## Learning parameters

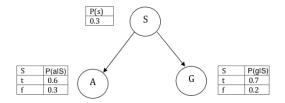
Qinfeng (Javen) Shi

1 May 2017

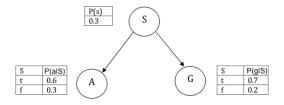
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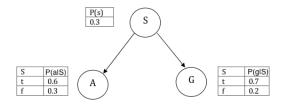


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- $\operatorname{argmax}_{G,A,S} P(G,A,S)$ ? (MAP inference)



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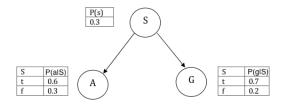
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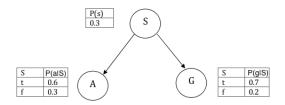


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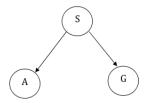
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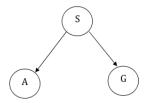
 $\rightarrow$  Learning structures

For bayesian networks,  $P(x_1, ..., x_n) = \prod_{i=1}^n P(x_i | Pa(x_i))$ . Parameters:  $P(x_i | Pa(x_i))$ .

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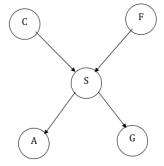


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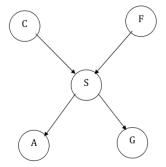


Parameters: P(S), P(A|S), P(G|S)

#### Parameters?

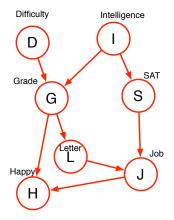


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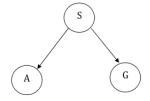


Parameters: P(C), P(F), P(S|C,F), P(A|S), P(G|S)

#### Parameters?



## How to learn the parameters from the data?



Parameters: P(S), P(A|S), P(G|S)

Data:

S	Α	G
1	0	0
0	0	1
1	1	0
1	1	0

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Parameters: P(S), P(A|S), P(G|S) Data:

$$P(S = 0) \approx \frac{N_{(S=0)}}{N_{(S=0)} + N_{(S=1)}} = \frac{1}{4}$$

$$P(S = 1) \approx \frac{N_{(S=1)}}{N_{(S=0)} + N_{(S=1)}} = \frac{3}{4}$$

$$P(A = 0|S = 0) \approx \frac{N_{(A=0,S=0)}}{N_{(S=0)}} = \frac{1}{1}$$

Parameters: P(S), P(A|S), P(G|S)What if we change the data only by one entry (instance)?

S. A. S. S. A. S.

S	Α	G		S	Α	G
1	0	0		1	0	0
0	0	1	$\rightarrow$	1	0	1
1	1	0		1	1	0
1	1	0		1	1	0

Parameters: P(S), P(A|S), P(G|S)

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 ?!

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Solution 1: set P(A|S=0) to be uniform distribution when  $N_{(S=0)}=0$ . Why?

Solution 2 (better): set (no need to check if the denominator)

$$P(A = 0|S = 0) \approx \frac{N_{(A=0,S=0)} + N_r}{N_{(S=0)} + (\#A) \times N_r}$$

Often  $N_r = 1$ . #A is the number of values of variable A can take.

### General solution when the denominator is 0

Let A, B, C, D, ... be the variables. To estimate P(A = 0|B = 0, C = 0).

$$P(A = 0|B = 0, C = 0) \approx \frac{N_{(A=0,B=0,C=0)}}{N_{(B=0,C=0)}}$$

What if  $N_{(B=0,C=0)}=0$ ? This means  $N_{(A=0,B=0,C=0)}=0$  and  $N_{(A=1,B=0,C=0)}=0$ .

Solution 1: When this happens, set P(A|B=0,C=0) to be uniform distribution.

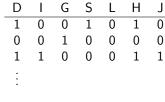
Solution 2 (better): Always set (no need to check if the denominator = 0 or not)

$$P(A = 0|B = 0, C = 0) \approx \frac{N_{(A=0,B=0,C=0)} + N_r}{N_{(B=0,C=0)} + (\#A) \times N_r}$$

Often  $N_r = 1$ . #A is the number of values of variable A can take.

# A general case (Student model)

### Data:



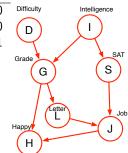
$$P(D=0) = \frac{N_{(D=0)}}{N_{total}}$$

$$P(D=1) = \frac{N_{(D=1)}}{N_{total}}$$

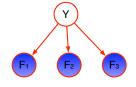
$$P(D = 0) = \frac{N_{(D=0)}}{N_{total}}$$

$$P(D = 1) = \frac{N_{(D=1)}}{N_{total}}$$

$$P(G = 0|D = 0, I = 1) = \frac{N_{(G=0,D=0,I=1)}}{N_{(D=0,I=1)}}$$

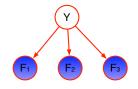


## A special case (Naive Bayes)

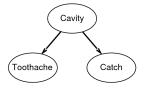


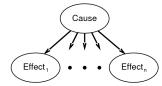
Parameters:  $P(F_1|Y), P(F_2|Y), ...$ 

## A special case (Naive Bayes)



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## More problems?

