Probabilistic graphical models
Directed (BNs) and undirected (MRFs) graphs

22-808: Generative models
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Fatemeh Seyyedsalehi

Probabilistic graphical models

▸ A framework to tackle with complex joint distributions

- ▸ Representation
  - ▸ Directed graphs: Bayesian network
  - ▸ Undirected graphs: Markov random fields
- ▸ Learning
- ▸ Inference

▸ This lecture

- ▸ Representation in PGMs

Probabilistic graphical models

▸ Searching in the fully generalized space of distributions even in a simple probabilistic problem is impossible!

▸ Learn an effective and general technique for parameterizing probability distributions using only a few parameters.

Probabilistic graphical models

- Independencies assumptions are useful
  - Simplify representation and alleviate inference complexities

- Enable us to incorporate domain knowledge and structures
  - Modular combination of heterogeneous parts
  - Combining data and knowledge (Bayesian philosophy)

Bayesian networks

▶ Directed graphical models are tools to present family of probability distributions that can be naturally described using a directed acyclic graph.

  ▶ Nodes as random variables

  ▶ Edges as dependencies

▶ The intention behind these parameterization is chain rule!

$$p(x_1, x_2, \ldots, x_n) = p(x_1)p(x_2 \mid x_1) \cdots p(x_n \mid x_{n-1}, \ldots, x_2, x_1)$$

Bayesian networks

▶ Bayesian networks represent a joint distribution in terms of **the graph structure** and **conditional probability distributions (CPD)**

$$G = (V, E)$$

- A random variable x_i for each node $i \in V$.

