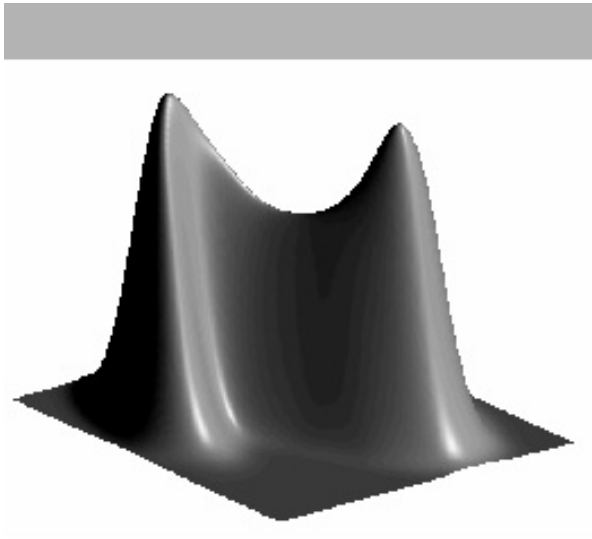
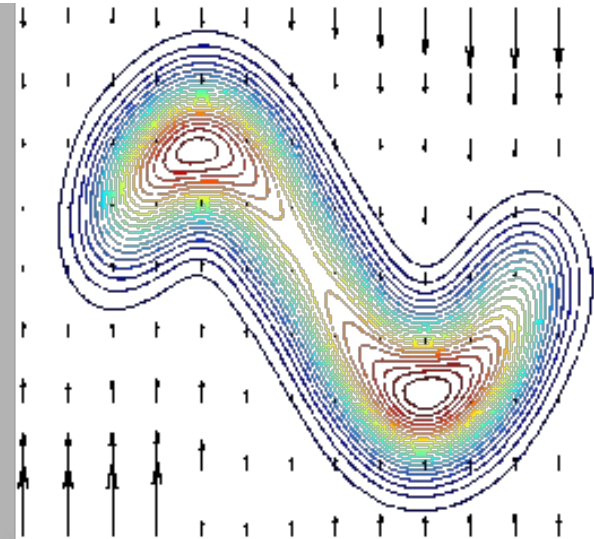


Markov Random Fields and Bayesian Image Analysis

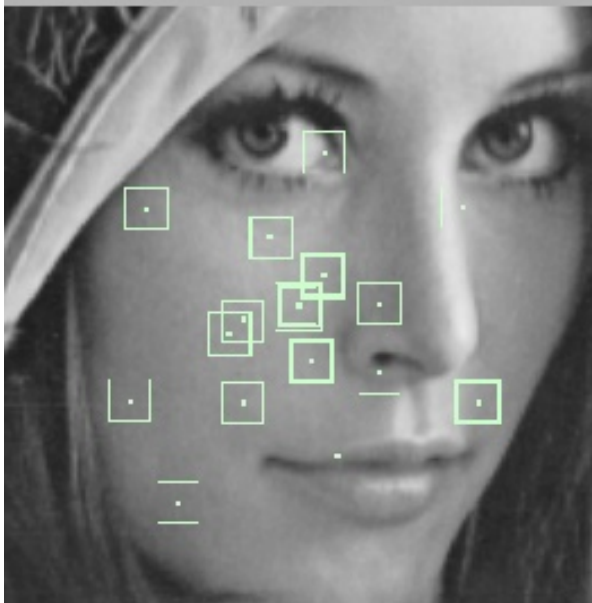
Wei Liu
Advisor: Tom Fletcher



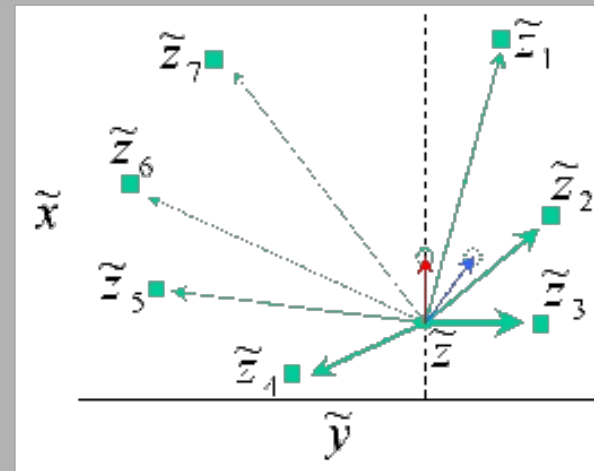
(a)



(b)



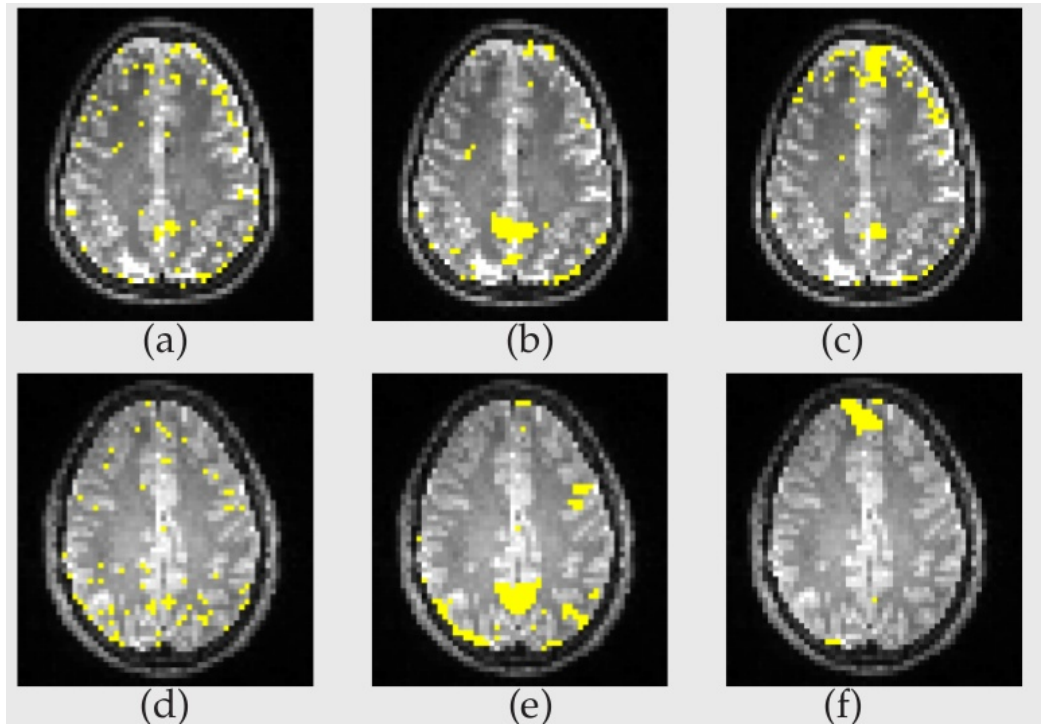
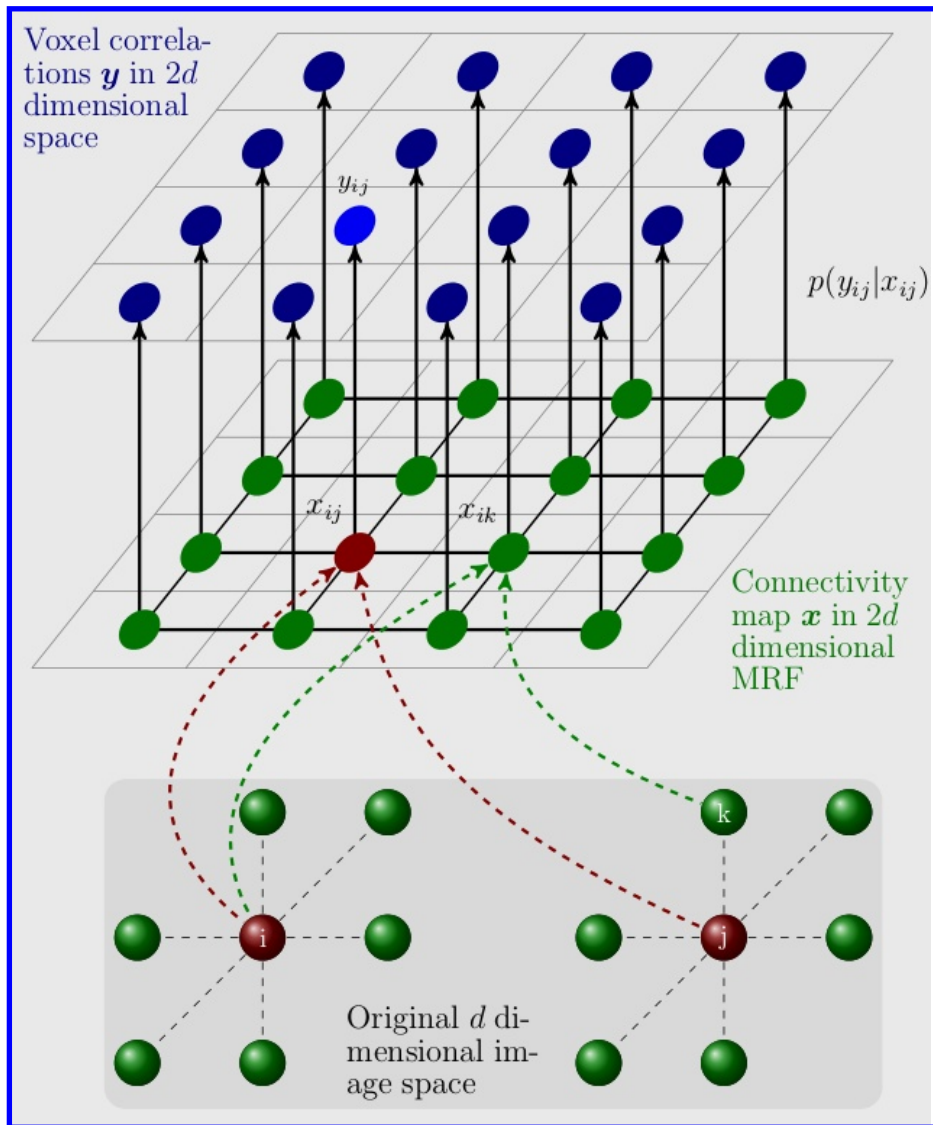
(c)

