

Readings:
K&F: 10.2, 10.3

Generalized Belief Propagation

Graphical Models – 10708
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November 12th, 2008

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More details on Loopy BP

Numerical problem:

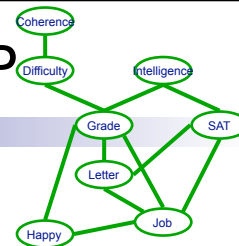
- messages < 1 get multiplied together as we go around the loops
- numbers can go to zero
- normalize messages to one:

$$\delta_{i \rightarrow j}(X_j) = \frac{1}{Z_{i \rightarrow j}} \sum_{x_i} \phi_i(x_i) \phi_{ij}(x_i, X_j) \prod_{k \in \mathcal{N}(i) - j} \delta_{k \rightarrow i}(x_i)$$

- $Z_{i \rightarrow j}$ doesn't depend on X_j , so doesn't change the answer

Computing node “beliefs” (estimates of probs.):

$$b_i(x_i) = \hat{P}(X_i) = \frac{1}{Z_i} \phi_i(X_i) \prod_{k \in \mathcal{N}(i)} \delta_{k \rightarrow i}(X_i)$$



sometimes important to compute in log space

$$\log \delta_{i \rightarrow j}(x_j) = \log$$

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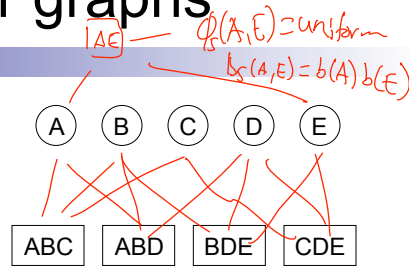
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Loopy BP in Factor graphs

- From node i to factor j :

- $F(i)$ factors whose scope includes X_i

$$\delta_{i \rightarrow j}(X_i) \propto \prod_{k \in F(i) - j} \delta_{k \rightarrow i}(X_i)$$



- From factor j to node i :

- $\text{Scope}[\phi_j] = Y \cup \{X_i\}$

$$\delta_{j \rightarrow i}(X_i) \propto \sum_{\mathbf{y}} \phi_j(X_i, \mathbf{y}) \prod_{X_k \in \text{Scope}[\phi_j] - X_i} \delta_{k \rightarrow j}(x_k)$$

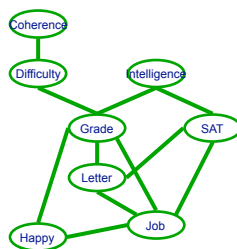
- Belief:

- Node: $P(x_i) \approx b_i(x_i) \propto \prod_{\phi_j, x_i \in \text{Scope}[\phi_j]} \delta_{j \rightarrow x_i}(x_i)$
- Factor: $\phi_j(Y) \approx b_{\phi_j}(Y) \propto \prod_{x_i \in Y} \delta_{x_i \rightarrow \phi_j}(x_i)$

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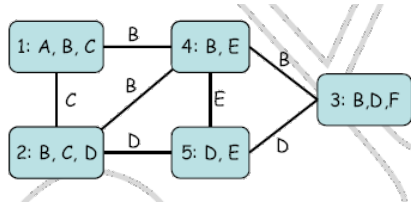
Loopy BP v. Clique trees: Two ends of a spectrum



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Generalize cluster graph



■ Generalized cluster graph:

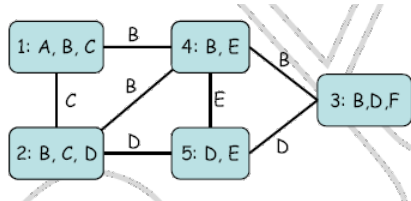
For set of factors F

- Undirected graph
- Each node i associated with a cluster C_i
- *Family preserving*: for each factor $f_j \in F$, \exists node i such that $\text{scope}[f_j] \subseteq C_i$
- Each edge $i - j$ is associated with a set of variables $S_{ij} \subseteq C_i \cap C_j$

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Running intersection property



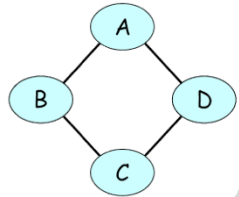
■ (Generalized) Running intersection property (RIP)

