

# Gibbs Random Fields

- A set of random variables  $F$  is said to be a Gibbs random field (GRF) on  $S$  with respect to  $\mathcal{N}$  if and only if its configurations obey a Gibbs distribution
- A Gibbs distribution takes the following form

$$P(f) = Z^{-1} \times e^{-\frac{U(f)}{T}} \quad Z = \sum_{f \in \mathcal{F}} e^{-\frac{U(f)}{T}}$$

is a normalizing constant called the partition function,  $T$  is a constant called the temperature which shall be assumed to be 1 and  $U(f)$  is the energy function.

- The energy

$$U(f) = \sum_{c \in \mathcal{C}} V_c(f) \quad (1)$$

is a sum of clique potentials  $V_c(f)$  over all possible cliques  $\mathcal{C}$

- The value of  $V_c(f)$  depends on the local configuration on the clique  $c$

# Gibbs Random Fields: Property

- The Gaussian distribution is a special member of this Gibbs distribution family
- A GRF is said to be homogeneous if  $V_c(f)$  is independent of the relative position of the clique  $c$  in  $S$
- It is said to be isotropic if  $V_c$  is independent of the orientation of  $c$
- To calculate a Gibbs distribution, it is necessary to evaluate the partition function  $Z$  which is the sum over all possible configuration in  $\mathbb{F}$
- There are a combinatorial number of elements in  $\mathbb{F}$  for a discrete  $\mathcal{L}$ , the evaluation is prohibitive even for problems of moderate sizes

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# Gibbs Random Fields: Configuration

- $P(f)$  measures the probability of the occurrence of a particular configuration or pattern,  $f$
- The more probable configurations are those with lower energies
- The temperature  $T$  controls the sharpness of the distribution
- When the temperature is high, all configurations tend to be equally distributed
- Near the zero temperature, the distribution concentrates around the global energy minima
- Given  $T$  and  $U(f)$ , a class of patterns can be generated by sampling the configuration space  $\mathbb{F}$  according to  $P(f)$
- For discrete labeling problem, a clique potential  $V_c(f)$  can be specified by a number of parameters
- For continuous labeling problem,  $V_c(f)$  is continuous function of  $f_c$

# Gibbs Random Fields: Energy

- The energy of a Gibbs distribution is expressed as the sum of several terms, each related to cliques of a certain size, i.e.,

$$U(f) = \sum_{\{i\} \in \mathcal{C}_1} V_1(f_i) + \sum_{\{i, i'\} \in \mathcal{C}_2} V_2(f_i, f_{i'}) + \sum_{\{i, i', i''\} \in \mathcal{C}_3} V_3(f_i, f_{i'}, f_{i''}) + \dots$$

- An important special case is when only cliques of size up to two are considered. In this case, the energy can also be written as

$$U(f) = \sum_{i \in S} V_1(f_i) + \sum_{i \in S} \sum_{i' \in N_i} V_2(f_i, f_{i'})$$

- $\{i, i'\}$  and  $\{i', i\}$  are two distinct cliques in  $\mathcal{C}_2$  because the sites in a clique are ordered

# GRF: Equivalence

- The conditional probability can be written as

$$P(f_i | f_{\mathcal{N}_i}) = \frac{e^{-[V_1(f_i) + \sum_{i' \in \mathcal{N}_i} V_2(f_i, f_{i'})]}}{\sum_{f_i \in \mathcal{C}} e^{-[V_1(f_i) + \sum_{i' \in \mathcal{N}_i} V_2(f_i, f_{i'})]}}$$

- Equivalence between Markov Random Field and Gibbs Random Field
- An MRF is characterized by its local property (the Markovianity)
- A GRF is characterized by its global property (the Gibbs distribution)
- The Hammersley-Clifford theorem establishes the equivalence of these two types of properties
- The theorem states that  $F$  is an MRF on  $S$  with respect to  $\mathcal{N}$  if and only if  $F$  is a GRF on  $S$  with respect to  $\mathcal{N}$ .

# Markov-Gibbs Equivalence

- This theorem provides a simple way of specifying the joint probability
- *One can specify the joint probability  $P(F = f)$  by specifying the clique potential functions  $V_c(f)$  and choose appropriate potential functions for desired system behavior. It encodes the a priori knowledge or preference about interactions between labels*
- *The MRF-Gibbs distribution has been used in solving optimization problems. In optimization problems, an objective function is in the form of an energy function and is to be minimized*
- *As the quantitative cost measure, an energy function defines the minimal solution as its minimum, usually a global one*
- *Formulate an energy function so that the correct solution is embedded as the minimum*

# MRF Image Modeling: Example

- image modeled as a random field  $Y = \{Y_s : s \in L\}$
- $L = \{(i, j) | 0 \leq i \leq N_1 - 1, 0 \leq j \leq N_2 - 1\}$  index set, a set of site indices on 2-D discrete  $N_1 \times N_2$  rectangular integer lattice
- for each lattice point or pixel  $s = (i, j) \in S$   $Y_s$  is a real-valued random variable
- random field  $Y$  is characterized by a joint probability distribution  $P_Y$  which may be characterized by an associated parameter set  $\theta_Y$
- random variables  $Y_s$  will take on sample values or realizations  $y_s$  from a common finite set of integers  $\{0, 1, 2, \dots, L_Y - 1\}$

