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# Machine Learning

## Chapter 8: Graphical Models

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## Introduction

- Visualize the structure of a probabilistic model
- Design and motivate new models
- Insights into the model's properties, in particular **conditional independence** obtained by inspection
- Complex computations = graphical manipulations

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## Terminology

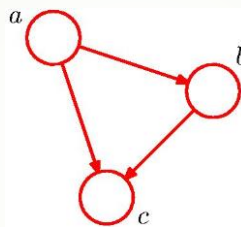
- Nodes (vertices) + links (arcs, edges)
- Node: a random variable
- Link: a probabilistic relationship
- **Directed graphical models** or **Bayesian networks** useful to express causal relationships between variables.
- **Undirected graphical models** or **Markov random fields** useful to express soft constraints between variables.
- **Factor graphs** convenient for solving inference problems

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## Bayesian Networks

- Directed Acyclic Graph (DAG)



$$p(a, b, c) = p(c|a, b)p(a, b) = p(c|a, b)p(b|a)p(a)$$

Generalization to  $K$  variables:

$$p(x_1, \dots, x_K) = p(x_K|x_1, \dots, x_{K-1}) \dots p(x_2|x_1)p(x_1)$$

- The associated graph is fully connected.
- The absence of links conveys important information.

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## Bayesian Networks

$$p(x_1, \dots, x_7) = p(x_1)p(x_2)p(x_3)p(x_4|x_1, x_2, x_3) \\ p(x_5|x_1, x_3)p(x_6|x_4)p(x_7|x_4, x_5)$$

is not a fully connected graph

General Factorization

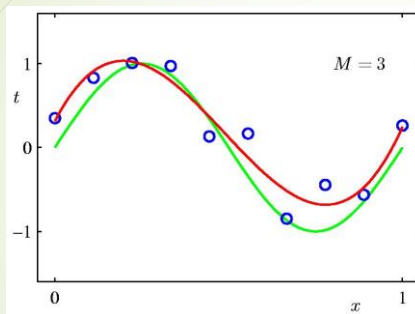
$$p(\mathbf{x}) = \prod_{k=1}^K p(x_k | \text{pa}_k)$$

- This key equation expresses the factorization properties of the joint distribution.
- There must be no directed cycles
- These graphs are also called **directed acyclic graphs (DAGs)**.

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## Bayesian Curve Fitting (1)



Polynomial

$$y(x, \mathbf{w}) = \sum_{j=0}^M w_j x^j$$

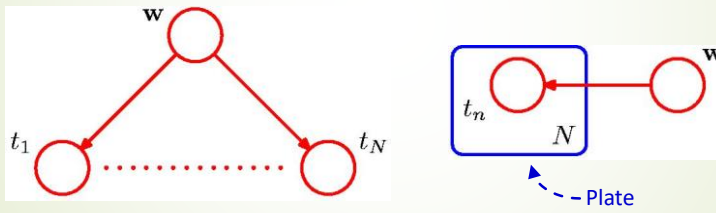
$$p(\mathbf{t}, \mathbf{w}) = p(\mathbf{w}) \prod_{n=1}^N p(t_n | y(\mathbf{w}, x_n))$$

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## Bayesian Curve Fitting (2)

$$p(\mathbf{t}, \mathbf{w}) = p(\mathbf{w}) \prod_{n=1}^N p(t_n | y(\mathbf{w}, x_n))$$



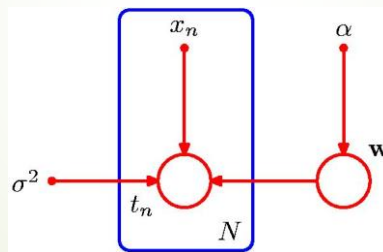
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## Bayesian Curve Fitting (3)

- Input variables and explicit hyperparameters

$$p(\mathbf{t}, \mathbf{w} | \mathbf{x}, \alpha, \sigma^2) = p(\mathbf{w} | \alpha) \prod_{n=1}^N p(t_n | \mathbf{w}, x_n, \sigma^2).$$



Deterministic parameters shown by small node

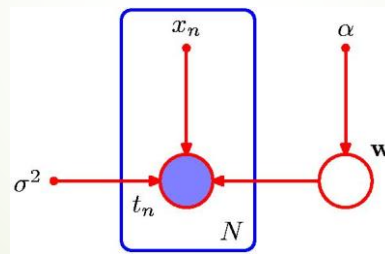
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## Bayesian Curve Fitting —Learning

► Condition on data

$$p(\mathbf{w}|\mathbf{t}) \propto p(\mathbf{w}) \prod_{n=1}^N p(t_n|\mathbf{w})$$



Latent (hidden) variable

Shaded nodes are set to observed values

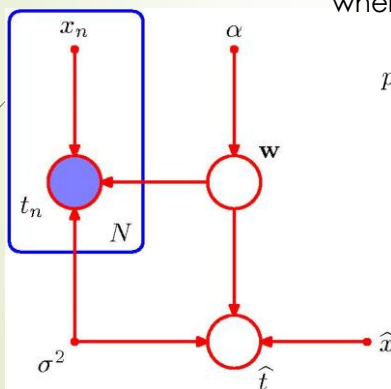
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## Bayesian Curve Fitting —Prediction

Predictive distribution:  $p(\hat{t}|\hat{x}, \mathbf{x}, \mathbf{t}, \alpha, \sigma^2) \propto \int p(\hat{t}, \mathbf{t}, \mathbf{w}|\hat{x}, \mathbf{x}, \alpha, \sigma^2) d\mathbf{w}$

where



$$p(\hat{t}, \mathbf{t}, \mathbf{w}|\hat{x}, \mathbf{x}, \alpha, \sigma^2) = \left[ \prod_{n=1}^N p(t_n|x_n, \mathbf{w}, \sigma^2) \right] p(\mathbf{w}|\alpha) p(\hat{t}|\hat{x}, \mathbf{w}, \sigma^2)$$

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## Generative Models

- Back to:

$$p(\mathbf{x}) = \prod_{k=1}^K p(x_k | \text{pa}_k)$$

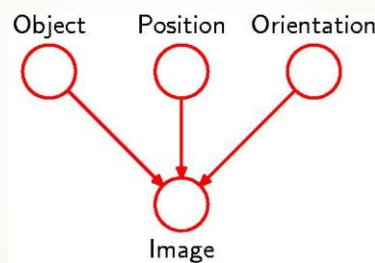
- Each node has a higher number than any of its parents.
- The factorization above corresponds to a DAG.
- Goal: draw a sample  $\hat{x}_1, \dots, \hat{x}_K$  from the joint distribution.
- Apply **ancestral sampling** starting from lower-numbered nodes, downwards through the graph's nodes.
- Generative graphical model captures the causal process that generated the observed data (object recognition example)

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## Generative Models

- Causal process for generating images



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## Discrete Variables (1)

- General joint distribution:  $K^2 - 1$  parameters



$$p(\mathbf{x}_1, \mathbf{x}_2 | \boldsymbol{\mu}) = \prod_{k=1}^K \prod_{l=1}^K \mu_{kl}^{x_{1k} x_{2l}}$$