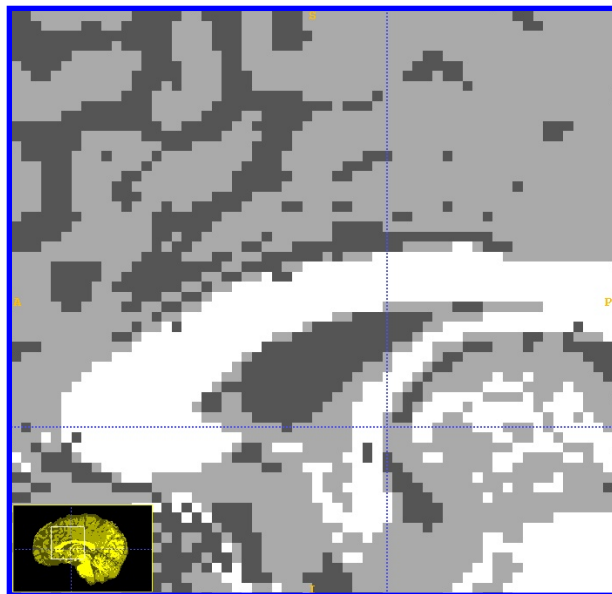
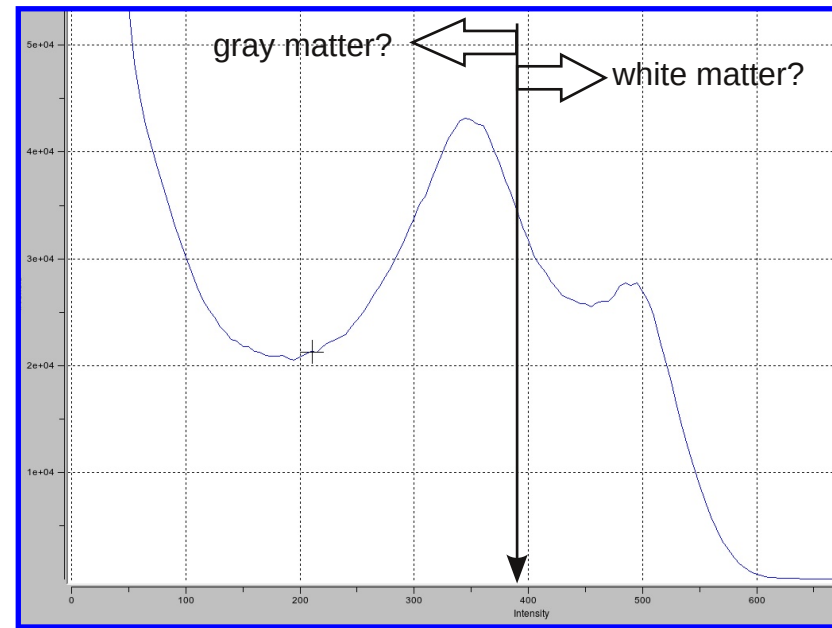
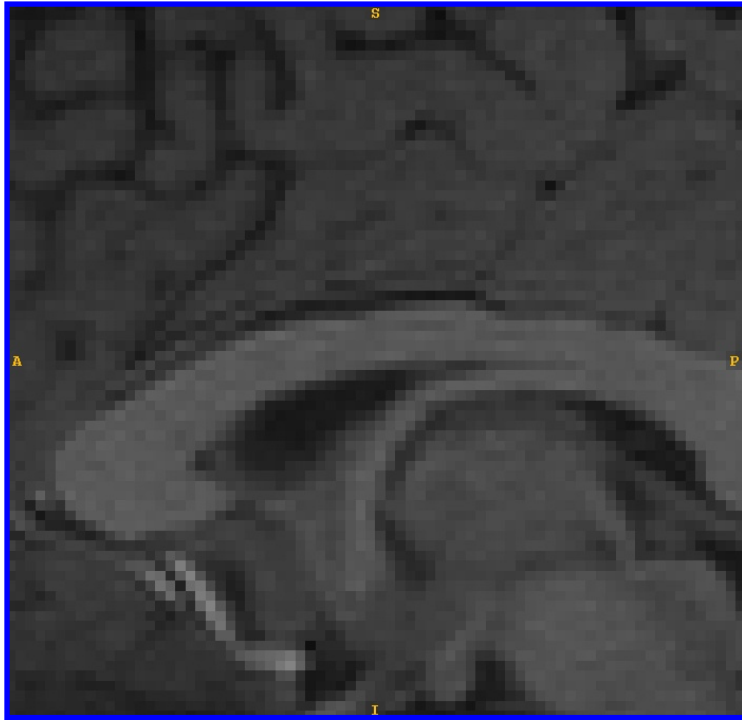


(d)

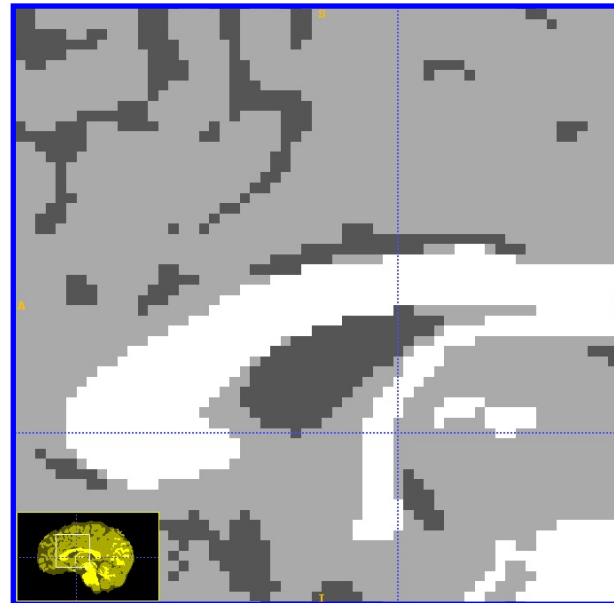
Awate and Whitaker 2006

Markov Random Field: Application Overview





without spatial MRF prior



with spatial MRF prior

- Unknown 'true' image X
- observed data Y
- Model $\mathcal{M}$ and parameter set θ

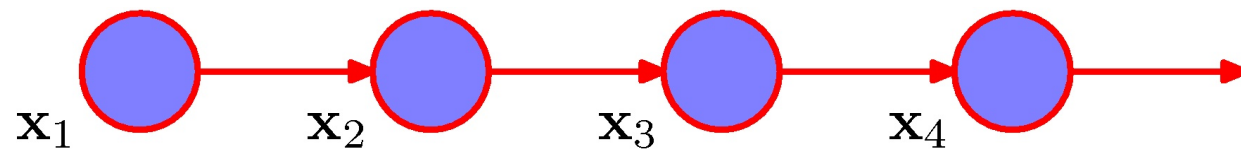
Goal: Estimate X from Y based on some objective function.

Definition 1. A markov chain is a sequence of random variables $X_1, X_2, X_3, \dots$ with the Markov property that given the present state, the future and past states are **conditionally** independent.

$$P(X_{n+1}|X_1, X_2, \dots, X_n) = P(X_{n+1}|x_n)$$

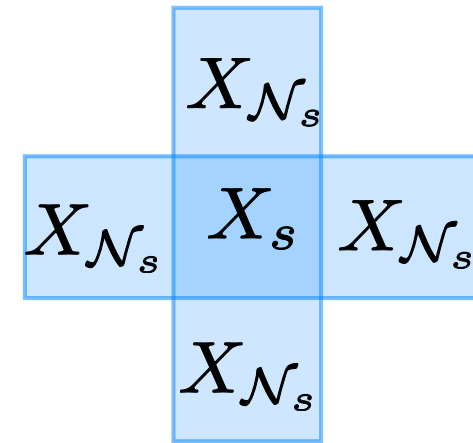
The joint probability of the sequence is given by

$$P(X) = P(X_0) \prod_{n=1}^N P(X_n|X_{n-1})$$



Define

- $\mathcal{S} = \{1, \dots, M\}$ the set of lattice points.
- $s \in \mathcal{S}$ a site in $\mathcal{S}$
- $\mathcal{L} = \{1, \dots, L\}$ the set of labels
- X_s the random variable at s . $X_s = x_s \in \mathcal{L}$
- $\mathcal{N}_s$ the set of sites neighboring s . Properties of neighboring sites:
 - $s \notin \mathcal{N}_s$
 - $s \in \mathcal{N}_t \Leftrightarrow t \in \mathcal{N}_s$
- $\mathcal{S}$ and neighbor system $\mathcal{N}$ together defines a graph $(\mathcal{S}, \mathcal{N}) = \mathcal{G}$.



Definition. X is called a random field if $X = \{X_1, \dots, X_N\}$ is a collection of random variables defined on the set $\mathcal{S}$, where each X_s takes a value x_s in $\mathcal{L}$. $x = \{x_1, \dots, x_N\}$ is called a configuration of the field.

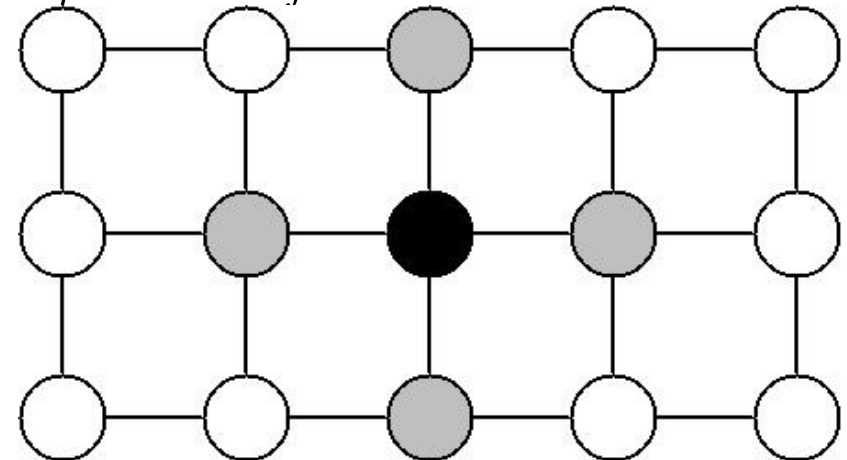
Definition. X is said to be a Markov random field on $\mathcal{S}$ with respect to a neighborhood system $\mathcal{N}$ if for all $s \in \mathcal{S}$

$$P(X_s | X_{\mathcal{S}-s}) = P(X_s | X_{\mathcal{N}_s})$$

Definition. X is homogeneous if $P(X_s | X_{\mathcal{N}_s})$ is independent of the relative location of site s in $\mathcal{S}$.

Generalization of Markov chain:

- unilateral $\rightarrow$ bilateral
- 1D $\rightarrow$ 2D
- time domain $\rightarrow$ space domain. No natural ordering on image pixels.