# MRF Image Modeling: Example

- let random field  $X = \{X_1, \dots, X_m\}$  be a Markov Random Field defined on  $L$  and  $m$  is the total number of the classes
- sites in  $L$  are related to each other via a neighborhood system  $\Psi = \{N_l, l \in L\}$   $N_l$  is the set of neighbor of site
- site is not a neighbor of itself  $\nexists$
- clique is a subset of sites in  $N_l$   $c \in N_l$  is a clique of every pair of distinct sites in  $c$  are neighbors
- random field  $X$  is considered to be an MRF on  $S$  if and only if  $P(X = x) > 0$  and 
$$P(X_l = x_l | X_r = x_r, r \neq s) = P(X_l = x_l | X_r = x_r, r \in N_l)$$
- difficult to determine the above characteristics in practice



# MRF Image Modeling: Example

- MRF has the form

$$P(X = x) = \frac{1}{Z} e^{-U(X)/T}$$

- $X = x$  is a realization from  $X = \{X_1, \dots, X_m\}$ , i.e.,  $x = \{x_1, \dots, x_m\}$  is a set of random field  $X$  and

$$U(X) = \sum_L^I V_c(x)$$

is global energy function and it is given by the sum of clique potentials  $V_c(x)$ , over all possible cliques

- choice of energy function is arbitrary

# MRF Image Modeling: Example

- general expression for the energy function

$$U(X) = \sum_L V_c(ci) + \sum_{c \in N_l} V_c(ci, cj)$$

- known as Potts model  $V_c(ci)$  external field that weighs the relative importance of different classes
- simplified Potts model with no external energy  $V_c(ci) = 0$
- local spatial transitions are taken into account and all the classes in the label segmentation,  $X_{opt}$ , as near as possible to the real image  $X^*$

$$V_c(ci, cj) = \begin{cases} -\beta & \text{if } ci = cj \\ 0 & \text{otherwise} \end{cases}$$

$$V_c(ci, cj) = \begin{cases} -\frac{\beta \sigma_i^2}{(\sigma_i^2 + (y_d - y_c)^2 \times d_{cd})} & \text{if } ci \neq cj \\ -\beta & \text{if } ci = cj \end{cases}$$

# MRF-MAP Segmentation

- $y_{ci}$  and  $y_{cj}$  are the pixel intensities of  $ci$  and  $cj$ ,  $d_{ci cj} = 1$  or  $\sqrt{2}$  represents the distance between the two pixels
- $\beta$  is constant that controls the classification
- decreasing of  $d_{ci cj}$  and  $(y_{ci} - y_{cj})^2$ ,  $V_c(ci, cj)$  decreased to  $-\beta$ , i.e.,  $ci$  and  $cj$  are right one pixel or their intensities are same
- image  $Y$  as rectangular lattice  $L$   $y_l$  denotes the intensity of the pixel at  $l$  and it correspond to the label  $x_l$  in  $X$
- Bayes theorem yields a complete model coupling intensities and labels

$$P(X|Y) = \frac{P(Y|X)P(X)}{P(Y)}$$



# MRF-MAP Segmentation

- the prior probability of the image  $P(Y)$  is independent of the labelling  $X$ , using MAP

$$X_{opt} = \max \arg_{X \in L} \{P(Y|X)P(X)\}$$

- assume that image data is obtained by adding an identical independently distributed (i.i.d) Gaussian noise

$$p(y_I|x_I) = \frac{1}{\sqrt{2\pi\sigma_I^2}} \exp \left\{ -\frac{(y_I - \mu_{xI})^2}{2\sigma_I^2} \right\}$$

- based on conditional independent assumption of  $Y$  the conditional density  $P(Y|X) = \prod_I P(y_I|x_I)$

## MRF-MAP Segmentation

$$P(Y|X) = \prod_L \left[ \frac{1}{\sqrt{2\pi\sigma_l^2}} \exp \left\{ -\frac{(y_l - \mu_{xl})^2}{2\sigma_l^2} \right\} \right] \propto \exp \left[ -\sum_L \frac{(y_l - \mu_{xl})^2}{2\sigma_l^2} \right]$$

- potential function of the conditional probability can be written as

$$U(Y|X) = \sum_L \frac{(y_l - \mu_{xl})^2}{2\sigma_l^2}$$

- similarly prior density of MRF takes the form of

$$P(X) = \frac{1}{Z} \exp \left[ -\sum_L \sum_{c \in N_l} \frac{V_c(c_i, c_j)}{T} \right]$$

$$P(X|Y) \propto \exp \left( -\sum_L \left[ \frac{(y_l - \mu_{xl})^2}{2\sigma^2} + \sum_{c \in N_l} \frac{V_c(c_i, c_j)}{T} \right] \right)$$

# MRF-MAP Segmentation

- image  $Y$  is then segmented by finding the field of labels  $X$

$$\begin{aligned} X_{opt} &= \min \arg_{x \in L} U(X|Y) \\ &= \min \arg \left( \sum_L \left[ \frac{(y_l - \mu_l)^2}{2\sigma^2} + \frac{1}{T} \sum_{c \in N_l} V_c(c_i, c_j) \right] \right) \end{aligned}$$

- optimization method: simulated annealing (SA), iterated conditional model (ICM) used to find the solution
- SA is slow but guarantee a global minimum solution
- ICM likely to reach local minima and no guarantee that a global minimum of energy function can be obtained, provides much faster convergence
- ICM iteratively decrease the energy by visiting and updating the pixel

# MRF-MAP Segmentation

- for each pixel  $l$ , given the observed image  $Y$  and current labels of all the pixels in the neighborhood, the label of  $X_l$  is replaced with one that can maximize the probability as

$$X_l^{(k+1)} = \arg \max P(X_l^{(k)} | Y, X_r^{(k)}, r = l)$$

- starting from the initial state, keep on running on the procedure above until either the predefined number of iterations is reached or the label of  $X$  does not change

Thank You