- Independent joint distribution:  $2(K - 1)$  parameters



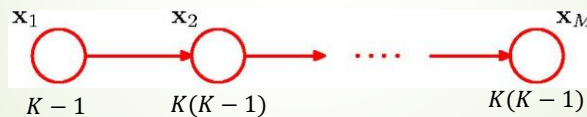
$$\hat{p}(\mathbf{x}_1, \mathbf{x}_2 | \boldsymbol{\mu}) = \prod_{k=1}^K \mu_{1k}^{x_{1k}} \prod_{l=1}^K \mu_{2l}^{x_{2l}}$$

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## Discrete Variables (2)

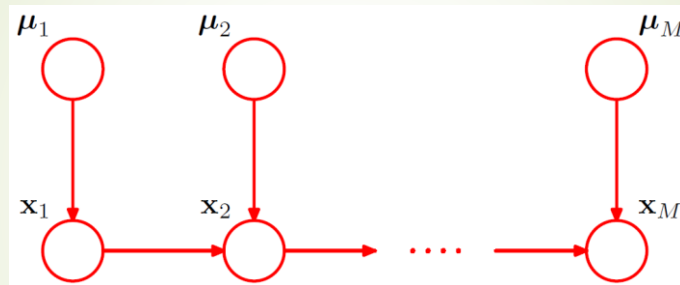
- General joint distribution over  $M$  variables:  $K^M - 1$  parameters
- $M$ -node Markov chain:  $K - 1 + (M - 1) K(K - 1)$  parameters



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## Discrete Variables: Bayesian Parameters (1)



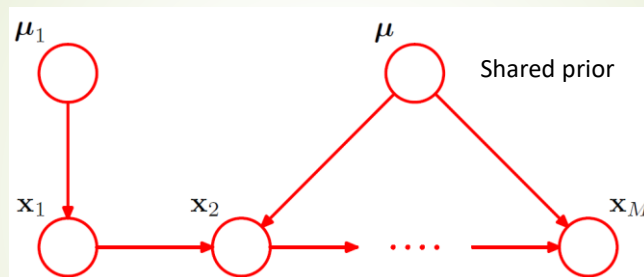
$$p(\{\mathbf{x}_m, \boldsymbol{\mu}_m\}) = p(\mathbf{x}_1 | \boldsymbol{\mu}_1) p(\boldsymbol{\mu}_1) \prod_{m=2}^M p(\mathbf{x}_m | \mathbf{x}_{m-1}, \boldsymbol{\mu}_m) p(\boldsymbol{\mu}_m)$$

$$p(\boldsymbol{\mu}_m) = \text{Dir}(\boldsymbol{\mu}_m | \boldsymbol{\alpha}_m)$$

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## Discrete Variables: Bayesian Parameters (2)



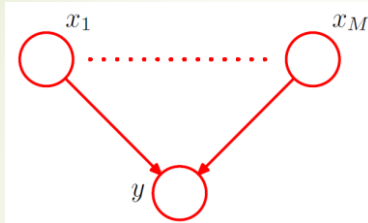
$$p(\{\mathbf{x}_m\}, \boldsymbol{\mu}_1, \boldsymbol{\mu}) = p(\mathbf{x}_1 | \boldsymbol{\mu}_1) p(\boldsymbol{\mu}_1) \prod_{m=2}^M p(\mathbf{x}_m | \mathbf{x}_{m-1}, \boldsymbol{\mu}) p(\boldsymbol{\mu})$$

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## Parameterized Conditional Distributions



If  $x_1, \dots, x_M$  are discrete,  $K$ -state variables,  $p(y = 1|x_1, \dots, x_M)$  in general has  $O(K^M)$  parameters.

The parameterized form

$$p(y = 1|x_1, \dots, x_M) = \sigma \left( w_0 + \sum_{i=1}^M w_i x_i \right) = \sigma(\mathbf{w}^T \mathbf{x})$$

requires only  $M + 1$  parameters

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## Linear-Gaussian Models

- Directed acyclic graph over  $D$  variables

$$p(x_i | \text{pa}_i) = \mathcal{N} \left( x_i \mid \sum_{j \in \text{pa}_i} w_{ij} x_j + b_i, v_i \right)$$

Each node is Gaussian, the mean is a linear function of the parents.

$$\begin{aligned} \ln p(\mathbf{x}) &= \sum_{i=1}^D \ln p(x_i | \text{pa}_i) \\ &= - \sum_{i=1}^D \frac{1}{2v_i} \left( x_i - \sum_{j \in \text{pa}_i} w_{ij} x_j - b_i \right)^2 + \text{const} \end{aligned}$$

$$x_i = \sum_{j \in \text{pa}_i} w_{ij} x_j + b_i + \sqrt{v_i} \epsilon_i$$

where  $\epsilon_i$  is a zero mean, unit variance Gaussian random variable

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## Linear-Gaussian Models

- Directed acyclic graph over  $D$  variables

$$p(x_i | \text{pa}_i) = \mathcal{N} \left( x_i \mid \sum_{j \in \text{pa}_i} w_{ij} x_j + b_i, v_i \right)$$

Each node is Gaussian, the mean is a linear function of the parents.

- Vector-valued Gaussian Nodes

$$p(\mathbf{x}_i | \text{pa}_i) = \mathcal{N} \left( \mathbf{x}_i \mid \sum_{j \in \text{pa}_i} \mathbf{W}_{ij} \mathbf{x}_j + \mathbf{b}_i, \Sigma_i \right)$$

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## Conditional Independence

- $a$  is **conditionally independent** of  $b$  given  $c$

$$p(a|b, c) = p(a|c)$$

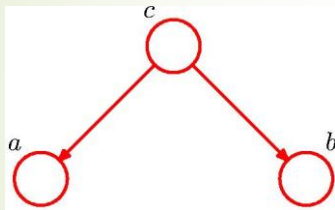
- Equivalently  $p(a, b|c) = \frac{p(a|b, c)p(b|c)}{p(a|c)p(b|c)}$

- Notation  $a \perp\!\!\!\perp b \mid c$

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## Conditional Independence: Example 1



$$p(a, b, c) = p(a|c)p(b|c)p(c)$$

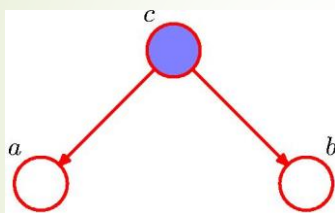
$$p(a, b) = \sum_c p(a|c)p(b|c)p(c)$$

$$a \not\perp b \mid \emptyset$$

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## Conditional Independence: Example 1



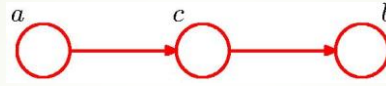
$$\begin{aligned} p(a, b|c) &= \frac{p(a, b, c)}{p(c)} \\ &= p(a|c)p(b|c) \end{aligned}$$

$$a \perp b \mid c$$

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## Conditional Independence: Example 2



$$p(a, b, c) = p(a)p(c|a)p(b|c)$$

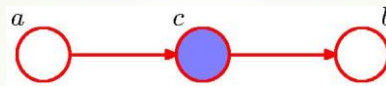
$$p(a, b) = p(a) \sum_c p(c|a)p(b|c) = p(a)p(b|a)$$

$$a \not\perp\!\!\!\perp b \mid \emptyset$$

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## Conditional Independence: Example 2



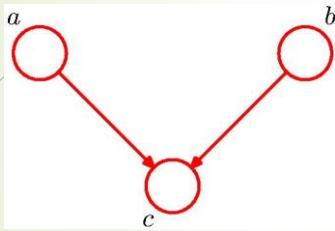
$$\begin{aligned} p(a, b|c) &= \frac{p(a, b, c)}{p(c)} \\ &= \frac{p(a)p(c|a)p(b|c)}{p(c)} \\ &= p(a|c)p(b|c) \end{aligned}$$

$$a \perp\!\!\!\perp b \mid c$$

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## Conditional Independence: Example 3



$$p(a, b, c) = p(a)p(b)p(c|a, b)$$

$$p(a, b) = p(a)p(b)$$

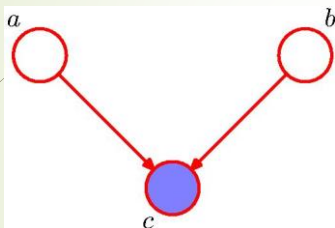
$$a \perp\!\!\!\perp b \mid \emptyset$$

Note: this is the opposite of Example 1, with  $c$  unobserved.