Advantage of MRF's

- Can be isotropic or anisotropic depending on the definition of neighbor system $\mathcal{N}$.
- Local dependencies

Disadvantages of MRF's

- difficult to compute $P(X)$ from local dependency $P(X_s|X_{\mathcal{N}_s})$
- Parameter estimation is difficult

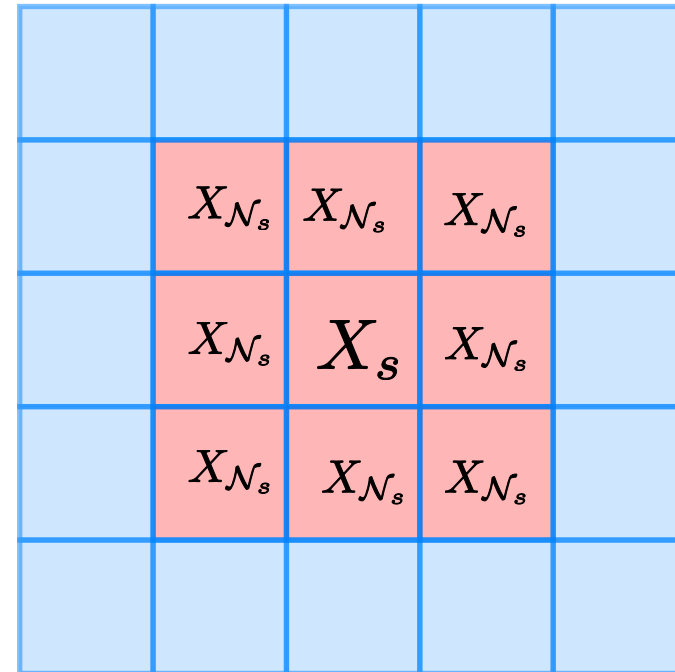
Hammersley-Clifford theorem build the relationship between local properties $P(X_s|X_{\mathcal{N}_s})$ and global properties $P(X)$.

Definition. A clique $\mathcal{C}$ is a set of points, which are all neighbors of each other

$$\mathcal{C}_1 = \{s | s \in \mathcal{S}\}$$

$$\mathcal{C}_2 = \{(s, t) | s \in \mathcal{N}_t, \quad t \in \mathcal{N}_s\}$$

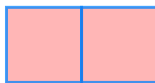
$$\mathcal{C}_3 = \dots$$



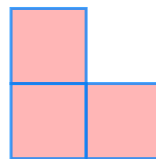
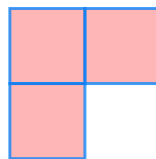
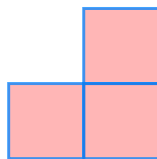
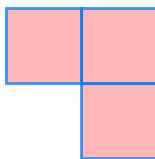
single site clique



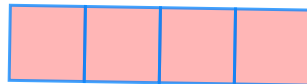
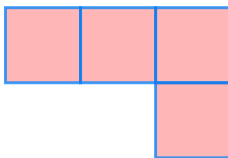
two-site clique



Three-site clique



not a clique



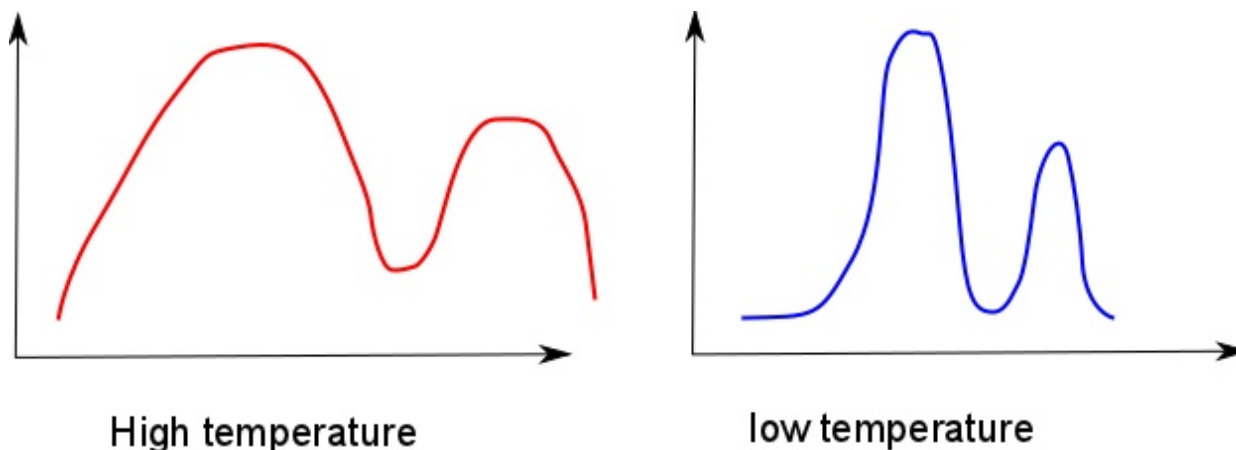
Definition. A set of random variable X is said to be a Gibbs random field (GRF) on $\mathcal{S}$ with respect to $\mathcal{N}$ if and only if its configurations obey a Gibbs distribution

$$P(X) = \frac{1}{Z} \exp\left\{-\frac{1}{T}U(X)\right\}$$

- $U(X)$ – energy function. Configurations with lower energy are more probable.

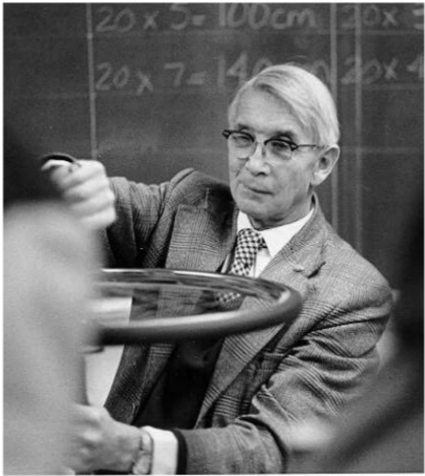
$$U(X) = \sum_{c \in \mathcal{C}} V_c(X)$$

- T – temperature. *Sharpness* of the distribution.
- Z – normalization constant. $Z = \sum_{X \in \mathcal{X}} \exp\left\{\frac{1}{T}U(X)\right\}$, $\mathcal{X} = \mathcal{L}^N$

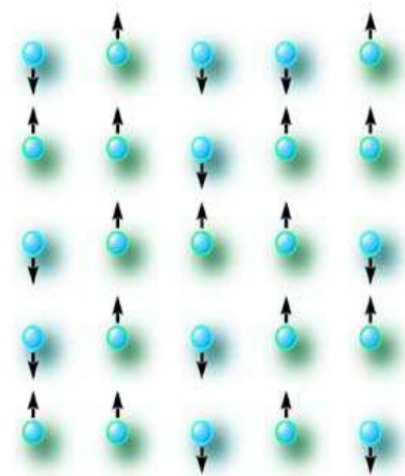


Theorem. *X is an Markov random field on $\mathcal{S}$ if and only if X is a Gibbs field on $\mathcal{S}$ with respect to $\mathcal{N}$.*

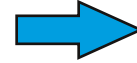
- Gives a method to specify joint probability by specifying the clique potential $V_c(X)$.
- Different clique potential gives different MRFs.
- Z is still difficult to compute.



(a)



(b)



1	1	1	1	1	1	1	1
1	1	1	1	1	1	1	1
1	1	1	1	1	-1	-1	1
1	1	1	1	-1	-1	-1	1
1	1	1	-1	-1	-1	-1	1
1	1	1	1	-1	-1	1	1
1	1	1	-1	-1	-1	-1	1
1	1	1	1	1	1	1	1

X_r	
X_s	
X_r	X_s

cliques

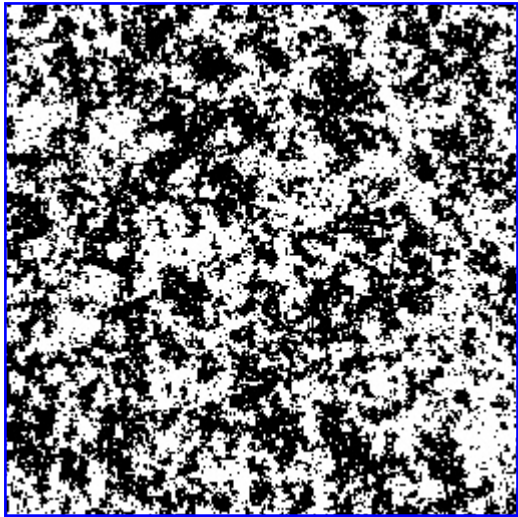
- Two state: $\mathcal{L} = \{-1, +1\}$
- Clique potential $V(X_r, X_s) = -\beta X_r X_s$

•

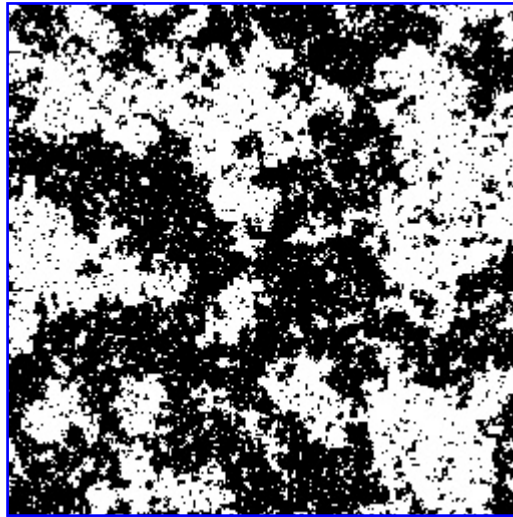
$$U(X) = \sum_{c \in \mathcal{C}} V_c(X) = -\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s, \quad P(X) = \frac{1}{Z} \exp\left\{-\frac{U(X)}{kT}\right\}$$

- Conditional distribution at site X_s :

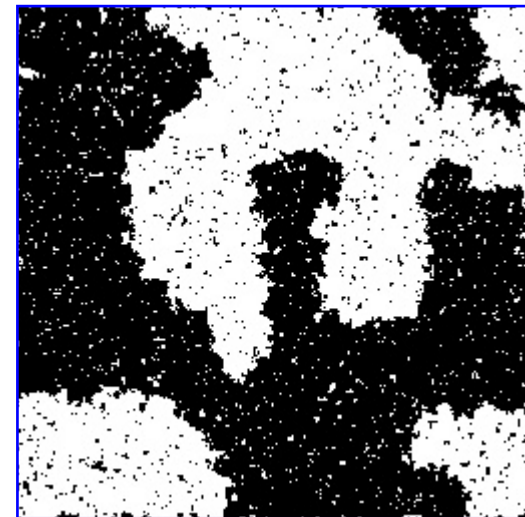
$$P(X_s | X_{\mathcal{N}_s}) = \frac{\exp\{\beta X_s \sum_{r \in \mathcal{N}_s} X_r\}}{2 \cosh(\beta \sum_{r \in \mathcal{N}_s} X_r)}$$



