

Graphical Models

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Factorization

Given a graph with nodes x_1, \dots, x_D , the corresponding joint density is given by

$$p_{\mathbf{x}}(\mathbf{x}) = \frac{1}{Z} \prod_C \psi_C(\mathbf{x}_C),$$

where $\mathbf{x} = (x_1, \dots, x_D)^T$ and ψ_C are functions over the maximal cliques of the graph, called *potential functions*.

Potential Functions

- What are potential functions? They are functions

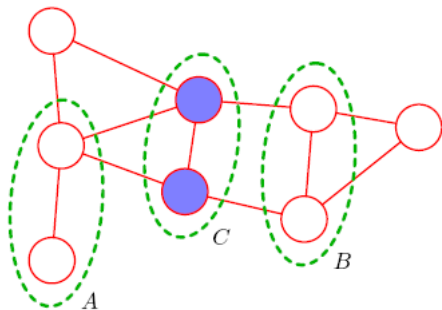
$$\psi_C : \mathbb{R}^{|C|} \rightarrow \mathbb{R},$$

where $|C|$ is the number of nodes in the maximal clique C .

- They represent some relationship between adjacent nodes (random variables) in the graph.
- They need not have a probabilistic interpretation. In image de-noising example, potential functions represented correlation between nodes of (two-node) cliques.

Conditional Independence

In a MRF, conditional independence is simpler: if all paths between node a and b pass through c , then $a \perp\!\!\!\perp b \mid c$. For example, in the following graph, it is true that $A \perp\!\!\!\perp B \mid C$.



Directed \rightarrow undirected

For every node x

- Add links between all parent nodes
- Drop arrows
- Set

$$\psi_{x, \text{pa}_x}(x, \text{pa}_x) = p(x | \text{pa}_x) \prod_{y \in \text{pa}_x} p(y)$$

Note that this process loses information about independence: in the directed graph, the parents of x were independent of each other and, in the undirected graph, they are not.

Thoughts on undirected graphs

It seems (to me) that – because of the “fuzziness” of the potential functions – there are two ways of getting an undirected graph:

- (i) generate it from a directed graph.
- (ii) generate it from a model that has a natural graphical representation (e.g. the image de-noising).

In contrast, directed graphs have a very precise probabilistic interpretation and can always be generated from models.

Using independence

Let x , y and z be discrete random variables that take on five values.

- In general, $p(x, y, z)$ is a $5 \times 5 \times 5$ table of values
- If x , y and z are independent, $p(x, y, z) = p(x)p(y)p(z)$; the joint is determined by the marginals

This basic principle will be used to speed up calculations of marginals and conditional marginals (like posteriors).

Chains

A chain is a graph of the form



and its joint distribution is given by

$$p(\mathbf{x}) = \psi_{1,2}(x_1, x_2) \cdots \psi_{N-1,N}(x_{N-1}, x_N),$$

where $\psi_{1,2}(x_1, x_2) = p(x_1)p(x_2|x_1)$ and

$$\psi_{k-1,k}(x_{k-1}, x_k) = p(x_k|x_{k-1}), \quad \text{for } k = 3, \dots, N.$$

Inference on chains

Given a chain of discrete random variables (with K possible values for each), compute $p(x_n)$, given by

$$\begin{aligned} p(x_n) &= \sum_{x_1} \cdots \sum_{x_{n-1}} \sum_{x_{n+1}} \cdots \sum_{x_N} p(\mathbf{x}) \\ &= \sum_{x_1} \cdots \sum_{x_{n-1}} \sum_{x_{n+1}} \cdots \sum_{x_N} \psi(x_1, x_2) \cdots \psi(x_{N-1}, x_N). \end{aligned}$$

Each x_i can take on K possible values; sum over K^{N-1} values!

We can take advantage of independence properties ($x_{k+1} \perp\!\!\!\perp x_{k-1} | x_k$) to make this computation more efficient.

Inference on chains

Consider the following:

$$\begin{aligned} p(x_n) &= \sum_{x_1} \cdots \sum_{x_{n-1}} \sum_{x_{n+1}} \cdots \sum_{x_N} p(\mathbf{x}) \\ &= \sum_{x_1} \cdots \sum_{x_{n-1}} \sum_{x_{n+1}} \cdots \sum_{x_{N-1}} \left[\psi(x_1, x_2) \cdots \psi(x_{N-1}, x_N^1) \right. \\ &\quad \left. + \psi(x_1, x_2) \cdots \psi(x_{N-1}, x_N^2) \right. \\ &\quad \vdots \\ &\quad \left. + \psi(x_1, x_2) \cdots \psi(x_{N-1}, x_N^K) \right] \\ &= \sum_{x_1} \cdots \sum_{x_{n-1}} \sum_{x_{n+1}} \cdots \sum_{x_{N-1}} \left[\psi(x_1, x_2) \cdots \psi(x_{N-2}, x_{N-1}) \sum_{x_N} \psi(x_{N-1}, x_N) \right] \end{aligned}$$

Inference on chains

Applying this “trick” repeatedly, and realizing that we can separate into two parts, we get that

$$\rho(x_n) = \left[\sum_{x_{n-1}} \psi(x_{n-1}, x_n) \cdots \left[\sum_{x_1} \psi(x_1, x_2) \right] \cdots \right] \\ \left[\sum_{x_{n+1}} \psi(x_n, x_{n+1}) \cdots \left[\sum_{x_N} \psi(x_{N-1}, x_N) \right] \cdots \right].$$

Computation reduced to $O(NK^2)$.

Factor Graphs

Express joint density using “factors”:

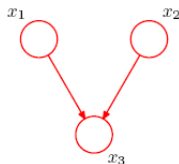
$$p(\mathbf{x}) = \prod_s f_s(\mathbf{x}_s)$$

Factor graphs:

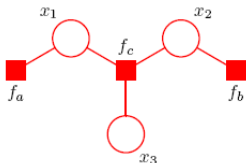
- Square nodes for “factors”
- Links between factor nodes and variable nodes used in factor
- Can convert between directed/undirected graphs to factor graphs

Example

Consider $p(\mathbf{x}) = p(x_1)p(x_2)p(x_3|x_1, x_2)$, with directed graph



Let $f_a(x_1) = p(x_1)$, $f_b(x_2) = p(x_2)$, $f_c(x_1, x_2, x_3) = p(x_3|x_1, x_2)$;
 the corresponding factor graph is



Summary

- Undirected graphs give a lot of flexibility; but not always applicable.
- Conditional independence is important! Taking advantage of it speeds up computation.
- Discuss!!