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## Conditional Independence: Example 3



$$\begin{aligned} p(a, b|c) &= \frac{p(a, b, c)}{p(c)} \\ &= \frac{p(a)p(b)p(c|a, b)}{p(c)} \end{aligned}$$

$$a \not\perp\!\!\!\perp b \mid c$$

Note: this is the opposite of Example 1, with  $c$  observed.

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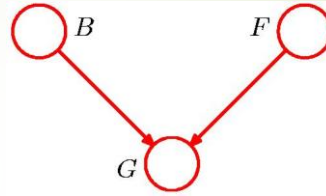
## “Am I out of fuel?”

$$p(G = 1|B = 1, F = 1) = 0.8$$

$$p(G = 1|B = 1, F = 0) = 0.2$$

$$p(G = 1|B = 0, F = 1) = 0.2$$

$$p(G = 1|B = 0, F = 0) = 0.1$$



$$p(B = 1) = 0.9$$

$$p(F = 1) = 0.9$$

and hence

$$p(F = 0) = 0.1$$

$B$  = Battery (0=flat, 1=fully charged)

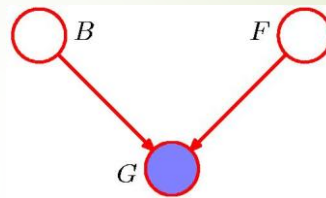
$F$  = Fuel Tank (0=empty, 1=full)

$G$  = Fuel Gauge Reading  
(0=empty, 1=full)

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## “Am I out of fuel?”



$$p(F = 0|G = 0) = \frac{p(G = 0|F = 0)p(F = 0)}{p(G = 0)} \simeq 0.257$$

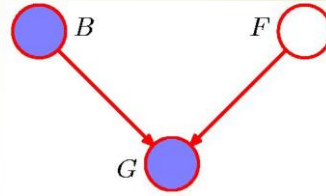
Probability of an empty tank increased by observing  $G = 0$ .

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# "Am I out of fuel?"



$$p(F = 0|G = 0, B = 0) = \frac{p(G = 0|B = 0, F = 0)p(F = 0)}{\sum_{F \in \{0,1\}} p(G = 0|B = 0, F)p(F)}$$

$$\simeq 0.111$$

Probability of an empty tank reduced by observing  $B = 0$ .  
This referred to as "explaining away".

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## D-separation

- **d-separation** (d connotes "directional") is a criterion for deciding, from a given causal graph, whether a set  $X$  of variables is independent of another set  $Y$ , given  $Z$ . The idea is to associate "dependence" with "connectedness" and "independence" with "unconnected-ness" or "separation".
- **Unconditional separation**
  - $x$  and  $y$  are d-connected if there is an unblocked path between them.
  - By "unblocked path" we mean a path that can be traced without traversing a pair of arrows that collide "head-to-head". In other words, arrows that meet head-to-head do not constitute a connection for the purpose of passing information, such a meeting will be called a "collider"

$$x \longrightarrow r \longrightarrow s \longrightarrow t \longleftarrow u \longleftarrow v \longrightarrow y$$

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## D-separation

- Unconditional separation

- $x$  and  $y$  are d-connected if there is an unblocked path between them.

$$x \longrightarrow r \longrightarrow s \longrightarrow t \longleftarrow u \longleftarrow v \longrightarrow y$$

- This graph contains one collider, at  $t$ . The path  $x-r-s-t$  is unblocked, hence  $x$  and  $t$  are d-connected. So is also the path  $t-u-v-y$ , hence  $t$  and  $y$  are d-connected. However,  $x$  and  $y$  are not d-connected (i.e., d-separated); there is no way of tracing a path from  $x$  to  $y$  without traversing the collider at  $t$ . (The ramification is that the covariance terms corresponding to these pairs of variables will be zero, for every choice of model parameters)

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## D-separation

- Blocking by conditioning

- $x$  and  $y$  are d-connected, conditioned on a set  $Z$  of nodes, if there is a collider-free path between  $x$  and  $y$  that traverses no member of  $Z$ . If no such path exists, we say that  $x$  and  $y$  are d-separated by  $Z$ . We also say then that every path between  $x$  and  $y$  is "blocked" by  $Z$ .

$$x \longrightarrow (r) \longrightarrow s \longrightarrow t \longleftarrow u \longleftarrow (v) \longrightarrow y$$

- Let  $Z = \{r, v\}$  (marked by circles).  $x$  and  $y$  are d-separated by  $Z$ , and so are  $x$  and  $s$ ,  $u$  and  $y$ ,  $s$  and  $u$  etc. The path  $x-r-s$  is blocked by  $Z$ . The only pairs of unmeasured nodes that remain d-connected in this example, conditioned on  $Z$ , are  $s$  and  $t$  and  $u$  and  $t$ .

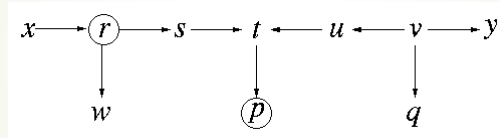
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## D-separation

- Conditioning on colliders

- If a collider is a member of the conditioning set  $Z$ , or has a descendant in  $Z$ , then it no longer blocks any path that traces this collider.



- Let  $Z = \{r, v\}$ .  $s$  and  $y$  are d-connected by  $Z$ , because the collider at  $t$  has a descendant ( $p$ ) in  $Z$ , which unblocks the path  $s-t-u-v-y$ . However,  $x$  and  $u$  are still d-separated by  $Z$ , because although the linkage at  $t$  is unblocked, the one at  $r$  is blocked since  $r$  is in  $Z$ .

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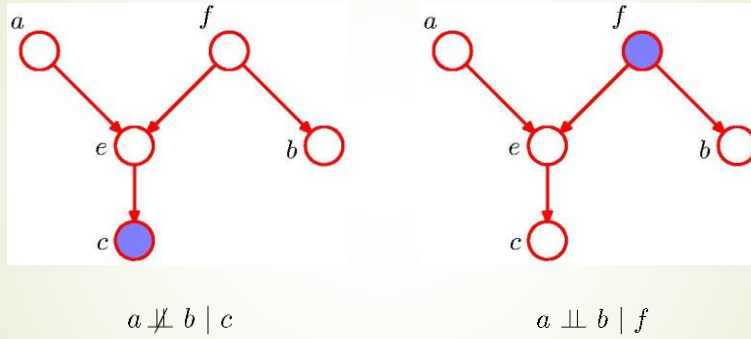
## D-separation

- $A$ ,  $B$ , and  $C$  are non-intersecting subsets of nodes in a directed graph.
- A path from  $A$  to  $B$  is blocked if it contains a node such that either
  - a) the arrows on the path meet either head-to-tail or tail-to-tail at the node, and the node is in the set  $C$ , or
  - b) the arrows meet head-to-head at the node, and neither the node, nor any of its descendants, is in the set  $C$ .
- If all paths from  $A$  to  $B$  are blocked,  $A$  is said to be d-separated from  $B$  by  $C$ .
- If  $A$  is d-separated from  $B$  by  $C$ , the joint distribution over all variables in the graph satisfies  $A \perp B | C$ .