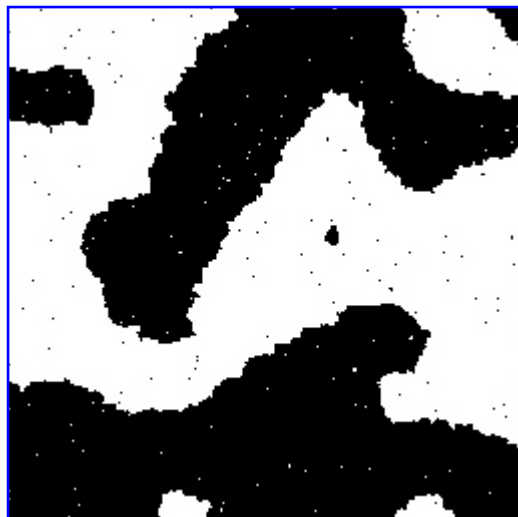
Beta = 0.8



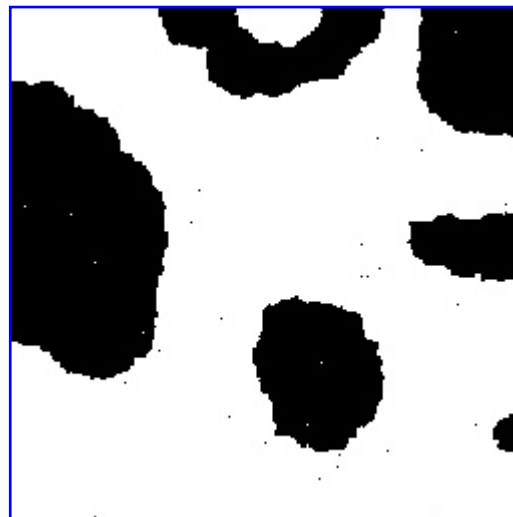
Beta = 0.88



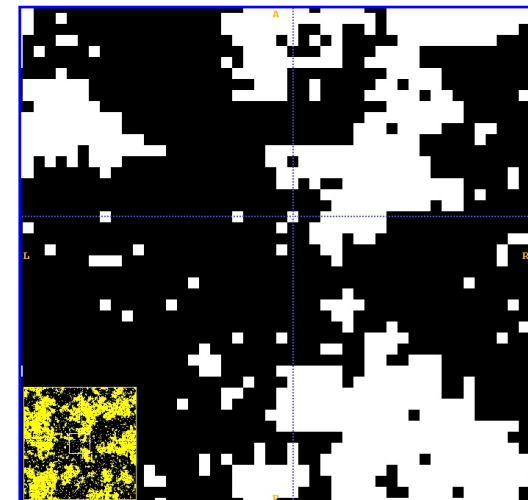
Beta = 1.0



Beta = 1.5

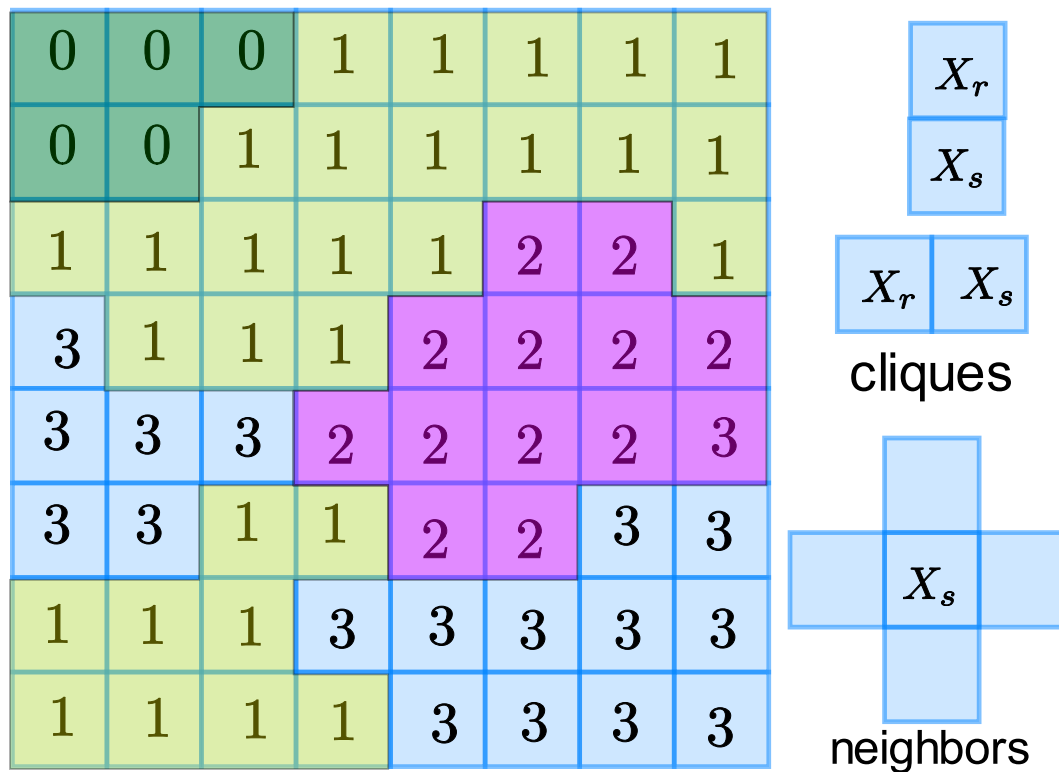


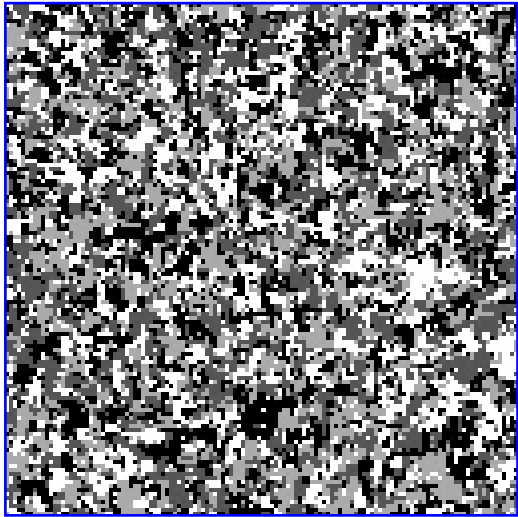
Beta = 2.0



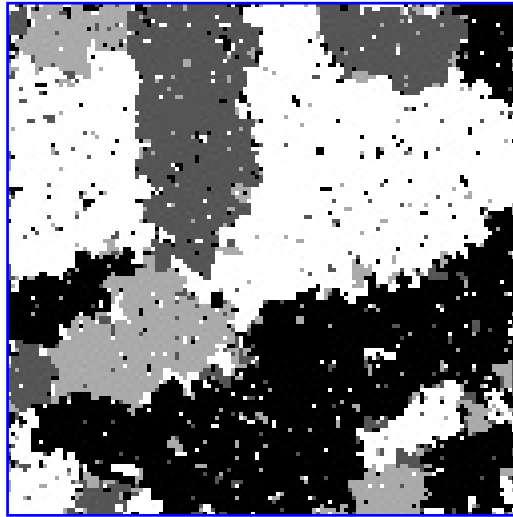
Beta = 0.88. detailed view

- Multiple state: $X_s = x_s \in \mathcal{L}$, $\mathcal{L} = \{1, 2, \dots, L\}$
- 4-neighbor or 8-neighbor system
- $V_1(X_s = l) = \alpha_l$, $l \in \mathcal{L}$
- $V_2(X_r, X_s) = \begin{cases} \beta & X_r \neq X_s \\ 0 & X_r = X_s \end{cases}$