- Cluster graph satisfies RIP if whenever $X \in C_i$ and $X \in C_j$ then \exists one and only one path from C_i to C_j where $X \in S_{uv}$ for every edge (u,v) in the path

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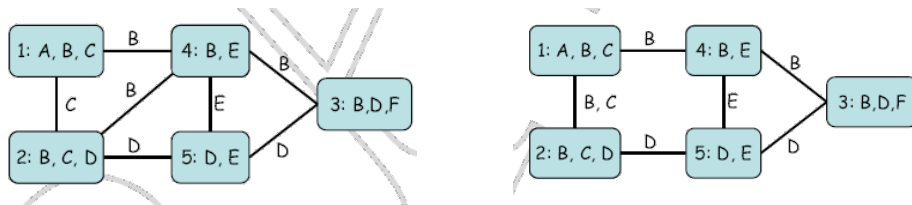
Examples of cluster graphs



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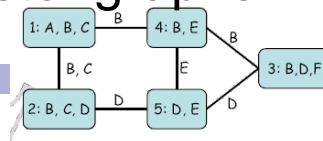
Two cluster graph satisfying RIP with different edge sets



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Generalized BP on cluster graphs satisfying RIP



Initialization:

- Assign each factor ϕ to a clique $\alpha(\phi)$, $\text{Scope}[\phi] \subseteq \mathbf{C}_{\alpha(\phi)}$
- Initialize cliques: $\psi_i^0(\mathbf{C}_i) \propto \prod_{\phi: \alpha(\phi)=i} \phi$
- Initialize messages: $\delta_{j \rightarrow i} = 1$

While not converged, send messages:

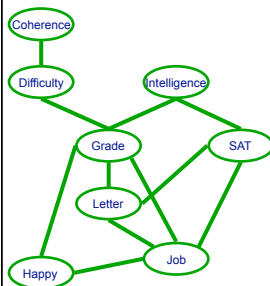
$$\delta_{i \rightarrow j}(\mathbf{S}_{ij}) \propto \sum_{\mathbf{C}_i - \mathbf{S}_{ij}} \psi_i^0(\mathbf{C}_i) \prod_{k \in \mathcal{N}(i) - j} \delta_{k \rightarrow i}(\mathbf{S}_{ik})$$

Belief:

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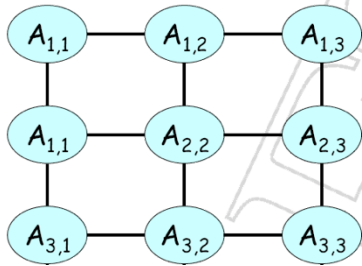
Cluster graph for Loopy BP



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What if the cluster graph doesn't satisfy RIP



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Region graphs to the rescue

- Can address generalized cluster graphs that don't satisfy RIP using *region graphs*:
 - Book: 10.3
- Example in your homework! ☺

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Revisiting Mean-Fields

$$\ln Z = F[P_{\mathcal{F}}, Q] + D(Q||P_{\mathcal{F}}) \quad F[P_{\mathcal{F}}, Q] = \sum_{\phi \in \mathcal{F}} E_Q[\ln \phi] + H_Q(\mathcal{X})$$

- Choice of Q:
- Optimization problem:

$$\max_Q F[P_{\mathcal{F}}, Q] = \sum_{\phi \in \mathcal{F}} E_Q[\ln \phi] + \sum_j H_{Q_j}(X_j), \quad \forall i, \sum_{x_i} Q_i(x_i) = 1$$

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Announcements

- Recitation tomorrow
- HW5 out soon
- Will not cover relational models this semester
 - Instead, recommend Pedro Domingos' tutorial on Markov Logic
 - Markov logic is one example of a relational probabilistic model
 - November 14th from 1:00 pm to 3:30 pm in Wean 4623

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Interpretation of energy functional

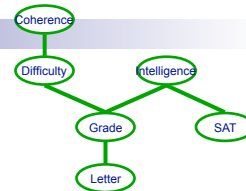
- Energy functional: $F[P_{\mathcal{F}}, Q] = \sum_{\phi \in \mathcal{F}} E_Q[\ln \phi] + H_Q(\mathcal{X})$
- Exact if $P=Q$: $\ln Z = F[P_{\mathcal{F}}, Q] + D(Q||P_{\mathcal{F}})$
- View problem as an approximation of entropy term:

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Entropy of a tree distribution

- Entropy term:
- Joint distribution:
- Decomposing entropy term:



- More generally: $H_P(\mathbf{X}) = \sum_{(i,j) \in E} H(X_i, X_j) - \sum_i (d_i - 1) H(X_i)$
 - d_i number neighbors of X_i

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Loopy BP & Bethe approximation

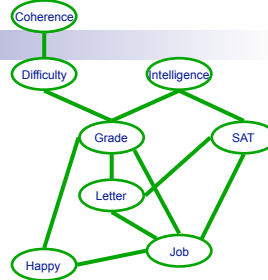
■ Energy functional: $F[P_{\mathcal{F}}, Q] = \sum_{\phi \in \mathcal{F}} E_Q[\ln \phi] + H_Q(\mathcal{X})$

■ Bethe approximation of Free Energy:

□ use entropy for trees, but loopy graphs:

$$\tilde{F}[P_{\mathcal{F}}, Q] = \sum_{(i,j) \in E} E_{b_{ij}}[\ln \phi_{ij}] + \sum_{(i,j) \in E} H_{b_{ij}}(X_i, X_j) - \sum_i (d_i - 1) H_{b_i}(X_i)$$

■ **Theorem:** If Loopy BP converges, resulting b_{ij} & b_i are stationary point (usually local maxima) of Bethe Free energy!



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GBP & Kikuchi approximation

■ Exact Free energy: Junction Tree

$$F[P_{\mathcal{F}}, Q] = \sum_{(i,j) \in E} E_{\pi_{ij}}[\ln \phi_{ij}] + \sum_i H_{\pi_{C_i}}(C_i) - \sum_{(i,j) \in T} H_{\pi_{S_{ij}}}(S_{ij})$$

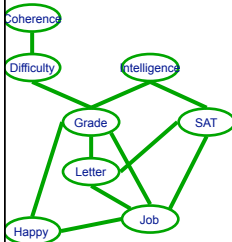
■ Bethe Free energy:

$$\tilde{F}[P_{\mathcal{F}}, Q] = \sum_{(i,j) \in E} E_{b_{ij}}[\ln \phi_{ij}] + \sum_{(i,j) \in E} H_{b_{ij}}(X_i, X_j) - \sum_i (d_i - 1) H_{b_i}(X_i)$$

■ Kikuchi approximation: Generalized cluster graph

□ spectrum from Bethe to exact

■ **Theorem:** If GBP converges, resulting b_{C_i} are stationary point (usually local maxima) of Kikuchi Free energy!



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What you need to know about GBP

- Spectrum between Loopy BP & Junction Trees:
 - More computation, but typically better answers
- If satisfies RIP, equations are very simple
- General setting, slightly trickier equations, but not hard
- Relates to variational methods: Corresponds to local optima of approximate version of energy functional

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Parameter learning in Markov nets

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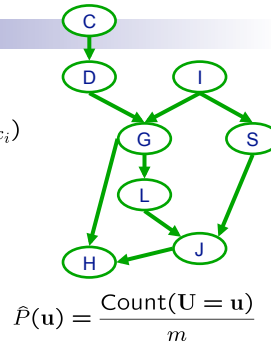
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Learning Parameters of a BN

- Log likelihood decomposes:

$$\ell(\mathcal{D} : \theta) = \log P(\mathcal{D} | \theta) = m \sum_i \sum_{x_i, \text{Pa}_{x_i}} \hat{P}(x_i, \text{Pa}_{x_i}) \log P(x_i | \text{Pa}_{x_i})$$

- Learn each CPT independently
- Use counts

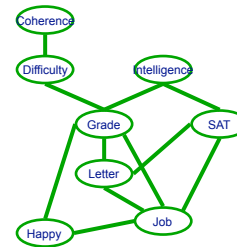


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Log Likelihood for MN

- Log likelihood of the data:



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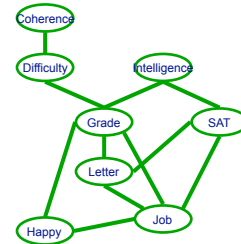
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Log Likelihood doesn't decompose for MNs

$$\hat{P}(\mathbf{u}) = \frac{\text{Count}(\mathbf{U} = \mathbf{u})}{m}$$

Log likelihood:

$$\ell(\mathcal{D} : \theta) = \log P(\mathcal{D} | \theta, \mathcal{G}) = m \sum_i \sum_{\mathbf{c}_i} \hat{P}(\mathbf{c}_i) \log \psi_i(\mathbf{c}_i) - m \log Z$$



A convex problem

- Can find global optimum!!

Term log Z doesn't decompose!!

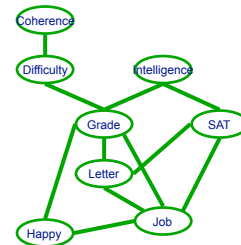
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Derivative of Log Likelihood for MNs

$$\hat{P}(\mathbf{u}) = \frac{\text{Count}(\mathbf{U} = \mathbf{u})}{m}$$

$$\ell(\mathcal{D} : \theta) = \log P(\mathcal{D} | \theta, \mathcal{G}) = m \sum_i \sum_{\mathbf{c}_i} \hat{P}(\mathbf{c}_i) \log \psi_i(\mathbf{c}_i) - m \log Z$$



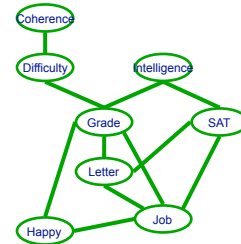
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Derivative of Log Likelihood for MNs 2

$$\hat{P}(u) = \frac{\text{Count}(U = u)}{m}$$

$$\ell(\mathcal{D} : \theta) = \log P(\mathcal{D} | \theta, \mathcal{G}) = m \sum_i \sum_{\mathbf{c}_i} \hat{P}(\mathbf{c}_i) \log \psi_i(\mathbf{c}_i) - m \log Z$$



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Derivative of Log Likelihood for MNs

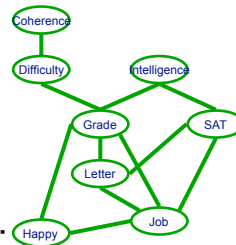
$$\hat{P}(u) = \frac{\text{Count}(U = u)}{m}$$

$$\ell(\mathcal{D} : \theta) = \log P(\mathcal{D} | \theta, \mathcal{G}) = m \sum_i \sum_{\mathbf{c}_i} \hat{P}(\mathbf{c}_i) \log \psi_i(\mathbf{c}_i) - m \log Z$$

- Derivative:

$$\frac{\partial \ell}{\partial \psi_i(\mathbf{c}_i)} = \frac{m \hat{P}(\mathbf{c}_i)}{\psi_i(\mathbf{c}_i)} - \frac{m P_{\mathcal{F}}^{\psi}(\mathbf{c}_i)}{\psi_i(\mathbf{c}_i)}$$

- Computing derivative requires inference:



- Can optimize using gradient ascent
 - Common approach
 - Conjugate gradient, Newton's method,...
- Let's also look at a simpler solution

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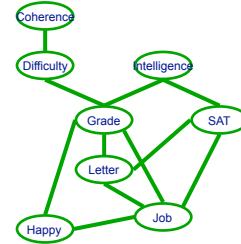
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Iterative Proportional Fitting (IPF)

$$\hat{P}(u) = \frac{\text{Count}(U = u)}{m}$$

$$\frac{\partial \ell}{\partial \psi_i(\mathbf{c}_i)} = \frac{m \hat{P}(\mathbf{c}_i)}{\psi_i(\mathbf{c}_i)} - \frac{m P_{\mathcal{F}}^{\psi}(\mathbf{c}_i)}{\psi_i(\mathbf{c}_i)}$$

- Setting derivative to zero:
- Fixed point equation:
- Iterate and converge to optimal parameters
 - Each iteration, must compute:



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What you need to know about learning MN parameters?

- BN parameter learning easy
- MN parameter learning doesn't decompose!
- Learning requires inference!
- Apply gradient ascent or IPF iterations to obtain optimal parameters

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