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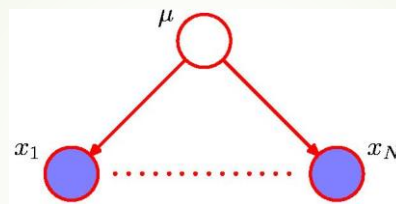
## D-separation: Example



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## D-separation: I.I.D. Data



$$p(\mathcal{D}|\mu) = \prod_{n=1}^N p(x_n|\mu)$$

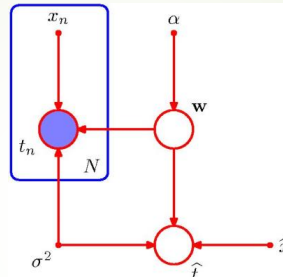
$$p(\mathcal{D}) = \int_{-\infty}^{\infty} p(\mathcal{D}|\mu)p(\mu) d\mu \neq \prod_{n=1}^N p(x_n)$$

Here  $\mu$  is a latent variable, because its value is not observed

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## D-Separation: Bayesian Polynomial Regression



- $w$  is a tail-to-tail node with respect to the path from  $\hat{t}$  to any one of the nodes  $\{t_n\}$ .
- Hence  $\hat{t} \perp t_n | w$
- Interpretation:
  - First use the training data to determine the posterior distribution over  $w$
  - Discard  $\{t_n\}$  and use posterior distribution for  $w$  to make predictions of  $\hat{t}$  for new input observations  $\hat{x}$

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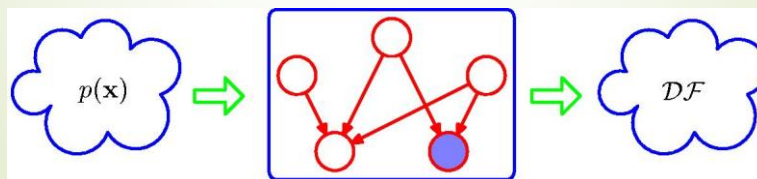
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## Interpretation as Filter

- Filter-I: allows a distribution to pass through if, and only if, it can be expressed in terms of the factorization implied by the graph

$$p(\mathbf{x}) = \prod_{k=1}^K p(x_k | \text{pa}_k)$$

- Filter-II: allows distributions to pass according to whether they respect all of the conditional independencies implied by the d-separation properties of the graph
- The set of all possible probability distributions  $p(\mathbf{x})$  that is passed by both the filters is precisely the same and are denoted by DF, for **directed factorization**



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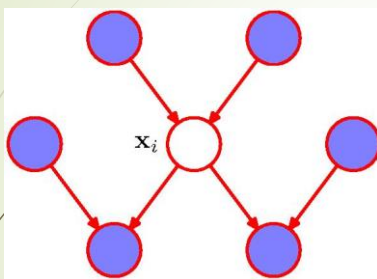
## Directed Graphs: Summary

- Represents specific decomposition of a joint probability distribution into a product of conditional probabilities
- Expresses a set of conditional independence statements through d-separation criterion
- Distributions satisfying d-separation criterion are denoted as DF
- Extreme Cases: DF can contain all possible distributions in case of fully connected graph or product of marginals in case fully disconnected graphs

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## The Markov Blanket



$$\begin{aligned}
 p(\mathbf{x}_i | \mathbf{x}_{\{j \neq i\}}) &= \frac{p(\mathbf{x}_1, \dots, \mathbf{x}_M)}{\int p(\mathbf{x}_1, \dots, \mathbf{x}_M) d\mathbf{x}_i} \\
 &= \frac{\prod_k p(\mathbf{x}_k | \text{pa}_k)}{\int \prod_k p(\mathbf{x}_k | \text{pa}_k) d\mathbf{x}_i}
 \end{aligned}$$

- Factors independent of  $\mathbf{x}_i$  cancel between numerator and denominator.
- Only factors remaining are
  - Parents and children  $\mathbf{x}_i$
  - Also co-parents: corresponding to parents of node  $\mathbf{x}_k$  (other than  $\mathbf{x}_i$ )

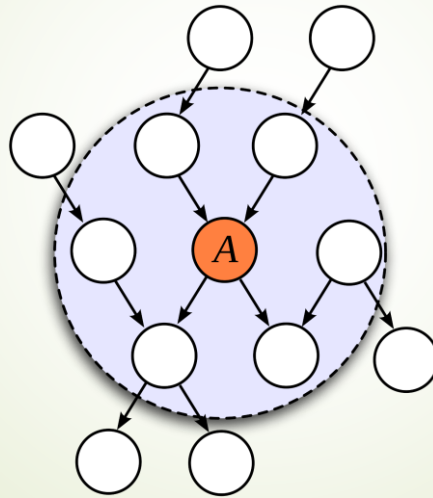
These remaining factors are referred to as The **Markov Blanket** of node  $\mathbf{x}_i$

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## The Markov Blanket



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## Factorization Properties

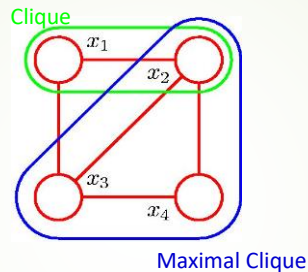
- Consider two nodes  $x_i$  and  $x_j$  that are not connected by a link then these are conditionally independent given all other nodes
- As there is no direct path between the nodes
- All other paths are blocked by nodes that are observed

$$p(x_i, x_j | \mathbf{x}_{\setminus \{i, j\}}) = p(x_i | \mathbf{x}_{\setminus \{i, j\}}) p(x_j | \mathbf{x}_{\setminus \{i, j\}})$$

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## Cliques and Maximal Cliques



- **Clique:** A set of fully connected nodes
- **Maximal Clique:** clique in which it is not possible to include any other nodes without it ceasing to be a clique
- Joint distribution can thus be factored in terms of the functions of maximal cliques
- Functions defined on maximal cliques include the subsets of maximal cliques

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## Joint Distribution

For clique  $C$  and the set of variables in that clique  $\mathbf{x}_C$ , The joint distribution

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

where  $\psi_C(\mathbf{x}_C)$  is the potential over clique  $C$  and the **partition function**  $Z$

$$Z = \sum_{\mathbf{x}} \prod_C \psi_C(\mathbf{x}_C)$$

is the normalization coefficient; note:  $M$   $K$ -state variables  $\rightarrow K^M$  terms in  $Z$ .

$$\psi_C(\mathbf{x}_C) = \exp \{-E(\mathbf{x}_C)\}$$

- For evaluating local marginal probabilities the unnormalized joint distribution can be used

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## Hammersley and Clifford Theorem

