditioned on its neighbors using the potential functions defined previously:

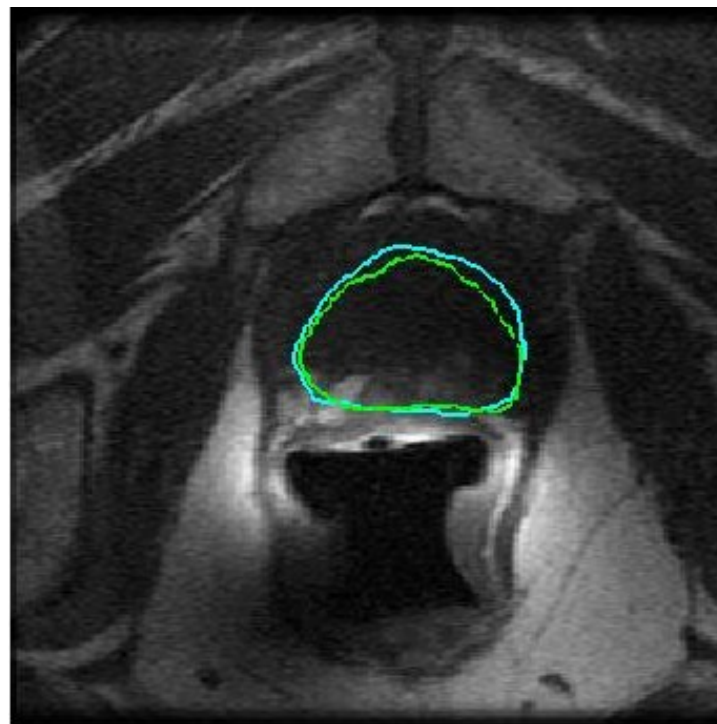
$$\ell(\vec{c}_2 | y_2, \vec{c}_1, \vec{c}_3) = \ell(y_2 | \vec{c}_2) + \alpha \left[ \Delta z \oint_{\vec{c}_2} ds + d_{\text{SAD}}(\vec{c}_2, \vec{c}_1) + d_{\text{SAD}}(\vec{c}_2, \vec{c}_3) \right]$$



# Results



Neighbor  
segmentations



Our result (cyan)  
with expert (green)



# Larger spacing

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- Apply local Metropolis-Hastings algorithm where we sample on a slice-by-slice basis
- Theory shows that asymptotic convergence is unchanged
- Unfortunately larger spacing similar to less regularization
- Currently have issues with poor data models that need to be resolved



# Full 3D

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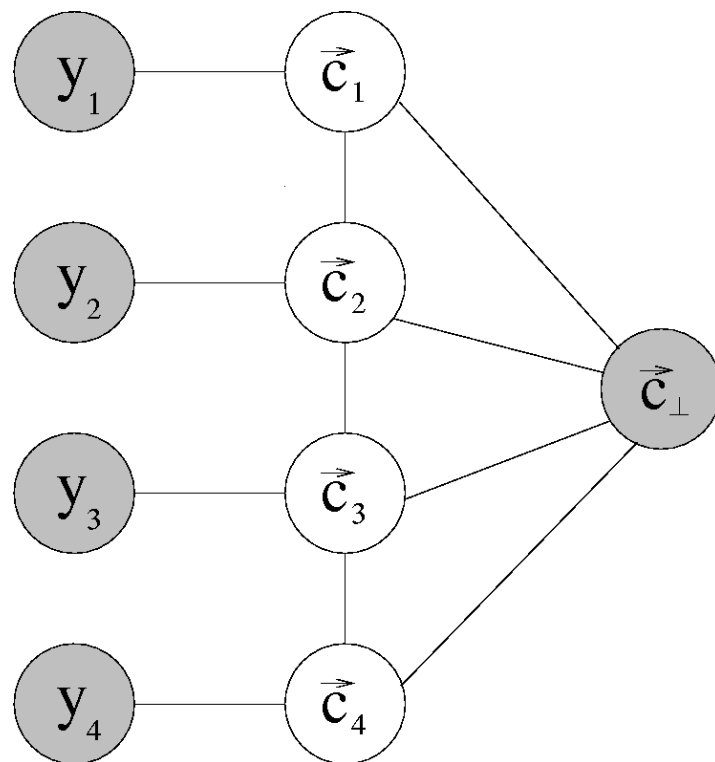
- In medical imaging, have multiple slice orientations (axial, sagittal, coronal)
- In seismic, vertical and horizontal shape structure expected
- With a 2.5D approach, this introduces complexity to the graph structure



# Incorporating perpendicular slices

- $c_{\perp}$  is now coupled to all of the horizontal slices
- $c_{\perp}$  only gives information on a subset of each slice
- Specify edge potentials as, e.g.:

$$\Psi_{i,\perp}(\vec{c}_i, \vec{c}_{\perp}) = \exp\left(-\int_0^{L_y} (\Gamma(x_0, y, i\Delta z) - \Gamma_{\perp}(x_0, y, i\Delta z))^2 dy\right)$$





# Other extensions

---

- Additional features can be added to the base model:
  - Uncertainty on the expert segmentations
  - Shape models (semi-local)
  - Exclusion/inclusion regions
  - Topological change (through level sets)



# Conclusion

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- Computationally feasible algorithm to sample from space of curves
- Approximate detailed balance
- Demonstrated utility for robustness to noise, multimodal distributions, displaying uncertainty
- Can generate arbitrary shapes with relatively complex geometry
- Conditional simulation provides a natural framework to incorporate partial user segmentations on a slice-by-slice level



## Further research

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- Multiple curves, known/unknown topology
- We would like  $q$  to naturally sample from space of smooth curves (different perturbation structure)
- Speed always an issue for MCMC approaches
  - Multiresolution perturbations
  - Parameterized perturbations (efficient basis)
  - Hierarchical models
- How to properly use samples?