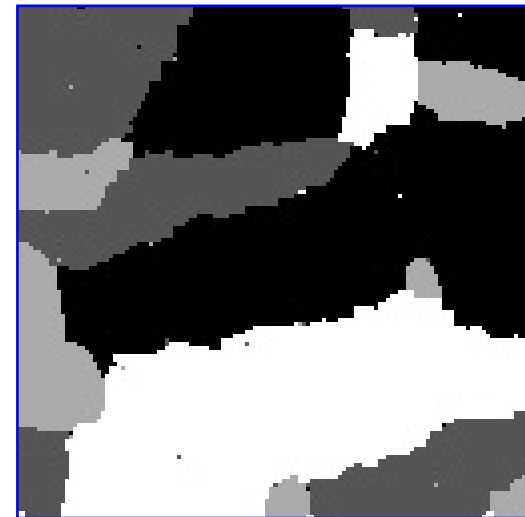




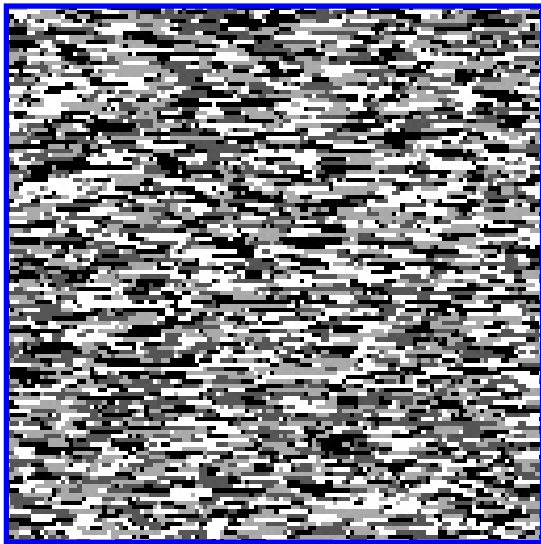
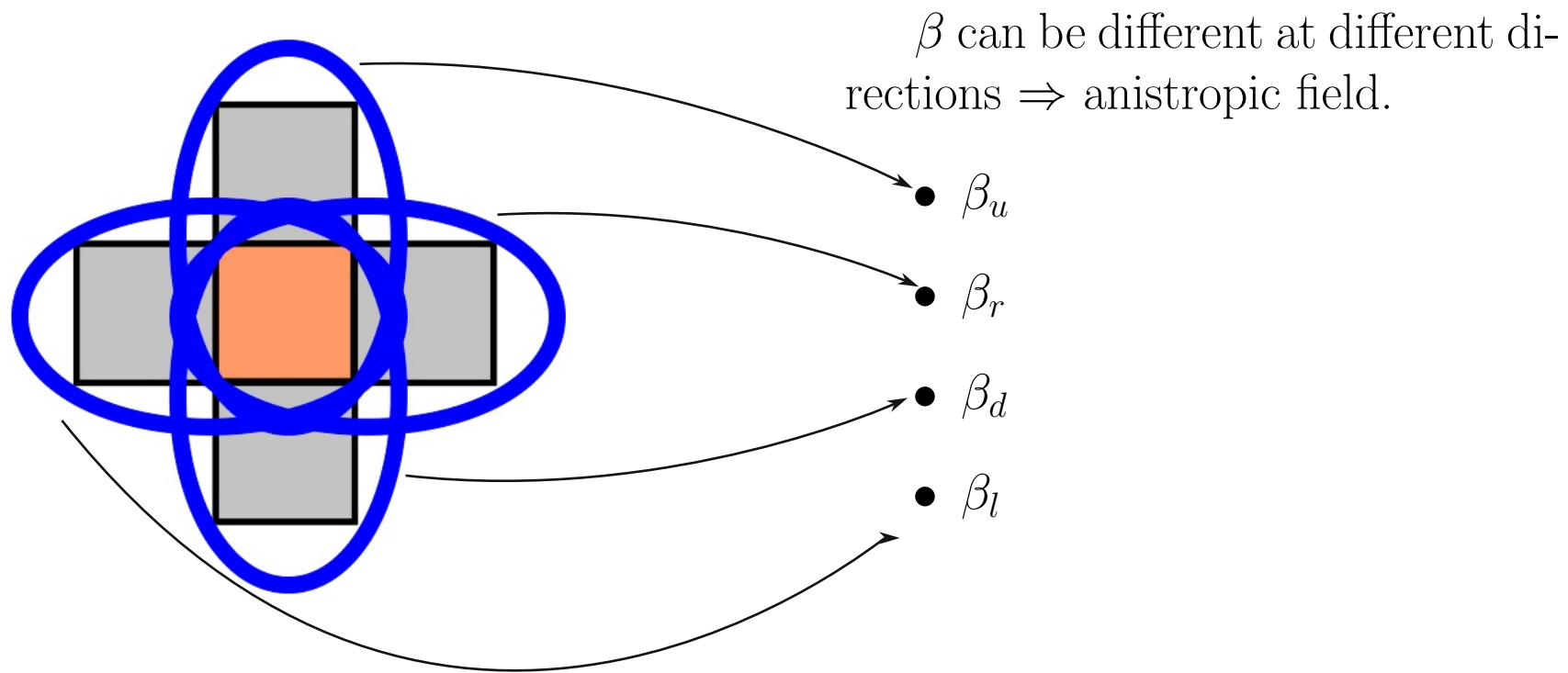
Beta = 0.88



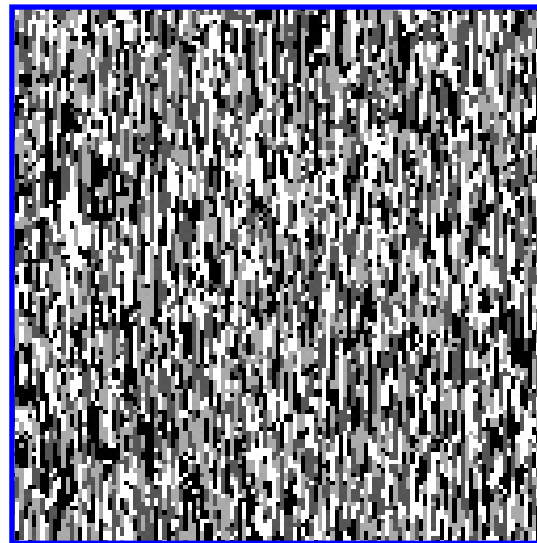
Beta = 1.2



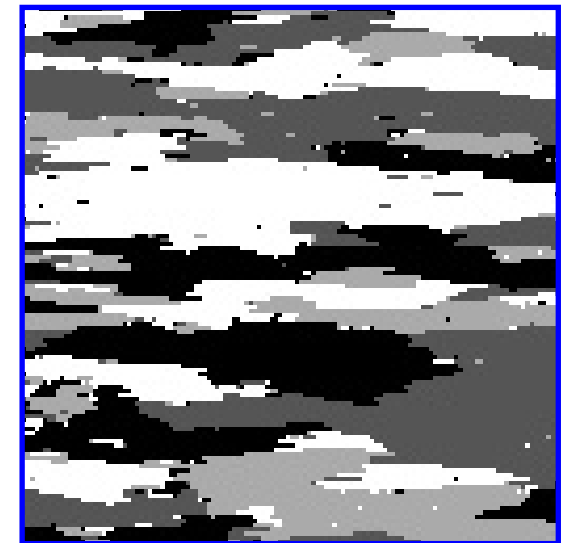
Beta = 2.0



$$\beta_u = \beta_d = 0, \beta_l = \beta_r = 2$$

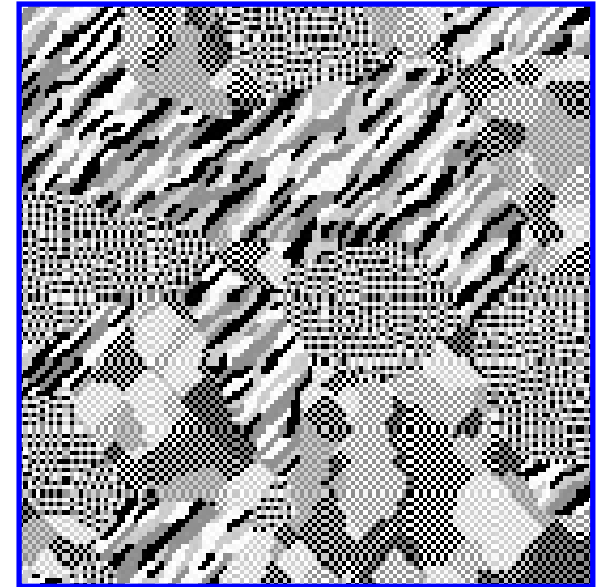
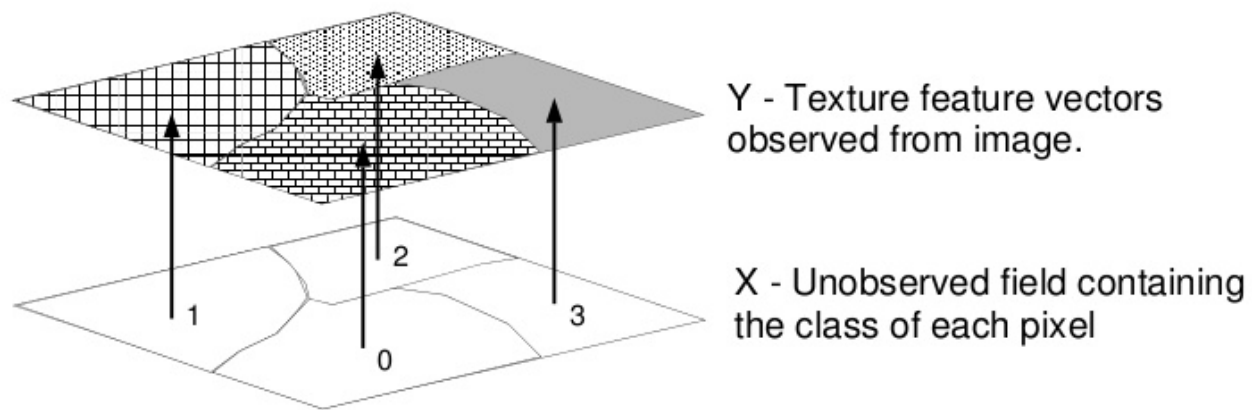


$$\beta_u = \beta_d = 2, \beta_l = \beta_r = 0$$



$$\beta_u = \beta_d = 1, \beta_l = 4, \beta_r = 2$$

Hierarchical MRF Model



- $X \in \mathcal{L}^N$ is MRF – region configuration.
- $P(Y_s|X_s)$ depends on $Y_{\mathcal{N}_s}$.
- Given X_s , $\{s, \mathcal{N}_s\}$ has same texture type.

Why do we want draw a sample of MRFs (or Gibbs distribution)?

$$P(X) = \frac{1}{Z} \exp\{-U(X)\}$$

- Compare simulated image with real image $\Rightarrow$ Model is good?
- Texture synthesis
- Model verification.
- Monte Carlo integration

Review of Monte Carlo integration. Consider the generic problem of evaluating the integral

$$\mathbb{E}_{f(x)} = \int_{\mathcal{X}} h(x) f(x) dx$$

We can use a set of samples $(x_1, x_2, \dots, x_M)$ generated from density $f(x)$ to approximate above integral by the empirical average

$$\bar{h} = \frac{1}{M} \sum_{m=1}^M h(x_m)$$

- Metropolis sampler. Used when we know $P(X)$ up to a constant
- Gibbs Sampler. Used when we know exactly $P(X)$

Goal: draw samples from some distribution $P(X)$ where $P(X) = f(X)/K$.

- Start with any initial value X_0 satisfying $f(X_0) > 0$.
- Sample a candidate point X^* from distribution $g(X)$ (proposal distribution).

- Calculate the

$$\alpha = \frac{P(X^*)}{P(X_{t-1})} = \frac{f(X^*)}{f(X_{t-1})}$$

- If $\alpha > 1$, accept candidate point and set $X_t = X^*$. Otherwise accept X^* with probability α .

We don't have to know the constant K !

Goal: draw samples from Gibbs distribution $P(X) = \frac{1}{Z} \exp\{-U(X)\}$.

1. Randomly init X^0 satisfying $f(X^0) > 0$ (X is the whole iamge)
2. For $s \in \mathcal{S}$ do step 3, 4, 5
3. Generate a *univariate* sample X_s^* from proposal probability $Q(X_s^*|X^{t-1})$ (Q can be uniform distribution), and replace X_s with X_s^* to get candidate X^* . X^{t-1} and X^* differs only at X_s .
4. Calculate the

$$\Delta U(X^*) = U(X^*) - U(X^{t-1}) = U(X_s^*) - U(X_s)$$

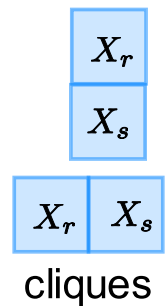
5. If $\Delta U(X^*) < 0$, accept cadidate point and set $X^t = X^*$. Otherwise accept X^* with probability $\exp\{-\Delta U(X^*)\}$.
6. Repeat above steps M times.

The sequence of random fields X^t (after burn-in period) is a Markov chain.

Goal: draw samples from Gibbs distribution
 $P(X) = \frac{1}{Z} \exp\{-U(X)\}$.

1. Randomly init X^0 satisfying $f(X^0) > 0$ (X is the whole iamge)
2. For $s \in \mathcal{S}$ do step 3, 4
3. Compute $P(X_s|X^{t-1}) = P(X_s|X_{\mathcal{N}_s}^{t-1})$ and draw sample X_s^* from it.
4. Accept X_s^* , i.e. replace X_s with X_s^* and obtain X^t .
5. Repeat above steps M times.