- Using filter analogy
- UI: the set of distributions that are consistent with the set of conditional independence statements read from the graph using graph separation
- UF: the set of distributions that can be expressed as a factorization described with respect to the maximal cliques
- The Hammersley-Clifford theorem states that the sets UI and UF are identical if  $\psi_C(\mathbf{x}_C)$  is strictly positive
- In such case

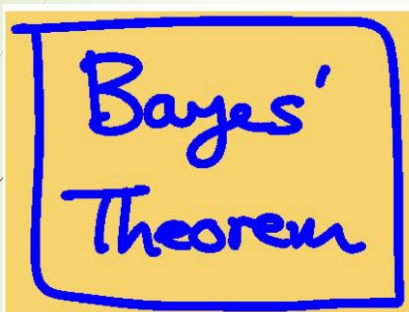
$$\psi_C(\mathbf{x}_C) = \exp\{-E(\mathbf{x}_C)\}$$

where  $E(\mathbf{x}_C)$  is called an energy function, and the exponential representation is called the Boltzmann distribution

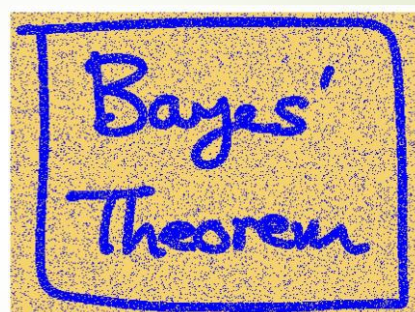
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## Illustration: Image De-Noising (1)



Original Image  $\mathbf{x}$

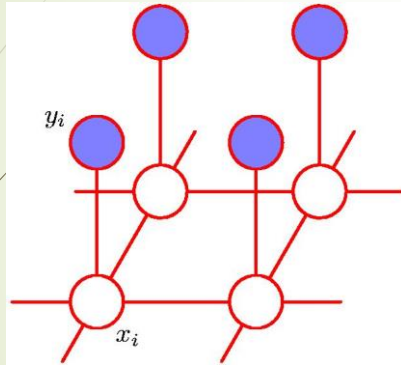


Noisy Image  $\mathbf{y}$  (10% noise)

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## Illustration: Image De-Noising (2)



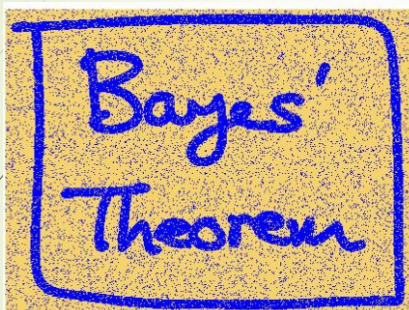
$$E(\mathbf{x}, \mathbf{y}) = h \sum_i x_i - \beta \sum_{\{i,j\}} x_i x_j - \eta \sum_i x_i y_i$$

$$p(\mathbf{x}, \mathbf{y}) = \frac{1}{Z} \exp\{-E(\mathbf{x}, \mathbf{y})\}$$

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## Illustration: Image De-Noising (3)



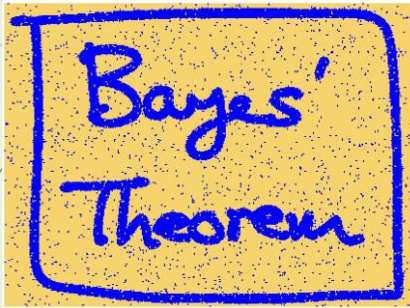
Noisy Image

Restored Image  
(Iterated Conditional Modes; ICM)

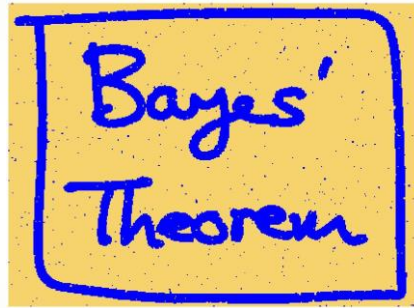
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## Illustration: Image De-Noising (4)



Restored Image (ICM)



Restored Image (Graph cuts)

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## Converting Directed to Undirected Graphs (1)



$$p(\mathbf{x}) = p(x_1)p(x_2|x_1)p(x_3|x_2)\cdots p(x_N|x_{N-1})$$

$$p(\mathbf{x}) = \frac{1}{Z} \psi_{1,2}(x_1, x_2) \psi_{2,3}(x_2, x_3) \cdots \psi_{N-1,N}(x_{N-1}, x_N)$$



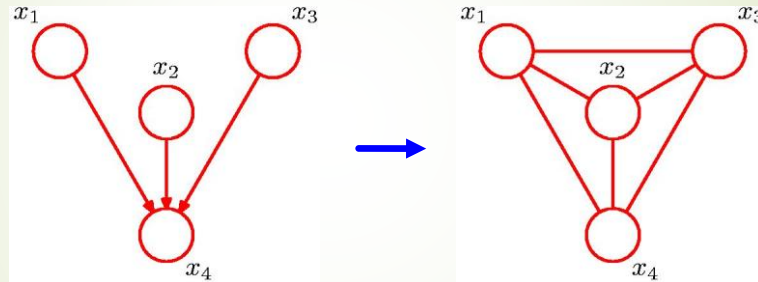
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## Converting Directed to Undirected Graphs (2)

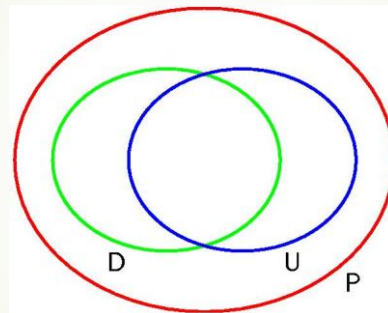
- Additional links between parent nodes (moralization)



$$\begin{aligned}
 p(\mathbf{x}) &= p(x_1)p(x_2)p(x_3)p(x_4|x_1, x_2, x_3) \\
 &= \frac{1}{Z} \psi_A(x_1, x_2, x_3) \psi_B(x_2, x_3, x_4) \psi_C(x_1, x_2, x_4)
 \end{aligned}$$

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## Directed vs. Undirected Graphs (1)



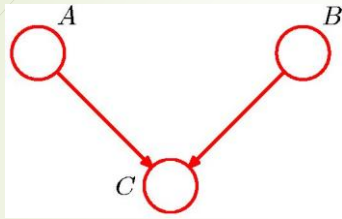
If every conditional independence property of the distribution is reflected in the graph, and vice versa, then the graph is said to be a [perfect map](#) for that distribution

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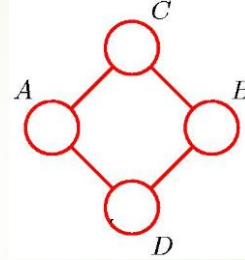
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## Directed vs. Undirected Graphs (2)



$$A \perp\!\!\!\perp B \mid \emptyset$$

$$A \not\perp\!\!\!\perp B \mid C$$



$$A \not\perp\!\!\!\perp B \mid \emptyset$$

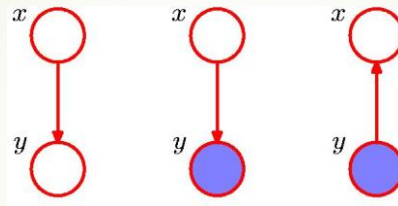
$$A \perp\!\!\!\perp B \mid C \cup D$$

$$C \perp\!\!\!\perp D \mid A \cup B$$

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## Inference in Graphical Models



$$p(y) = \sum_{x'} p(y|x')p(x')$$

$$p(x|y) = \frac{p(y|x)p(x)}{p(y)}$$

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## Inference on a Chain



$$p(\mathbf{x}) = \frac{1}{Z} \psi_{1,2}(x_1, x_2) \psi_{2,3}(x_2, x_3) \cdots \psi_{N-1,N}(x_{N-1}, x_N)$$