- One conditional probability distribution (CPD) $p(x_i \mid x_{A_i})$ per node, specifying the probability of x_i conditioned on its parents' values.
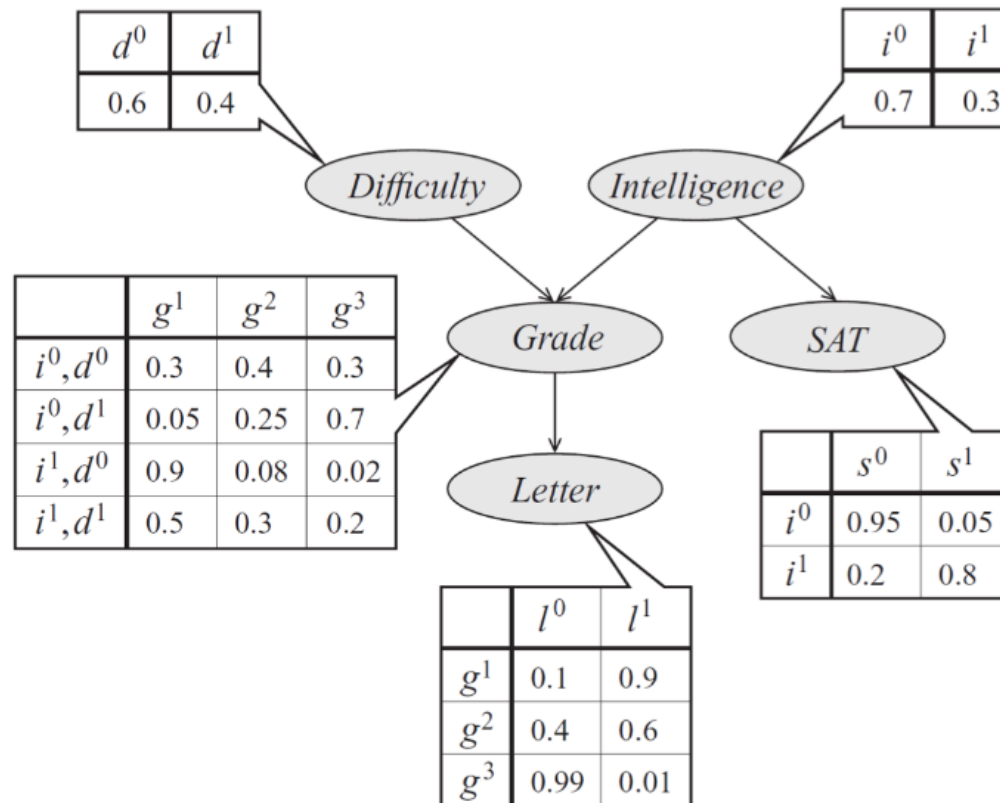
$$P(I, D, G, L, S) = P(I)P(D)P(G \mid I, D)P(L \mid G)P(S \mid I).$$
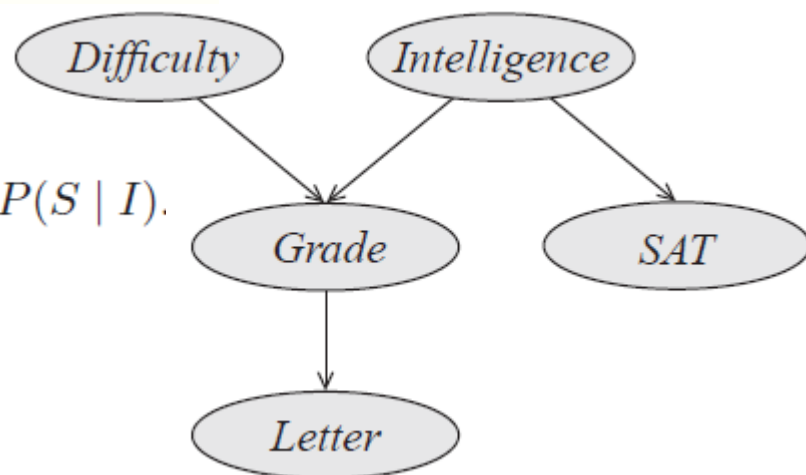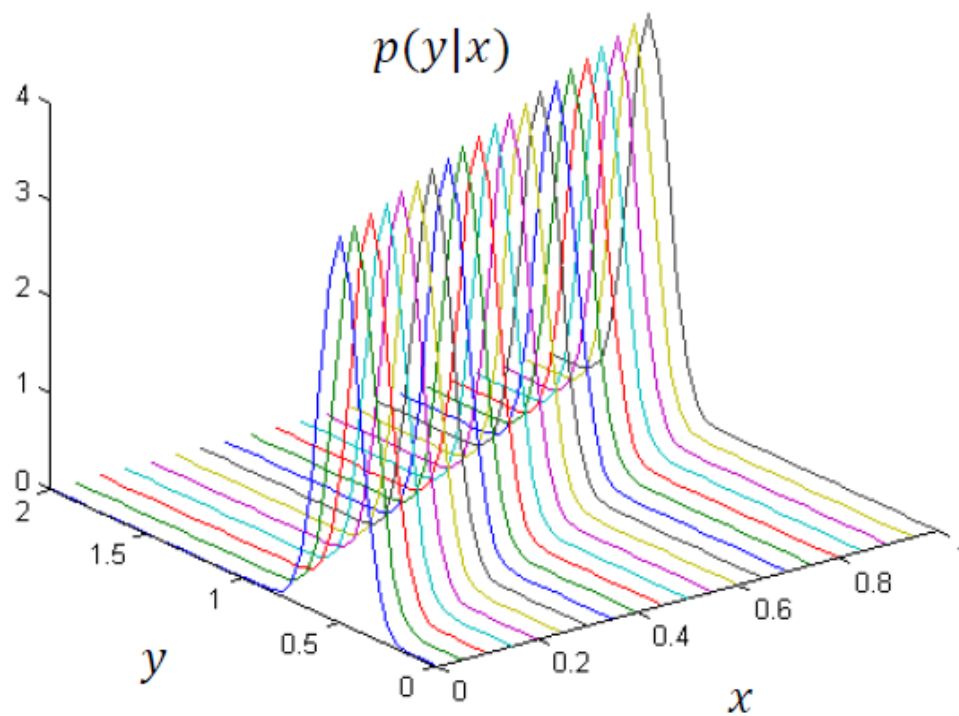
Bayesian networks
Discrete example

▸ When the variables are discrete, we may think of the factors (CPDs) as *probability tables*, in which rows correspond to assignments to parents and columns correspond to values of the node.
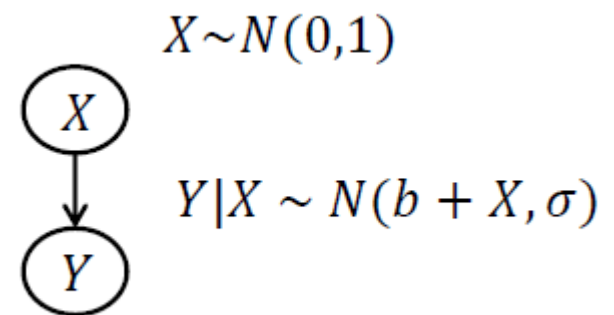
d^0	d^1
0.6	0.4