1	1	1	1	1	1	1	1
1	1	1	1	1	1	1	1
1	1	1	1	1	-1	-1	1
1	1	1	1	-1	-1	-1	1
1	1	1	-1	-1	-1	-1	1
1	1	1	1	-1	-1	1	1
1	1	1	-1	-1	-1	-1	1
1	1	1	1	1	1	1	1



The sequence of random fields X^t (after burn-in period) is a Markov chain.

Gibbs:

- Always accepted.
- Have to compute $P(X_s = l | X_{\mathcal{N}_s})$ for all $l \in \mathcal{L}$.

Metropolis:

- Expected acceptance rate is $1/L$ – low when L is large $\Rightarrow$ more burn-in time.
- No need to compute $P(X_s = l | X_{\mathcal{N}_s})$ for all $l \in \mathcal{L}$. Only compute $U(X_s^* | X_{\mathcal{N}_s})$ for candidate X^* .

Image Segmentation:

- $X \in \mathcal{L}^N$: Image labels we're interested
- $Y \in \mathcal{R}^N$: noise data (observed image)

Goal: Estimate X from Y .

Image Denoising:

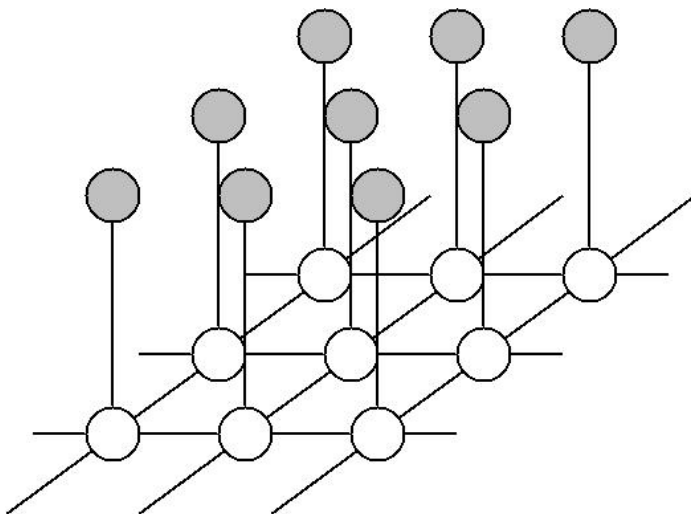
- $X \in \mathcal{R}^N$: True image intensity.
- $Y \in \mathcal{R}^N$: noise data (observed image)

Goal: Recover X from Y .

$$P(X|Y) \propto P(X) + P(Y|X)$$

MRF prior

$$P(X) = \frac{1}{Z} \exp U(X)$$



Conditional likelihood

For Segmentation:

$$P(Y|X) = \sum_{s \in \mathcal{S}} P(Y_s | X_s = l) = \mathcal{N}(\mu_l, \sigma_l^2)$$

For denoising:

$$P(Y|X) = \sum_{s \in \mathcal{S}} P(Y_s | X_s = x_s) = \mathcal{N}(x_s, \sigma_l^2)$$

- Define a model.

$$P(X) = \frac{1}{Z} \exp\{U(X)\}$$
$$P(Y|X) = \sum_{s \in \mathcal{S}} P(Y_s | X_s = l) = \mathcal{N}(\mu_l, \sigma_l^2)$$

- Formulation of objective function. Optimal Criteria.
- Search solution in the admissible space.

- Bayesian Risk is defined as

$$R(X^*) = \int_{X \in \mathcal{X}} C(X, X^*) P(X|Y) dX$$

- $C(X, X^*)$: cost function. X : true value. X^* : estimated value.

- $C(X, X^*) = ||X - X^*||^2 \Rightarrow X^* = \int_{X \in \mathcal{X}} P(X|Y) dX$ (Posterior mean)

- $C(X, X^*) = \begin{cases} 0 & ||X - X^*|| \leq \delta \\ 1 & \text{otherwise} \end{cases} \Rightarrow X^* = \operatorname{argmax}_{X \in \mathcal{X}} P(X|Y) =$
 $\operatorname{argmax}_{X \in \mathcal{X}} (P(X) + P(Y|X))$. This is mode of posterior.

Image Segmentation. Two classes $\mathcal{L} = \{-1, 1\}$

- Prior is Ising model

- $P(X) = \frac{1}{Z} \exp\{U(X)\}$, $U(X) = -\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s$. Assume T and K is 1.

- $P(X_s | X_{\mathcal{N}_s}) = \frac{\exp\{\beta X_s \sum_{r \in \mathcal{N}_s} X_r\}}{2 \cosh(\beta \sum_{r \in \mathcal{N}_s} X_r)}$

- Conditional likelihood $P(Y|X) = \prod_{s \in \mathcal{S}} P(Y_s | X_s)$, $P(Y_s | X_s = l) = \mathcal{N}(\mu_l, \sigma_l^2)$

- objective function:

$$\begin{aligned} \log P(X|Y) &\propto \log P(X) + \log P(Y|X) \\ &= -\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s - \log(Z) + \sum_{s \in \mathcal{S}} \frac{(Y_s - \mu_l)^2}{2\sigma_l^2} - \log(\sigma_l) + \text{const} \end{aligned}$$

- Combinatorial optimization problem. NP hard.

(Approximation) Optimization method:

- Iterated Conditional Modes
- Simulated Annealing
- Graph-cuts

Strategy:

- constrained minimization $\Rightarrow$ unconstrained minimization (Lagrange multiplier).
- discrete labels $\Rightarrow$ continuous labels (Relaxation labeling).

1. Init X by maximum likelihood $X^0 = \operatorname{argmax}_{X \in \mathcal{X}} P(Y|X)$
2. For $s \in \mathcal{S}$, update X_s by

$$X_s^{t+1} = \operatorname{argmax}_{X_s \in \mathcal{L}} \log P(X_s | X_{\mathcal{N}_s}^t, Y_s).$$

For the Ising-Gaussian case, this is

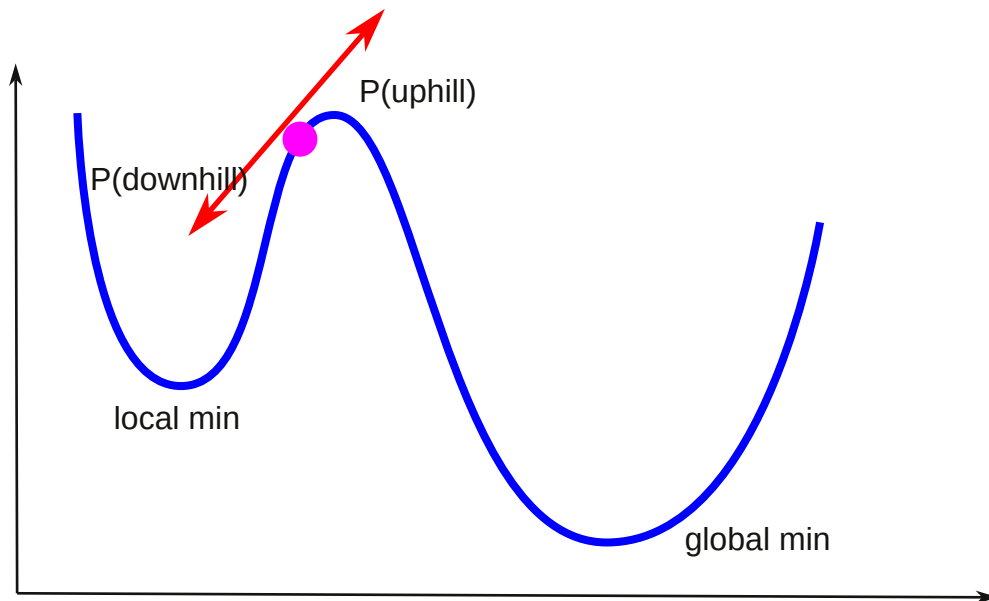
$$\begin{aligned} X_s^{t+1} &= \operatorname{argmax}_{X_s \in \mathcal{L}} \log P(X_s | X_{\mathcal{N}_s}) + \log P(Y_s | X_s) \\ &= \operatorname{argmin}_{X_s \in \mathcal{L}} \left\{ -\beta X_s \sum_{r \in \mathcal{N}_s} X_r^t + \frac{(Y_s - \mu_l)^2}{2\sigma^2} + \log(\sigma_l) \right\}. \end{aligned}$$

Note μ_l and σ_l is function of X_s , and $\log(Z_s) = \log(2 \cosh(\beta \sum_{r \in \mathcal{N}_s} X_r^t))$ is not a function of X_s .

3. Do above step for all $s \in \mathcal{S}$.
4. Repeat 2 and 3 until converge.

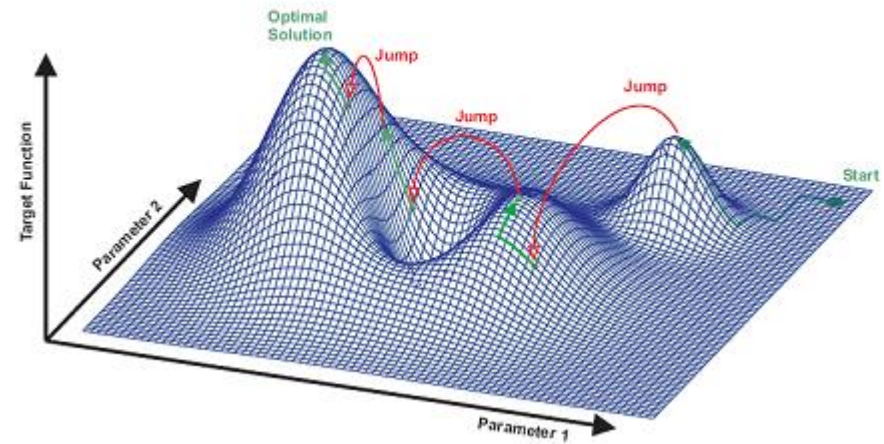
- Greedy algorithm $\Rightarrow$ local minimum.
- Sensitive to initialization.
- Quick convergence.

Simulated Annealing



- Not always downhill moving.
- Global minimum with enough scan.