- not minimise classification error or other measure of your task.
- not much flexibility on the features nor the parameters.

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- not minimise classification error or other measure of your task.
- not much flexibility on the features nor the parameters.

Alternatives: using MRFs or factor graphical models.

### Parameters for MRFs

For MRFs, let V be the set of nodes, and C be the set of clusters c.

$$P(\mathbf{x};\theta) = \frac{\exp(\sum_{c \in \mathcal{C}} \theta_c(\mathbf{x}_c))}{Z(\theta)},$$
 (1)

where normaliser  $Z(\theta) = \sum_{\mathbf{x}} \exp\{\sum_{c'' \in \mathcal{C}} \theta_{c''}(\mathbf{x}_{c''})\}$ . Parameters:  $\{\theta_c\}_{c \in \mathcal{C}}$  or  $\mathbf{w}$ .

- Often assume  $\theta_c(\mathbf{x}_c) = \langle \mathbf{w}, \Phi_c(\mathbf{x}_c) \rangle$ .
- w ← empirical risk minimisation (ERM).

### Inference:

- MAP inference  $\mathbf{x}^* = \operatorname{argmax}_{\mathbf{x}} \sum_{c \in \mathcal{C}} \theta_c(\mathbf{x}_c)$ (hint:  $\log P(\mathbf{x}) \propto \sum_{c \in \mathcal{C}} \theta_c(\mathbf{x}_c)$ )
- Marginal inference  $P(\mathbf{x}_c) = \sum_{\mathbf{x}_{v/c}} P(\mathbf{x})$

### Parameters for MRFs

In learning, we look for a F that predicts labels well via

$$\mathbf{y}^* = \max_{\mathbf{y} \in \mathcal{Y}} F(\mathbf{x}_i, \mathbf{y}; \mathbf{w}).$$

Given graph G = (V, E), one often assume

$$\mathbf{w} = \begin{bmatrix} \mathbf{w}_1 \\ \mathbf{w}_2 \end{bmatrix}, \qquad \Phi(\mathbf{x}, \mathbf{y}) = \begin{bmatrix} \sum_{i \in V} \Phi_i(y^{(i)}, \mathbf{x}) \\ \sum_{(i,j) \in E} \Phi_{i,j}(y^{(i)}, y^{(j)}, \mathbf{x}) \end{bmatrix}$$

$$F(\mathbf{x}, \mathbf{y}; \mathbf{w}) = \langle \mathbf{w}, \Phi(\mathbf{x}, \mathbf{y}) \rangle$$

$$= \sum_{i \in V} \langle \mathbf{w}_1, \Phi_i(y^{(i)}, \mathbf{x}) \rangle + \sum_{(i,j) \in E} \langle \mathbf{w}_2, \Phi_{i,j}(y^{(i)}, y^{(j)}, \mathbf{x}) \rangle$$

$$= \sum_{i \in V} \theta_i(y^{(i)}, \mathbf{x}) + \sum_{(i,i) \in E} \theta_{i,j}(y^{(i)}, y^{(j)}, \mathbf{x})$$

# Max Margin Approaches

A gap between  $F(\mathbf{x}_i, \mathbf{y}_i; \mathbf{w})$  and best  $F(\mathbf{x}_i, \mathbf{y}; \mathbf{w})$  for  $\mathbf{y} \neq \mathbf{y}_i$ , that is

$$F(\mathbf{x}_i, \mathbf{y}_i; \mathbf{w}) - \max_{\mathbf{y} \in \boldsymbol{\vartheta}, \mathbf{y} \neq \mathbf{y}_i} F(\mathbf{x}_i, \mathbf{y}; \mathbf{w})$$

### Structured SVM - 1

Primal:

$$\min_{\mathbf{w},\xi} \ \frac{1}{2} \|\mathbf{w}\|^2 + C \sum_{i=1}^m \xi_i \quad \text{s.t.}$$
 (2a)

$$\forall i, \mathbf{y} \neq \mathbf{y}_i, \langle \mathbf{w}, \Phi(\mathbf{x}_i, \mathbf{y}_i) - \Phi(\mathbf{x}_i, \mathbf{y}) \rangle \ge \Delta(\mathbf{y}_i, \mathbf{y}) - \xi_i.$$
 (2b)

Dual is a quadratic programming (QP) problem similar to binary SVM's dual.