i^0	i^1
0.7	0.3

	g^1	g^2	g^3
i^0,d^0	0.3	0.4	0.3
i^0,d^1	0.05	0.25	0.7
i^1,d^0	0.9	0.08	0.02
i^1,d^1	0.5	0.3	0.2

	s^0	s^1
i^0	0.95	0.05
i^1	0.2	0.8

	l^0	l^1
g^1	0.1	0.9
g^2	0.4	0.6
g^3	0.99	0.01

Difficulty *Intelligence* *Grade* *SAT* *Letter*

Bayesian networks
Continues example

$$p(y|x)$$

$$X \sim N(0,1)$$

$$Y|X \sim N(b + X, \sigma)$$

$$b = 0.5$$
$$\sigma = 0.1$$

Bayesian networks

▶ A probability distribution is factorized over a *DAG G* if it can be decomposed into a product of factors specified by G.

▶ A Bayesian network represent distributions via products of smaller, local conditional probability distributions.

  ▶ Introduces independency assumptions over variables

▶ $I(p)$: denote the set of all independencies that hold for a joint distribution p.

  ▶ $p(x, y) = p(x)p(y) \rightarrow x \perp y \in I(p)$

9

Bayesian networks

▸ Let G be a graph over $x_1, x_2, \ldots, x_n$ distribution p factorizes over G if:

$$p(x_1, x_2, \ldots, x_n) = \prod_{i=1}^{n} p(x_i | pa(x_i))$$

▸ $pa(.)$: parents of a node

▸ Factorization $\Leftrightarrow$ Independence
  ▸ If p factorizes over G, then any variable in p is independent of its non-descendants given its parents (in G)
  ▸ If any variable in the distribution p is independent of its non-descendants given its parents (in the graph G) then p factorizes over G

Independencies described by directed graphs

- *Common parent*. If G is of the form $X \leftarrow Z \rightarrow Y$, and Z is observed, then $X \perp Y \mid Z$. However, if Z is unobserved, then $X \not\perp Y$. Intuitively this stems from the fact that Z contains all the information that determines the outcomes of X and Y; once it is observed, there is nothing else that affects these variables' outcomes.

- *Cascade*: If G equals $X \rightarrow Z \rightarrow Y$, and Z is again observed, then, again $X \perp Y \mid Z$. However, if Z is unobserved, then $X \not\perp Y$. Here, the intuition is again that Z holds all the information that determines the outcome of Y; thus, it does not matter what value X takes.

- *V-structure* (also known as *explaining away*): If G is $X \rightarrow Z \leftarrow Y$, then knowing Z couples X and Y. In other words, $X \perp Y$ if Z is unobserved, but $X \not\perp Y \mid Z$ if Z is observed.
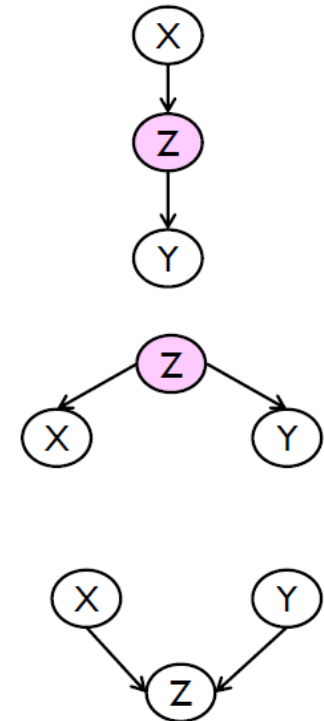
Independencies described by directed graphs
D-separation

- Considering three disjoint sets of nodes:
  - *A, B, C*

- *A* is **d-separated** from *B* by *C* if all paths between *A* and *B* are blocked by *C*
  - There is no **active path** between A and B

- A is d-separated from B by C if $\boldsymbol{A} \perp \boldsymbol{B} | \boldsymbol{C}$

Path blocking

▶ Head to tail during path

$X \in A$ $Z \in C$ $Y \in B$
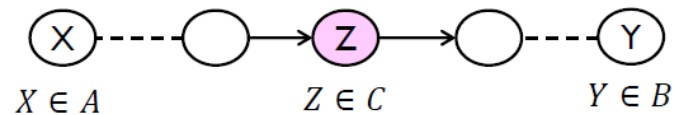
▶ Tail to tail during path

$X \in A$ $Z \in C$ $Y \in B$
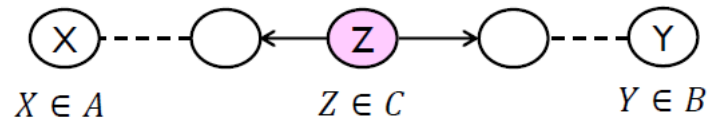
▶ Head to head, visiting a v-structure