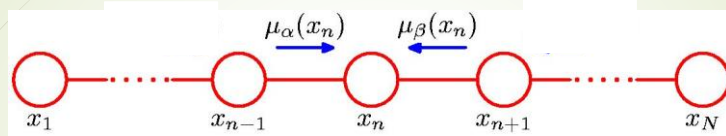
Each discrete variables have  $K$  states, in which case each potential function  $\psi_{n-1,n}(x_{n-1}, x_n)$  comprises a  $K \times K$  table, and so the joint distribution has  $(N - 1)K^2$  parameters

$$p(x_n) = \sum_{x_1} \cdots \sum_{x_{n-1}} \sum_{x_{n+1}} \cdots \sum_{x_N} p(\mathbf{x}) \quad \text{Complexity: } O(K^N)$$

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## Inference on a Chain

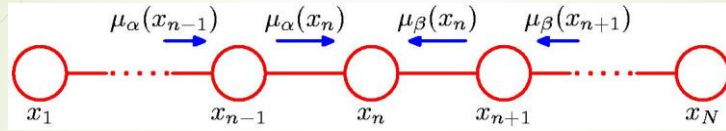


$$p(x_n) = \frac{1}{Z} \underbrace{\left[ \sum_{x_{n-1}} \psi_{n-1,n}(x_{n-1}, x_n) \cdots \left[ \sum_{x_1} \psi_{1,2}(x_1, x_2) \right] \cdots \right]}_{\mu_\alpha(x_n)} \underbrace{\left[ \sum_{x_{n+1}} \psi_{n,n+1}(x_n, x_{n+1}) \cdots \left[ \sum_{x_N} \psi_{N-1,N}(x_{N-1}, x_N) \right] \cdots \right]}_{\mu_\beta(x_n)}$$

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## Inference on Markov Chain: Message Passing

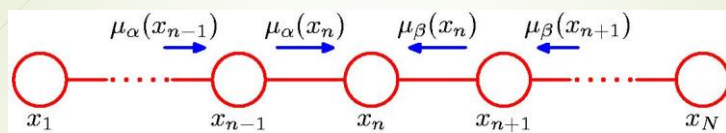


$$\begin{aligned}
 \mu_{\alpha}(x_n) &= \sum_{x_{n-1}} \psi_{n-1,n}(x_{n-1}, x_n) \left[ \sum_{x_{n-2}} \cdots \right] \\
 &= \sum_{x_{n-1}} \psi_{n-1,n}(x_{n-1}, x_n) \mu_{\alpha}(x_{n-1}). \\
 \mu_{\beta}(x_n) &= \sum_{x_{n+1}} \psi_{n,n+1}(x_n, x_{n+1}) \left[ \sum_{x_{n+2}} \cdots \right] \\
 &= \sum_{x_{n+1}} \psi_{n,n+1}(x_n, x_{n+1}) \mu_{\beta}(x_{n+1}).
 \end{aligned}$$

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## Inference on Markov Chain: Message Passing



$$\mu_{\alpha}(x_2) = \sum_{x_1} \psi_{1,2}(x_1, x_2) \qquad \mu_{\beta}(x_{N-1}) = \sum_{x_N} \psi_{N-1,N}(x_{N-1}, x_N)$$

$$Z = \sum_{x_n} \mu_{\alpha}(x_n) \mu_{\beta}(x_n)$$

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## Linear-Complexity Inference on Markov Chain

To compute local marginals:

- Recursively compute and store all forward messages,  $\mu_\alpha(x_n)$ .
- Recursively compute and store all backward messages,  $\mu_\beta(x_n)$ .

- Compute  $Z = \sum_{x_n} \mu_\alpha(x_n) \mu_\beta(x_n)$       Complexity:  $O(K)$

- Compute 
$$p(x_n) = \frac{1}{Z} \mu_\alpha(x_n) \mu_\beta(x_n)$$

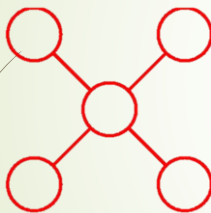
for all variables required.

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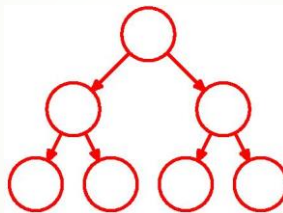
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## Trees

Undirected Tree



Directed Tree



Polytree

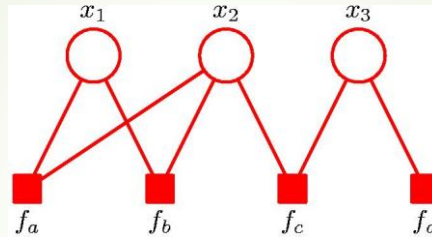


The **sum-product algorithm** provides an efficient framework for exact inference in tree-structured graphs

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## Factor Graphs



Factor graphs are said to be **bipartite** because they consist of two distinct kinds of nodes, and all links go between nodes of opposite type.

- Both directed and undirected graphs allow a global function of several variable to be expressed as a product of factors over subsets of those variables.
- Factor graphs make this decomposition explicit by introducing additional nodes for the factors themselves in addition to the nodes representing the variables.

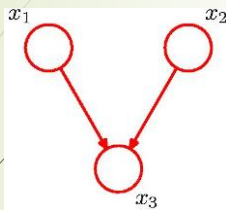
$$p(\mathbf{x}) = f_a(x_1, x_2) f_b(x_1, x_2) f_c(x_2, x_3) f_d(x_3)$$

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

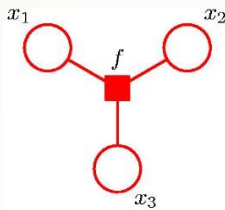
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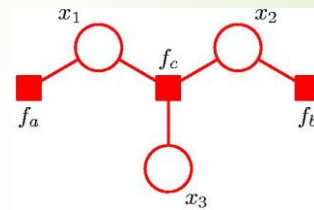
## Factor Graphs from Directed Graphs



$$p(\mathbf{x}) = p(x_1)p(x_2) \\ p(x_3|x_1, x_2)$$



$$f(x_1, x_2, x_3) = \\ p(x_1)p(x_2)p(x_3|x_1, x_2)$$

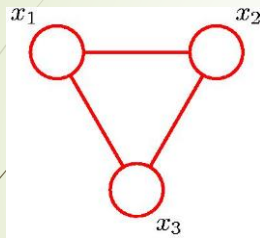


$$f_a(x_1) = p(x_1) \\ f_b(x_2) = p(x_2) \\ f_c(x_1, x_2, x_3) = p(x_3|x_1, x_2)$$

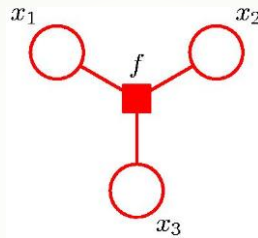
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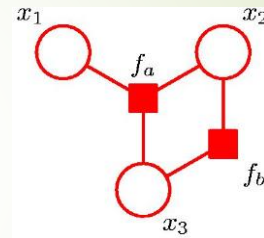
## Factor Graphs from Undirected Graphs



$$\psi(x_1, x_2, x_3)$$



$$\begin{aligned} f(x_1, x_2, x_3) \\ = \psi(x_1, x_2, x_3) \end{aligned}$$



$$\begin{aligned} f_a(x_1, x_2, x_3) f_b(x_2, x_3) \\ = \psi(x_1, x_2, x_3) \end{aligned}$$

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## The Sum-Product Algorithm (1)

### Objective:

- i. to obtain an efficient, exact inference algorithm for finding marginals;
- ii. in situations where several marginals are required, to allow computations to be shared efficiently.