### Structured SVM - 2

Cutting plane method needs to find the label for the most violated constraint in (2b)

$$\mathbf{y}_{i}^{\dagger} = \operatorname*{argmax}_{\mathbf{y} \in \mathcal{Y}} \Delta(\mathbf{y}_{i}, \mathbf{y}) + \langle \mathbf{w}, \Phi(\mathbf{x}_{i}, \mathbf{y}) \rangle. \tag{3}$$

With  $\mathbf{y}_{i}^{\dagger}$ , one can solve following relaxed problem (with much fewer constraints)

$$\min_{\mathbf{w},\xi} \ \frac{1}{2} \|\mathbf{w}\|^2 + C \sum_{i=1}^m \xi_i \quad \text{s.t.}$$
 (4a)

$$\forall i, \left\langle \mathbf{w}, \Phi(\mathbf{x}_i, \mathbf{y}_i) - \Phi(\mathbf{x}_i, \mathbf{y}_i^{\dagger}) \right\rangle \ge \Delta(\mathbf{y}_i, \mathbf{y}_i^{\dagger}) - \xi_i.$$
 (4b)

### Structured SVM - 3

```
Input: data x_i, labels y_i, sample size m, number of iterations T
Initialise S_0 = \emptyset, \mathbf{w}_0 = 0 (or a random vector), and t = 0.
for t = 0 to T do
    for i = 1 to m do
         \mathbf{y}_{i}^{\dagger} = \operatorname{argmax}_{\mathbf{y} \in \mathcal{Y}, \mathbf{y} \neq \mathbf{y}_{i}} \langle \mathbf{w}_{t}, \Phi(\mathbf{x}_{i}, \mathbf{y}) \rangle + \Delta(\mathbf{y}_{i}, \mathbf{y}),
         \boldsymbol{\xi_i} = \left[ \Delta(\mathbf{y}_i, \mathbf{y}) + \left\langle \mathbf{w}_t, \left( \Phi(\mathbf{x}_i, \mathbf{y}_i^{\dagger}) - \Phi(\mathbf{x}_i, \mathbf{y}_i) \right) \right\rangle \right]_+,
         if \xi_i > 0 then
              Increase constraint set S_t \leftarrow S_t \cup \{\mathbf{y}_t^{\dagger}\}\
         end if
    end for
    \mathbf{w}_t recovered using dual variables.
    \alpha \leftarrow optimise dual QP with constraint set S_t.
end for
```

## Other Max Margin Approaches

Other approaches using Max Margin principle such as Max Margin Markov Network (M3N), ...

## Probabilistic Approaches

### Main types:

- Maximum Entropy (MaxEnt)
- Maximum a Posteriori (MAP)
- Maximum Likelihood (ML)

## Maximum Entropy

Maximum Entropy (ME) estimates **w** by maximising the entropy. That is,

$$\mathbf{w}^* = \operatorname*{argmax}_{\mathbf{w}} \sum_{\mathbf{x} \in \mathcal{X}, \mathbf{y} \in \mathcal{Y}} - \mathbf{P}_{\mathbf{w}}(\mathbf{x}, \mathbf{y}) \ln \mathbf{P}_{\mathbf{w}}(\mathbf{x}, \mathbf{y}).$$

Duality between maximum likelihood, and maximum entropy, subject to moment matching constraints on the expectations of features.

### MAP

Let likelihood function  $\mathcal{L}(\mathbf{w})$  be the modelled probability or density for the occurrence of a sample configuration  $(\mathbf{x}_1, \mathbf{y}_1), \dots, (\mathbf{x}_m, \mathbf{y}_m)$  given the probability density  $\mathbf{P}_{\mathbf{w}}$  parameterised by  $\mathbf{w}$ . That is,

$$\mathcal{L}(\mathbf{w}) = \mathbf{P}_{\mathbf{w}} ((\mathbf{x}_1, \mathbf{y}_1), \dots, (\mathbf{x}_m, \mathbf{y}_m)).$$

Maximum a Posteriori (MAP) estimates  $\mathbf{w}$  by maximising  $\mathcal{L}(\mathbf{w})$  times a prior  $P(\mathbf{w})$ . That is

$$\mathbf{w}^* = \operatorname*{argmax}_{\mathbf{w}} \mathcal{L}(\mathbf{w}) P(\mathbf{w}). \tag{5}$$