   ▶ Z and none of its descendants are observed

$X \in A$ $Y \in B$
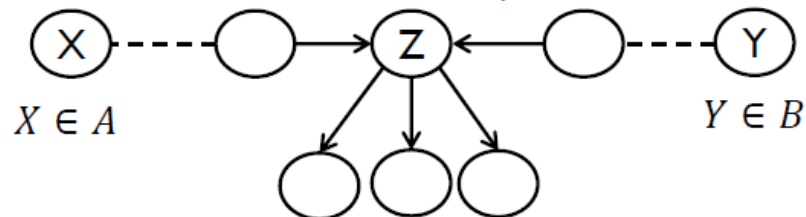
Independencies described by directed graphs

For example, in the graph below, X_1 and X_6 are d-separated given X_2, X_3. However, X_2, X_3 are not d-separated given X_1, X_6, because we can find an active path (X_2, X_6, X_5, X_3)
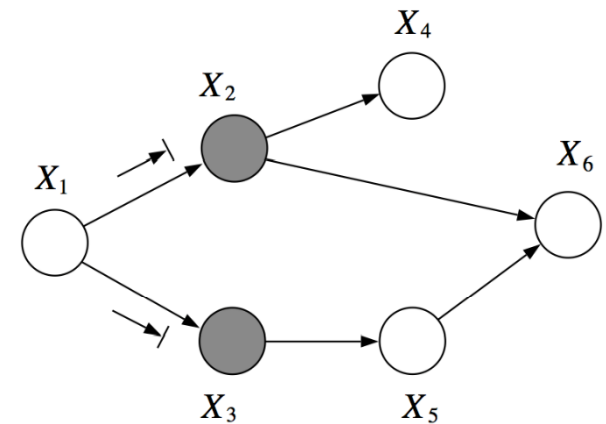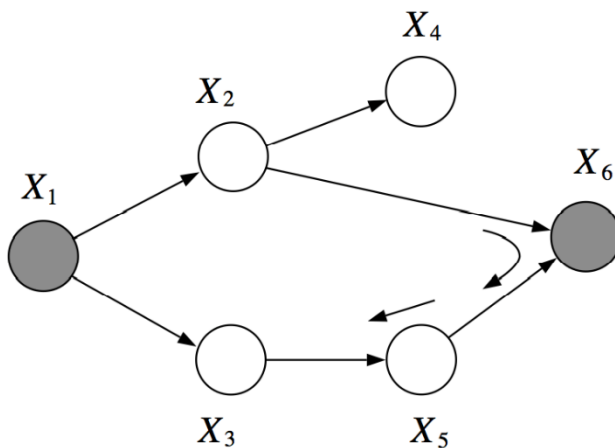
A simple d-separation simulator ☺

Markov blanket of a node

▶ A variable is conditionally independent of all other variables given its **Markov blanket**

▶ Markov blanket if a set *A* is *U* when:
  ▶ The minimal set of nodes such that *A* is independent from the rest of the graph if *U* is observed

▶ Markov blanket of a node:
  ▶ All parents
  ▶ All children
  ▶ Co-parents of children
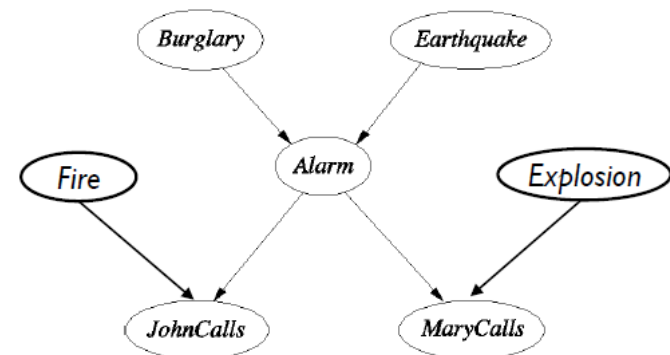
Markov random networks

▶ Undirected graphs for representation of joint distributions

  ▶ Unlike in the directed case, we are not saying anything about how one variable is generated from another set of variables (as a conditional probability distribution would do).