### Key idea: Distributive Law

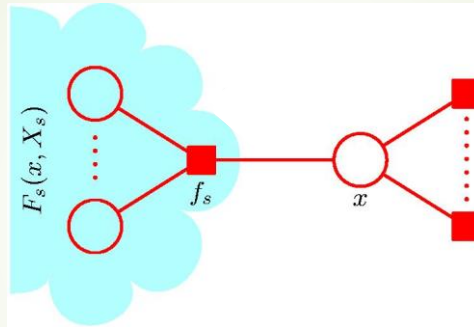
$$ab + ac = a(b + c)$$

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## The Sum-Product Algorithm (2)



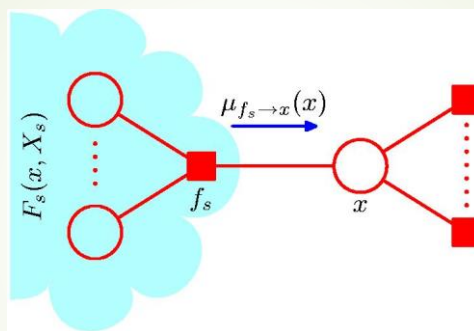
$$p(x) = \sum_{\mathbf{x} \setminus x} p(\mathbf{x})$$

$$p(\mathbf{x}) = \prod_{s \in \text{ne}(x)} F_s(x, X_s)$$

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## The Sum-Product Algorithm (3)



$$p(x) = \prod_{s \in \text{ne}(x)} \left[ \sum_{X_s} F_s(x, X_s) \right]$$

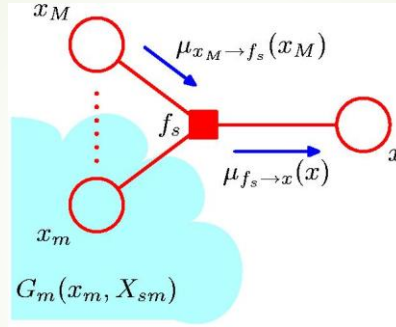
$$= \prod_{s \in \text{ne}(x)} \mu_{f_s \rightarrow x}(x).$$

$$\mu_{f_s \rightarrow x}(x) \equiv \sum_{X_s} F_s(x, X_s)$$

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## The Sum-Product Algorithm (4)

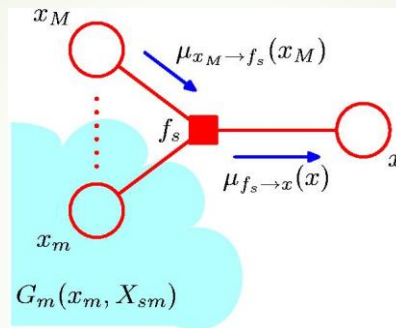


$$F_s(x, X_s) = f_s(x, x_1, \dots, x_M) G_1(x_1, X_{s1}) \dots G_M(x_M, X_{sM})$$

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## The Sum-Product Algorithm (5)

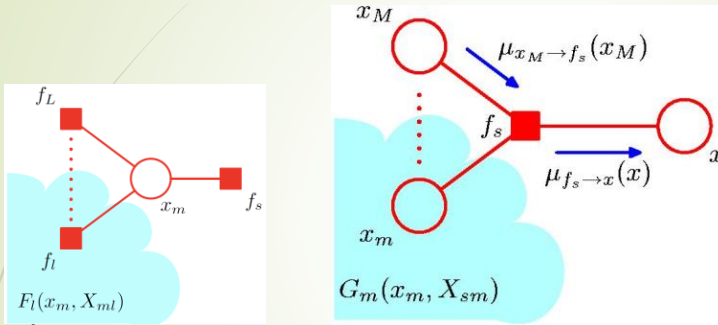


$$\begin{aligned} \mu_{f_s \rightarrow x}(x) &= \sum_{x_1} \dots \sum_{x_M} f_s(x, x_1, \dots, x_M) \prod_{m \in \text{ne}(f_s) \setminus x} \left[ \sum_{X_{sm}} G_m(x_m, X_{sm}) \right] \\ &= \sum_{x_1} \dots \sum_{x_M} f_s(x, x_1, \dots, x_M) \prod_{m \in \text{ne}(f_s) \setminus x} \mu_{x_m \rightarrow f_s}(x_m) \end{aligned}$$

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## The Sum-Product Algorithm (6)



$$\begin{aligned}\mu_{x_m \rightarrow f_s}(x_m) &\equiv \sum_{X_{sm}} G_m(x_m, X_{sm}) = \sum_{X_{sm}} \prod_{l \in \text{ne}(x_m) \setminus f_s} F_l(x_m, X_{ml}) \\ &= \prod_{l \in \text{ne}(x_m) \setminus f_s} \mu_{f_l \rightarrow x_m}(x_m)\end{aligned}$$

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## The Sum-Product Algorithm (7)

Initialization



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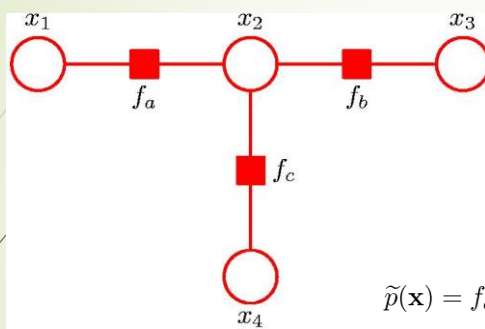
## The Sum-Product Algorithm (8)

- To compute local marginals:
  - Pick an arbitrary node as root
  - Compute and propagate messages from the leaf nodes to the root, storing received messages at every node.
  - Compute and propagate messages from the root to the leaf nodes, storing received messages at every node.
  - Compute the product of received messages at each node for which the marginal is required, and normalize if necessary.