Simulated Annealing



Simulated Annealing

- Assuming Ising+Gaussian model

$$\begin{aligned} P(X|Y) &\propto P(X) \cdot P(Y|X) \\ &= \frac{1}{Z} \exp\left\{\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s\right\} \cdot \prod_{s \in \mathcal{S}} \exp\left\{-\frac{(X_s - \mu_l(X_s))^2}{2\sigma_l^2(X_s)} - \log(\sqrt{2\pi}\sigma_l(X_s))\right\} \\ &= \frac{1}{Z_P} \exp\{-U_P(X|Y)\} \\ U_P(X|Y) &= -\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s + \frac{(X_s - \mu_l(X_s))^2}{2\sigma_l^2(X_s)} + \sqrt{2\pi}\sigma_l(X_s) \end{aligned}$$

Posterior distribution $P(X|Y)$ is also Gibbs.

Simulated Annealing cont.

Goal: Find $\operatorname{argmax}_{X \in \mathcal{X}} P(X|Y)$

- Introduce temperature T :

$$P(X|Y) = \frac{1}{Z_P} \exp \{U_P(X|Y)\} \Rightarrow P(X|Y) = \frac{1}{Z_P} \exp \left\{ \frac{U_P(X|Y)}{T} \right\}$$

1. Init with X^0 and a high temperature T .
2. Draw samples from $P(X|Y)$ by Gibbs Sampling or Metropolis Sampling (by sample from $P(X_s|Y_s), \forall s \in \mathcal{S}$).
3. Decrease T and repeat step 2.
4. Repeat step 2 and 3 until T is low enough.

Why this works?

Energy minimization for Segmentation

	Variational methods (optimization in $\mathcal{R}^\infty$)	Combinatorial methods (optimization in $\mathcal{Z}^n$)
explicit boundary representation	<i>Snakes & Balloons</i> (variational formulations) e.g. (Kass et al., 1988; Cohen, 1991)	<i>Dynamic Programming</i> and “ <i>path-based</i> ” <i>graph methods</i> (2D only) (e.g., Amir et al., 1990; Geiger et al., 1995; Mortensen and Barrett, 1998; Falcão et al., 1998; Jermyn and Ishikawa, 1999)
implicit boundary representation	<i>Level-sets</i> (e.g., Sethian, 1999; Osher and Fedkiw, 2002; Sapiro, 2001; Osher and Paragios, 2003)	Combinatorial Graph Cuts (as originally outlined in Boykov and Jolly, 2001)

Boykov et. al. 2006

Graph Cuts for Ising Model

- Different with Normalized Cuts.
- For two class labeling problem, find the *global* minimum of

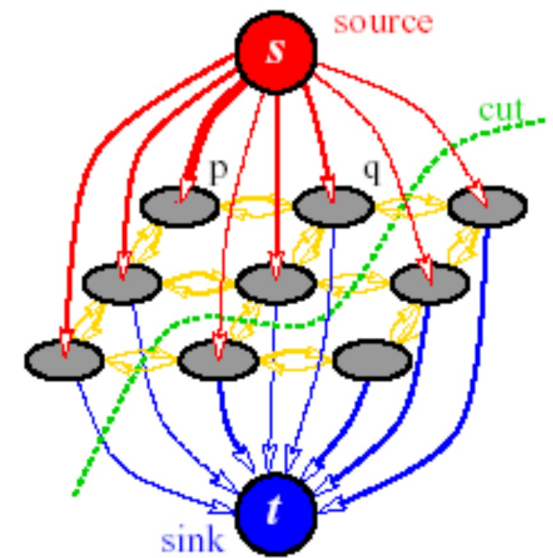
$$\log P(X|Y) = \sum_{s \in \mathcal{S}} \lambda_s X_s + \sum_{(r,s) \in \mathcal{C}_2} \beta_{(r,s)} (X_s X_r + (1 - X_s)(1 - X_r)),$$

where $\lambda_s = \log P(Y_s | X_s = 1) / P(Y_s | X_s = 0)$.

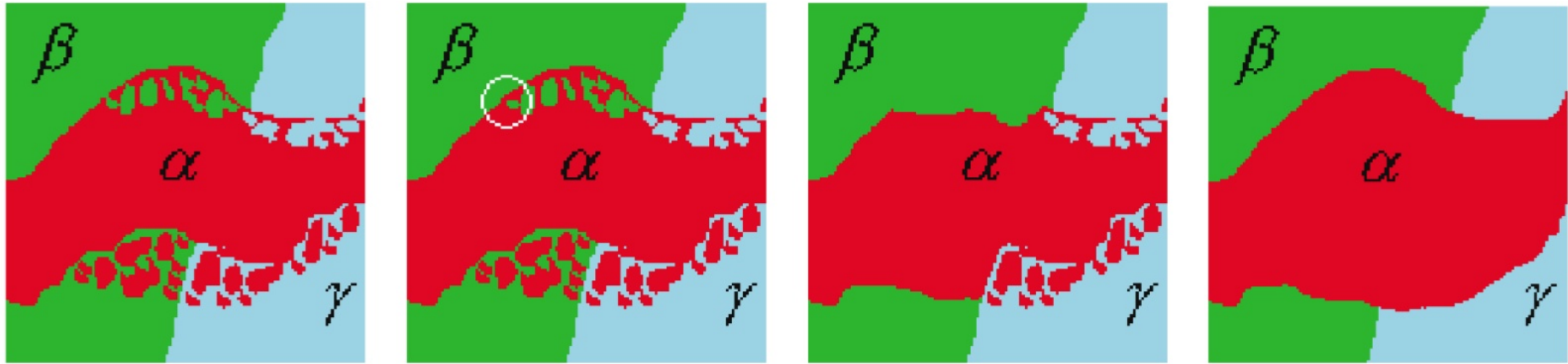
- Define a graph $\mathcal{G} = (\mathcal{V}, \mathcal{E})$. $\mathcal{V} = \{\mathcal{S}, u, t\}$

$$c_{ws} = \begin{cases} \lambda_s & \lambda_s > 0 \\ -\lambda_s & \lambda_s < 0 \end{cases}, \quad c_{sr} = \beta_{(s,r)}$$

- Define partition $B = \{u\} \cup \{s : X_s = 1\}$, $W = \{t\} \cup \{s : X_s = 0\}$ and cut $C(X) = \sum_{s \in B} \sum_{r \in W} c_{rs}$.
- It can be proved $C(x) = \log P(X|Y) + \text{const}$. In words, finding a min-cut is equivalent to find the minimum of posterior $P(X|Y)$.
- Ford-Fulkerson algorithm and Push-Relabeling method can be used to find such a cut quickly.

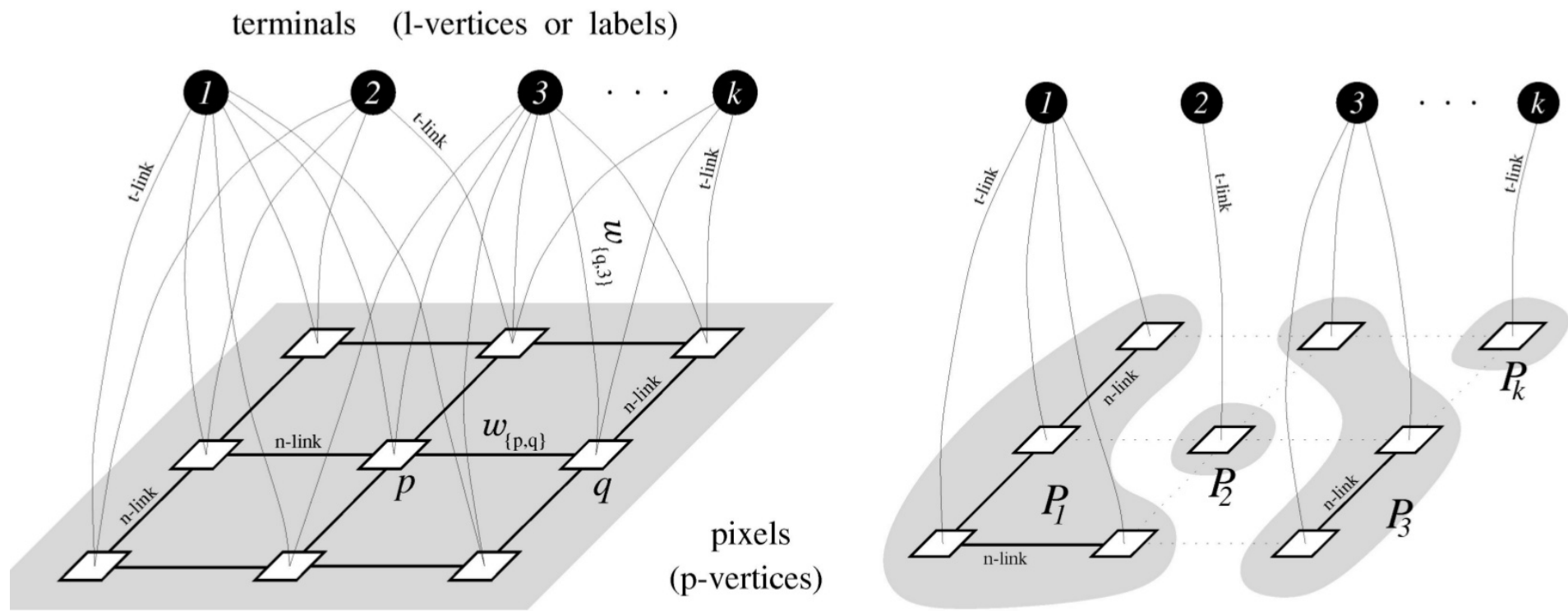


Boykov ICCV, 2005



From Left 1. Initial image. 2. standard move (ICM), 3. strong moves of alpha-beta swap. 4. strong moves of alpha expansion. (Boykov 2002).

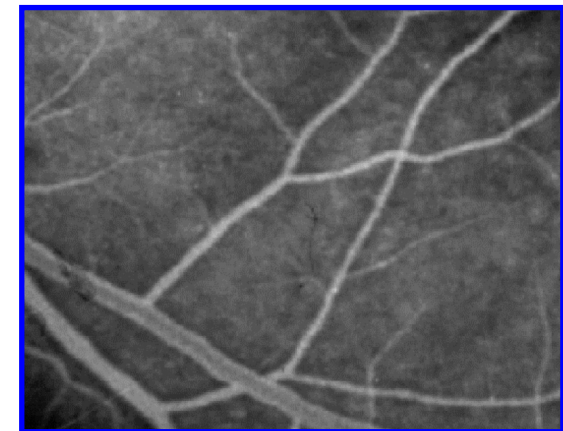
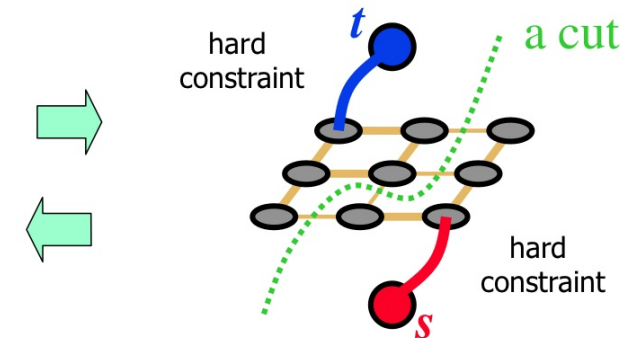
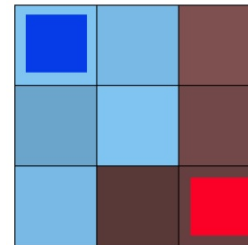
- Convert the multi-labeling problem to 2-labeling problem by $\alpha - \beta$ swap and α expansion.
- Approximation method, but with strong sense of local minima.
- Answer questions like: if results is not good, is that due to bad modeling or bad optimization algorithm?