$$\tilde{p}(A, B, C, D) = \phi(A, B)\phi(B, C)\phi(C, D)\phi(D, A)$$

$$\phi(X, Y) = \begin{cases} 10 & \text{if } X = Y = 1 \\ 5 & \text{if } X = Y = 0 \\ 1 & \text{otherwise.} \end{cases}$$

$$p(A, B, C, D) = \frac{1}{Z}\tilde{p}(A, B, C, D)$$
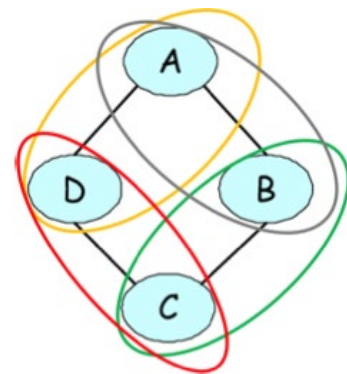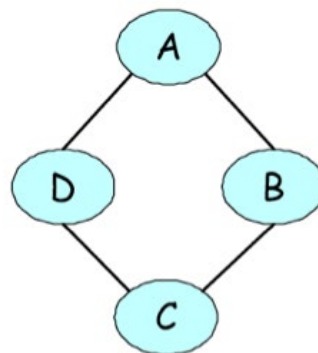
Markov random networks

▸ They specify dependent variables (but no causality relations) and define the strength of their interactions.

▸ This defines an energy landscape over the space of possible assignments and we convert this energy to a probability via the normalization constant.
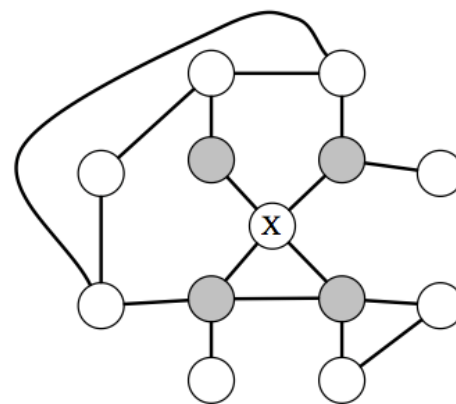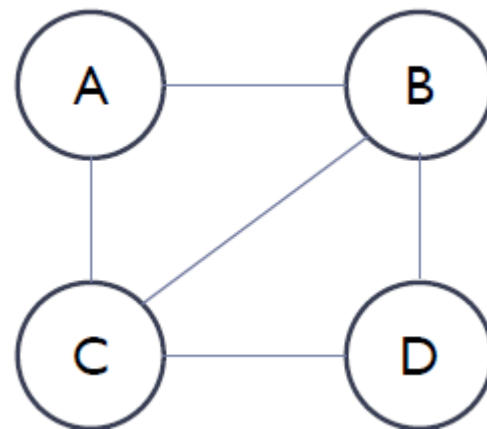
MRF factorization

▶ **Clique**: subsets of nodes in the graph that are fully connected (complete subgraph)

▶ **Maximal clique**: no superset of the nodes in a clique are also compose a clique

▶ Factors are functions of the variables in cliques

  ▶ To reduce the number of factors we allow factors only for maximal cliques

Cliques: {A,B,C}, {B,C,D}, {A,B}, {A,C}, {B,C}, {B,D}, {C,D}, {A}, {B}, {C}, {D}

Max-cliques: {A,B,C}, {B,C,D}

MRF factorization

▸ A distribution $p(.)$ is factorized over an MRF G if it can be parameterized as follows,

$$p(x_1, x_2, \ldots, x_n) = \frac{1}{Z} \prod_{i=1}^{k} \phi_i(D_i)$$

$$Z = \sum_{X} \prod_{i=1}^{k} \phi_i(D_i)$$

where each D_i is a **complete subgraph** of G

▸ When there is no direct edge between two nodes, x_i and x_j, there exist at least the following conditional independency between them:

$$x_i \perp x_j \mid X/\{x_i, x_j\}$$

  ▸ To hold this independency in $p(.)$, these two variables are not appeared in the domain of a same factor

MRF factorization

▶ Potential functions:
  ▶ The function over each clique (factor)

▶ Potential functions and cliques in the graph completely determine the joint distribution.

▶ Potentials are not necessarily marginal or conditional distributions

Markov random networks

▶ Formal definition

A Markov Random Field (MRF) is a probability distribution p over variables $x_1, \ldots, x_n$ defined by an *undirected* graph G in which nodes correspond to variables x_i. The probability p has the form

$$p(x_1, \ldots, x_n) = \frac{1}{Z} \prod_{c \in C} \phi_c(x_c),$$

where C denotes the set of *cliques* (i.e., fully connected subgraphs) of G, and each *factor* ϕ_c is a non-negative function over the variables in a clique. The *partition function*

$$Z = \sum_{x_1, \ldots, x_n} \prod_{c \in C} \phi_c(x_c)$$

Independencies in MRFs

▸ A simple rule:

  ▸ Variables x and y are dependent if they are connected by a path of unobserved variables.