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## Sum-Product: Example (1)

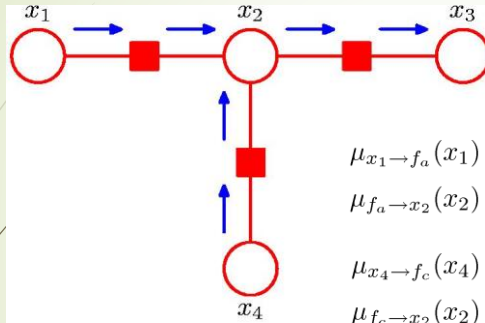


$$\tilde{p}(\mathbf{x}) = f_a(x_1, x_2) f_b(x_2, x_3) f_c(x_2, x_4)$$

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## Sum-Product: Example (2)

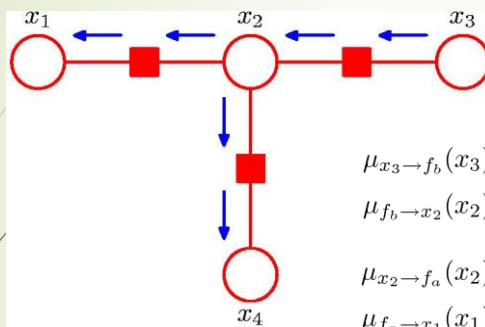


$$\begin{aligned}
 \mu_{x_1 \rightarrow f_a}(x_1) &= 1 \\
 \mu_{f_a \rightarrow x_2}(x_2) &= \sum_{x_1} f_a(x_1, x_2) \\
 \mu_{x_4 \rightarrow f_c}(x_4) &= 1 \\
 \mu_{f_c \rightarrow x_2}(x_2) &= \sum_{x_4} f_c(x_2, x_4) \\
 \mu_{x_2 \rightarrow f_b}(x_2) &= \mu_{f_a \rightarrow x_2}(x_2) \mu_{f_c \rightarrow x_2}(x_2) \\
 \mu_{f_b \rightarrow x_3}(x_3) &= \sum_{x_2} f_b(x_2, x_3) \mu_{x_2 \rightarrow f_b}(x_2)
 \end{aligned}$$

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## Sum-Product: Example (3)

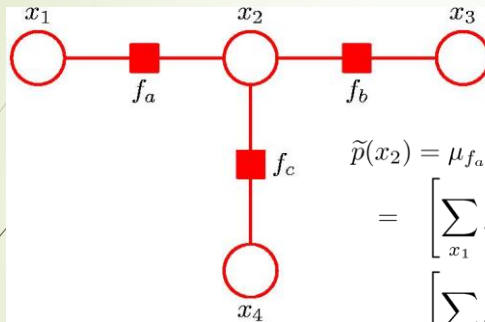


$$\begin{aligned}
 \mu_{x_3 \rightarrow f_b}(x_3) &= 1 \\
 \mu_{f_b \rightarrow x_2}(x_2) &= \sum_{x_3} f_b(x_2, x_3) \\
 \mu_{x_2 \rightarrow f_a}(x_2) &= \mu_{f_b \rightarrow x_2}(x_2) \mu_{f_c \rightarrow x_2}(x_2) \\
 \mu_{f_a \rightarrow x_1}(x_1) &= \sum_{x_2} f_a(x_1, x_2) \mu_{x_2 \rightarrow f_a}(x_2) \\
 \mu_{x_2 \rightarrow f_c}(x_2) &= \mu_{f_a \rightarrow x_2}(x_2) \mu_{f_b \rightarrow x_2}(x_2) \\
 \mu_{f_c \rightarrow x_4}(x_4) &= \sum_{x_2} f_c(x_2, x_4) \mu_{x_2 \rightarrow f_c}(x_2)
 \end{aligned}$$

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## Sum-Product: Example (4)



$$\begin{aligned}
 \tilde{p}(x_2) &= \mu_{f_a \rightarrow x_2}(x_2) \mu_{f_b \rightarrow x_2}(x_2) \mu_{f_c \rightarrow x_2}(x_2) \\
 &= \left[ \sum_{x_1} f_a(x_1, x_2) \right] \left[ \sum_{x_3} f_b(x_2, x_3) \right] \\
 &\quad \left[ \sum_{x_4} f_c(x_2, x_4) \right] \\
 &= \sum_{x_1} \sum_{x_3} \sum_{x_4} f_a(x_1, x_2) f_b(x_2, x_3) f_c(x_2, x_4) \\
 &= \sum_{x_1} \sum_{x_3} \sum_{x_4} \tilde{p}(\mathbf{x})
 \end{aligned}$$

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## The Max-Sum Algorithm (1)

- Objective: an efficient algorithm for finding
  - i. the value  $\mathbf{x}^{\max}$  that maximises  $p(\mathbf{x})$ ;
  - ii. the value of  $p(\mathbf{x}^{\max})$ .
- In general, maximum marginals  $\neq$  joint maximum.

|         | $x = 0$ | $x = 1$ |
|---------|---------|---------|
| $y = 0$ | 0.3     | 0.4     |
| $y = 1$ | 0.3     | 0.0     |

$$\arg \max_x p(x, y) = 1 \quad \arg \max_x p(x) = 0$$

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## The Max-Sum Algorithm (2)

- Maximizing over a chain (max-product)



$$\begin{aligned}
 p(\mathbf{x}^{\max}) &= \max_{\mathbf{x}} p(\mathbf{x}) = \max_{x_1} \dots \max_{x_N} p(\mathbf{x}) \\
 &= \frac{1}{Z} \max_{x_1} \dots \max_{x_N} [\psi_{1,2}(x_1, x_2) \dots \psi_{N-1,N}(x_{N-1}, x_N)] \\
 &= \frac{1}{Z} \max_{x_1} \left[ \max_{x_2} \left[ \psi_{1,2}(x_1, x_2) \left[ \dots \max_{x_N} \psi_{N-1,N}(x_{N-1}, x_N) \right] \dots \right] \right]
 \end{aligned}$$

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## The Max-Sum Algorithm (3)

- Generalizes to tree-structured factor graph

$$\max_{\mathbf{x}} p(\mathbf{x}) = \max_{x_n} \prod_{f_s \in \text{ne}(x_n)} \max_{X_s} f_s(x_n, X_s)$$

- maximizing as close to the leaf nodes as possible

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## The Max-Sum Algorithm (4)

■ Max-Product  $\rightarrow$  Max-Sum

■ For numerical reasons, use

$$\ln \left( \max_{\mathbf{x}} p(\mathbf{x}) \right) = \max_{\mathbf{x}} \ln p(\mathbf{x}).$$

■ Again, use distributive law

$$\max(a + b, a + c) = a + \max(b, c).$$

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## The Max-Sum Algorithm (5)

Initialization (leaf nodes)

$$\mu_{x \rightarrow f}(x) = 0 \qquad \mu_{f \rightarrow x}(x) = \ln f(x)$$

Recursion

$$\begin{aligned} \mu_{f \rightarrow x}(x) &= \max_{x_1, \dots, x_M} \left[ \ln f(x, x_1, \dots, x_M) + \sum_{m \in \text{ne}(f_s) \setminus x} \mu_{x_m \rightarrow f}(x_m) \right] \\ \phi(x) &= \arg \max_{x_1, \dots, x_M} \left[ \ln f(x, x_1, \dots, x_M) + \sum_{m \in \text{ne}(f_s) \setminus x} \mu_{x_m \rightarrow f}(x_m) \right] \\ \mu_{x \rightarrow f}(x) &= \sum_{l \in \text{ne}(x) \setminus f} \mu_{f_l \rightarrow x}(x) \end{aligned}$$

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## The Junction Tree Algorithm

- *Exact* inference on general graphs.
- Works by turning the initial graph into a *junction tree* and then running a sum-product-like algorithm.
- *Intractable* on graphs with large cliques.

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## Loopy Belief Propagation

- Sum-Product on general graphs.
- Initial unit messages passed across all links, after which messages are passed around until convergence (not guaranteed!).
- *Approximate* but *tractable* for large graphs.
- Sometime works well, sometimes not at all.

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