- For label $\{\alpha, \beta\} \in \mathcal{L}$
 - Find $\hat{X} = \operatorname{argmin} E(X')$ among X' within one $\alpha - \beta$ swap of X .
 - If $E(\hat{X}) < E(X)$, accept $\hat{X}$
- Repeat above step for all pair of labels $\{\alpha, \beta\}$.

Pros:

- Break the multi-cut problem to a sequence of binary $s - t$ cuts by $\alpha - \beta$ swap and α expansion.
- Approximation method, but with strong sense of local minima.
- Easy to add hard constraints.
- Answer questions like: if results is not good, is that due to bad modeling or bad optimization algorithm?
- Parallel algorithm $\Rightarrow$ Push-Relabel algorithm.

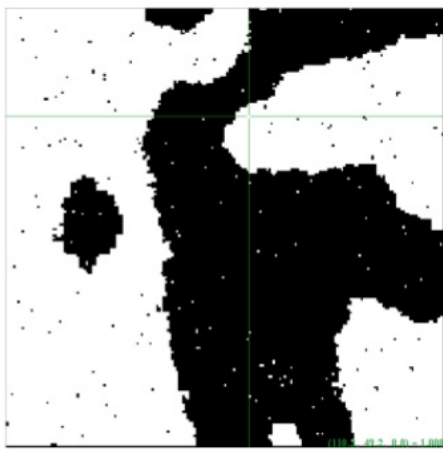


Vessels and aneurism. (kolmogorov, ICCV 2006)

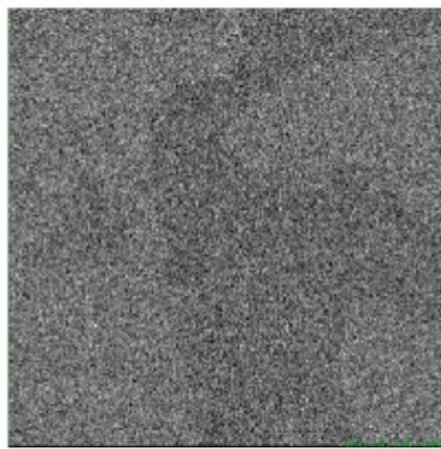
Cons:

- minimize boundary $\Rightarrow$ tends to fail for structures that are not blob shape, like vessels,

- Correct model and correct parameters $\Rightarrow$ good result.
- Correct model, and incorrect parameters $\Rightarrow$ bad result.



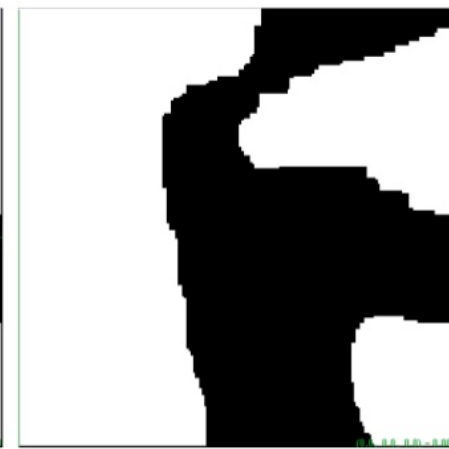
(a) true



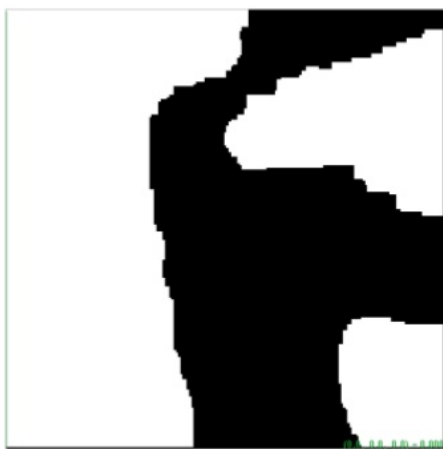
(b) with noise



(c) recovered, $\beta = 0.7$



(d) recovered, $\beta = 0.5$



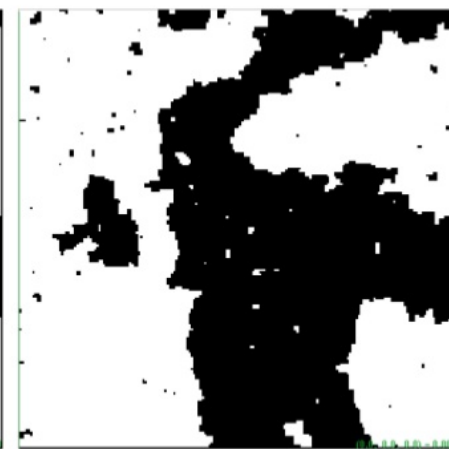
(e) recovered, $\beta = 0.4$



(f) recovered, $\beta = 0.3$



(g) recovered, $\beta = 0.2$



(h) recovered, $\beta = 0.15$

Problem 1:

Given data $X \sim MRF$, we assume model $\mathcal{M}$ with unknown parameter set θ .

Goal: Estimate θ .

Problem 2:

Given noised data Y , we assume model $\mathcal{M}$ with unknown parameter set θ .

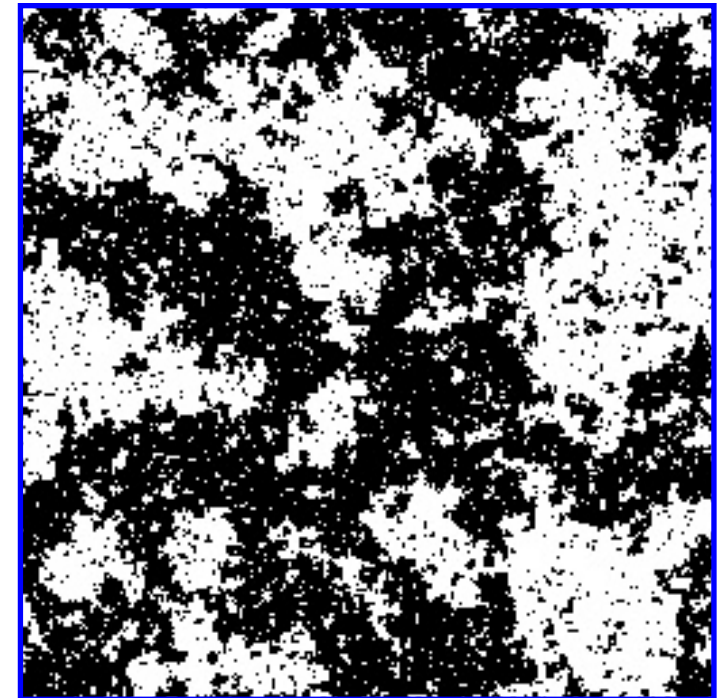
Goal: Estimate θ and hidden MRF X simultaneously.

Problem 2 is significantly harder and for now we focus on problem 1.

Given an image shown on the right and suppose we know it is generated from Ising model

$$P(X) = \frac{1}{Z} \exp\left\{-\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s\right\}.$$

Question: what is the best estimation of β ?



- Least square estimation
- Pseudo-likelihood
- Coding method

For Ising model,

- $U(X) = -\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s, \quad P(X) = \frac{1}{Z} \exp\{-U(X)\}, \quad P(X_s | X_{\mathcal{N}_s}) = \frac{\exp\{\beta X_s \sum_{r \in \mathcal{N}_s} X_r\}}{2 \cosh(\beta \sum_{r \in \mathcal{N}_s} X_r)}$

- The ratio of observed states

$$\log \left(\frac{P(X_s = 1 | X_{\mathcal{N}_s})}{P(X_s = 0 | X_{\mathcal{N}_s})} \right) = 2\beta \sum_{r \in \mathcal{N}_s} X_r$$

- For each set of neighboring pixel value $\mathcal{N}_s$, we compute
 - The observed rate of $\log \left(\frac{P(X_s=1|X_{\mathcal{N}_s})}{P(X_s=0|X_{\mathcal{N}_s})} \right)$
 - The value of $\sum_{r \in \mathcal{N}_s} X_r$.
- We have a set of over-determined linear equations and β can be solved with standard least square method.
- Easy implementation.

- Review ML estimation.
- ML estimation of θ : $\theta = \operatorname{argmax} P(X; \theta) = \operatorname{argmax} \frac{1}{Z(\theta)} \exp\{U(X; \theta)\}$. Intractable $Z(\theta)$

- Pseudo-likelihood:

$$PL(X) = \prod_{s \in \mathcal{S}} P(X_s | X_{\mathcal{N}_s})$$

does not have $Z(\theta)$ anymore.

- Solve θ by standard method, $\frac{\partial \ln PL(X; \theta)}{\partial \theta} = 0$
- For full Bayesian, if we know $P(\theta)$, the estimation is

$$\hat{\theta} = \arg \max P(\theta | X) \propto P(\theta) \cdot P(X | \theta)$$

For Ising model,

- $U(X) = -\beta \sum_{(r,s) \in \mathcal{C}_2} X_r X_s, \quad P(X) = \frac{1}{Z} \exp\{-U(X)\}, \quad P(X_s | X_{\mathcal{N}_s}) = \frac{\exp\{\beta X_s \sum_{r \in \mathcal{N}_s} X_r\}}{2 \cosh(\beta \sum_{r \in \mathcal{N}_s} X_r)}$

- The ratio of observed states

$$\log \left(\frac{P(X_s = 1 | X_{\mathcal{N}_s})}{P(X_s = 0 | X_{\mathcal{N}_s})} \right) = 2\beta \sum_{r \in \mathcal{N}_s} X_r$$

- For each set of neighboring pixel value $\mathcal{N}_s$, we compute
 - The observed rate of $\log \left(\frac{P(X_s=1|X_{\mathcal{N}_s})}{P(X_s=0|X_{\mathcal{N}_s})} \right)$
 - The value of $\sum_{r \in \mathcal{N}_s} X_r$.
- We have a set of over-determined linear equations and β can be solved with standard least square method.
- Easy implementation.