▸ Markov blanket in MRFs:

  ▸ In both BNs and MRFs

  ▸ In MRFs: simply all neighbors of a node
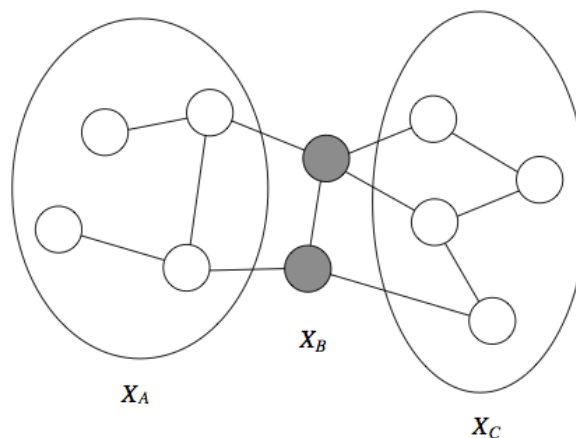
MRF example:
Image denoising

- Pixels are noisy observed variables: y_i
- We assume the noise free image as a latent behind the observed pixels: x_i
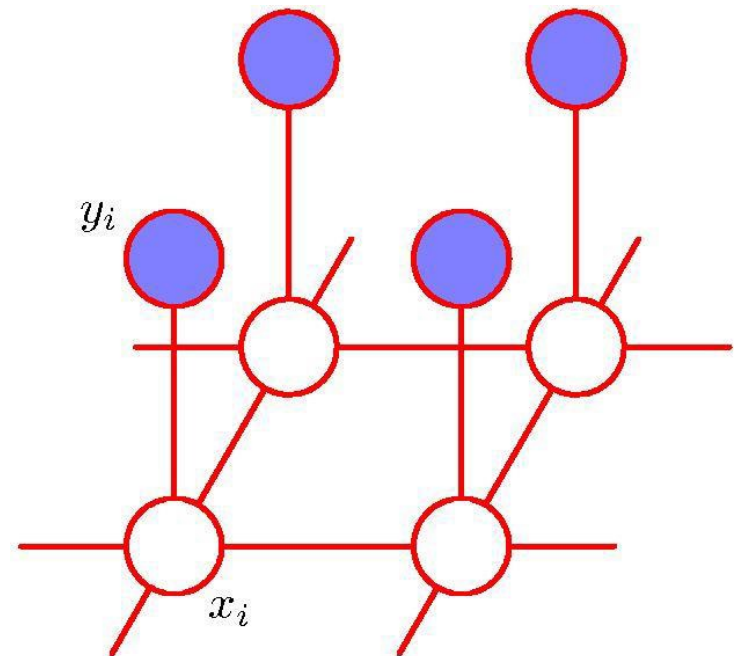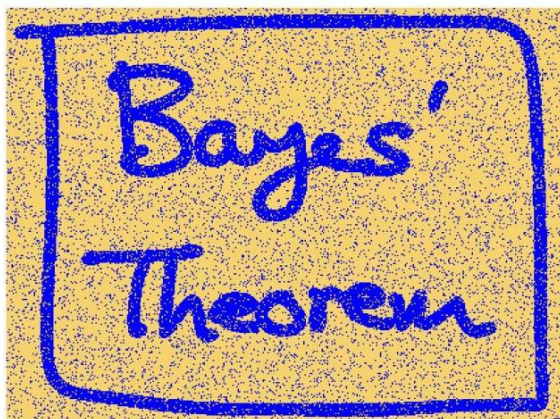
MRFs compared to BNs

▸ Pros.
  ▸ They can be applied to a wider range of problems in which there is no natural directionality associated with variable dependencies.
  ▸ Undirected graphs can succinctly express certain dependencies that Bayesian nets cannot easily describe (although the converse is also true)

▸ Cons.
  ▸ Computing the normalization constant Z requires summing over a potentially exponential number of assignments.
    ▸ NP-hard; many undirected models will be intractable and will require approximation techniques.
  ▸ Difficult to interpret.
  ▸ It is much easier to generate data from a Bayesian network

Hybrid graphs

▶ Partially directed acyclic graphs

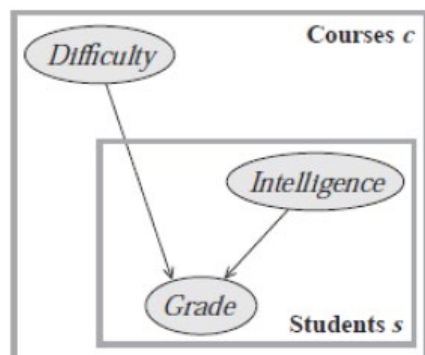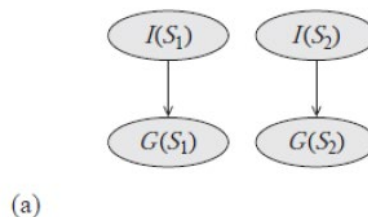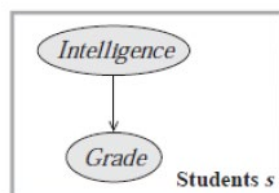 ▶ A combination of both directed and undirected graphs
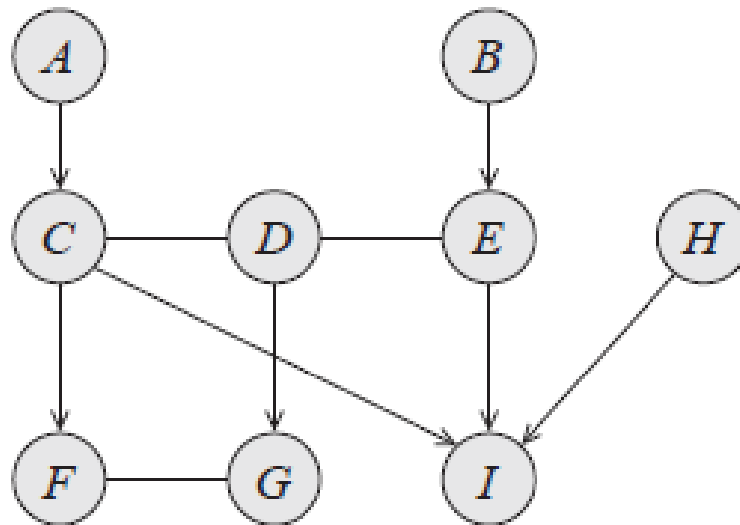
Plate notation

▶ Plate notation is a rectangle in graphical model representation which shows random variables generated from the same distribution

▶ Plate notation present a replication of random variables that share same parameters
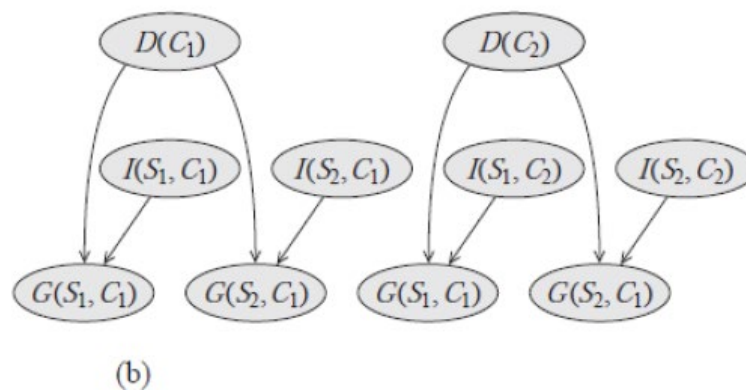
Generative vs. discriminative models

- In generative models we describe the generation process of observed variables

- In discriminative models, we learn how samples are discriminated
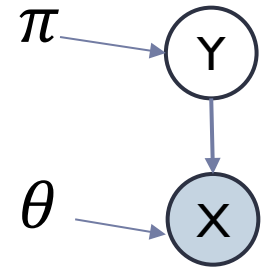
  - Decision boundaries in classifiers
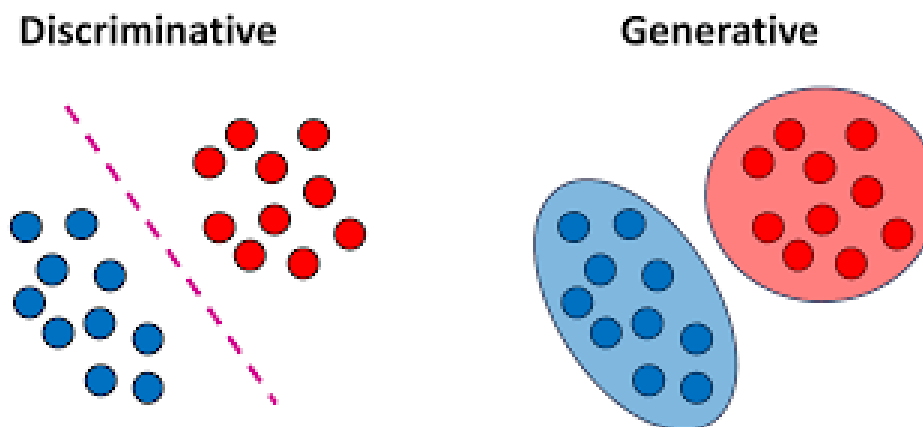
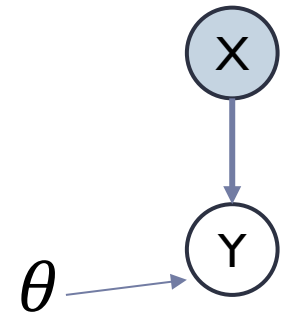# Generative vs. discriminative models Example

- Generative classifier
  - We should learn $p(y)$, $p(x|y)$

- Discriminative classifier
  - We should learn $p(x)$, $p(y|x)$
  - However, for classification task

$p(y|x)$ is the only thing we need.

  - Less parameters are needed to be learned
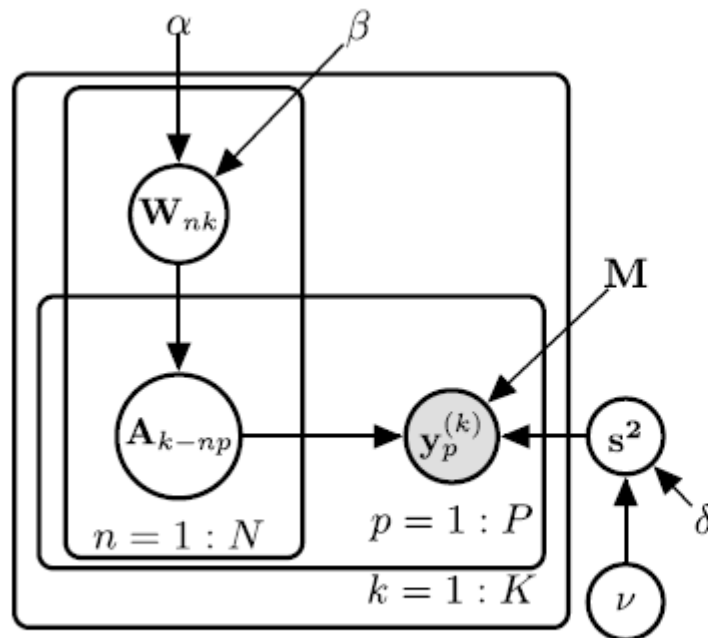
- When we only need to discriminate
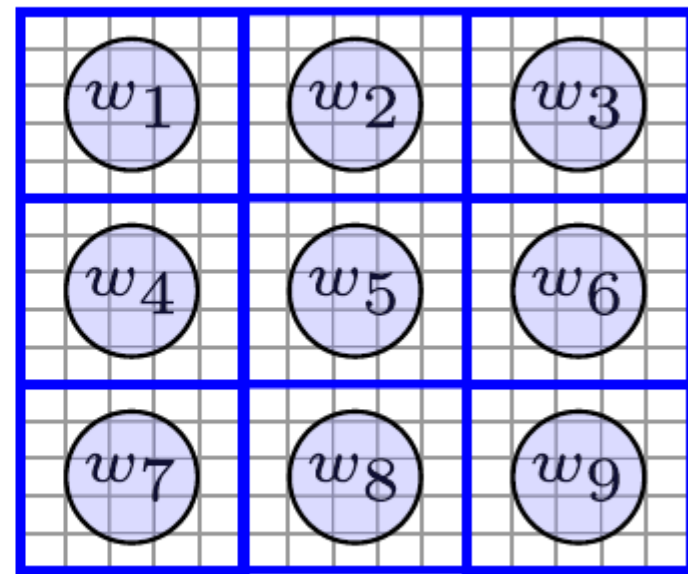
Between samples, discriminative models are preferred.

Generative PGM example
Hyperspectral unmixing with PGMs

▸ A generative model

  ▸ K = number of patches

  ▸ P = number of pixels in each patch

  ▸ N = the dimension of vector A

(a)

(b)

Next topic

- Probabilistic graphical models
  - Exact and approximate inference