Assuming  $\{(\mathbf{x}_i, \mathbf{y}_i)\}_{1 \leq i \leq m}$  are I.I.D. samples from  $\mathbf{P}_{\mathbf{w}}(\mathbf{x}, \mathbf{y})$ , (5) becomes

$$\begin{split} \mathbf{w}^* &= \underset{\mathbf{w}}{\operatorname{argmax}} \prod_{1 \leq i \leq m} \mathbf{P}_{\mathbf{w}}(\mathbf{x}_i, \mathbf{y}_i) P(\mathbf{w}) \\ &= \underset{\mathbf{w}}{\operatorname{argmin}} \sum_{1 \leq i \leq m} -\ln \mathbf{P}_{\mathbf{w}}(\mathbf{x}_i, \mathbf{y}_i) - \ln P(\mathbf{w}). \end{split}$$

### Maximum Likelihood

Maximum Likelihood (ML) is a special case of MAP when  $P(\mathbf{w})$  is uniform which means

$$\mathbf{w}^* = \underset{\mathbf{w}}{\operatorname{argmax}} \prod_{1 \leq i \leq m} \mathbf{P}_{\mathbf{w}}(\mathbf{x}_i, \mathbf{y}_i)$$
$$= \underset{\mathbf{w}}{\operatorname{argmin}} \sum_{1 \leq i \leq m} - \ln \mathbf{P}_{\mathbf{w}}(\mathbf{x}_i, \mathbf{y}_i).$$

Alternatively, one can replace the joint distribution  $P_w(x,y)$  by the conditional distribution  $P_w(y \mid x)$  that gives a discriminative model called Conditional Random Fields (CRFs)

## Conditional Random Fields (CRFs) - 1

Assume the conditional distribution over  $\mathcal{Y} \mid \mathcal{X}$  has a form of exponential families, *i.e.*,

$$P(\mathbf{y} \mid \mathbf{x}; \mathbf{w}) = \frac{\exp(\langle \mathbf{w}, \Phi(\mathbf{x}, \mathbf{y}) \rangle)}{Z(\mathbf{w}, \mathbf{x})},$$
 (6)

where

$$Z(\mathbf{w}, \mathbf{x}) = \sum_{\mathbf{y}' \in \mathcal{Y}} \exp(\langle \mathbf{w}, \Phi(\mathbf{x}, \mathbf{y}') \rangle), \tag{7}$$

and

$$\mathbf{w} = \begin{bmatrix} \mathbf{w}_1 \\ \mathbf{w}_2 \end{bmatrix}, \qquad \Phi(\mathbf{x}, \mathbf{y}) = \begin{bmatrix} \sum_{i \in V} \Phi_i(y^{(i)}, \mathbf{x}) \\ \sum_{(i,j) \in E} \Phi_{i,j}(y^{(i)}, y^{(j)}, \mathbf{x}) \end{bmatrix}$$

More generally speaking, the global feature can be decomposed into local features on cliques (fully connected subgraphs).

Denote  $(\mathbf{x}_1,\ldots,\mathbf{x}_m)$  as  $\mathbf{X}$ ,  $(\mathbf{y}_1,\ldots,\mathbf{y}_m)$  as  $\mathbf{Y}$ . The classical approach is to maximise the conditional likelihood of  $\mathbf{Y}$  on  $\mathbf{X}$ , incorporating a prior on the parameters. This is a Maximum a Posteriori (MAP) estimator, which consists of maximising

$$P(w | X, Y) \propto P(w) P(Y | X; w).$$

From the i.i.d. assumption we have

$$\mathbf{P}(\mathbf{Y} \mid \mathbf{X}; \mathbf{w}) = \prod_{i=1}^{m} \mathbf{P}(\mathbf{y}_{i} \mid \mathbf{x}_{i}; \mathbf{w}),$$

and we impose a Gaussian prior on w

$$P(\mathbf{w}) \propto \exp\left(rac{-||\mathbf{w}||^2}{2\sigma^2}
ight).$$

Maximising the posterior distribution can also be seen as minimising the negative log-posterior, which becomes our risk function  $R(\mathbf{w}, \mathbf{X}, \mathbf{Y})$ 

$$R(\mathbf{w}, \mathbf{X}, \mathbf{Y}) = -\ln(P(\mathbf{w}) \mathbf{P}(\mathbf{Y} \mid \mathbf{X}; \mathbf{w})) + c$$

$$= \frac{||\mathbf{w}||^2}{2\sigma^2} - \sum_{i=1}^{m} \underbrace{\left(\langle \Phi(\mathbf{x}_i, \mathbf{y}_i), \mathbf{w} \rangle - \ln(Z(\mathbf{w}, \mathbf{x}_i))\right)}_{:=\ell_L(\mathbf{x}_i, \mathbf{y}_i, \mathbf{w})} + c,$$

where c is a constant and  $\ell_L$  denotes the log loss *i.e.* negative log-likelihood. Now learning is equivalent to

$$\mathbf{w}^* = \underset{\mathbf{w}}{\operatorname{argmin}} R(\mathbf{w}, \mathbf{X}, \mathbf{Y}).$$

Above is a convex optimisation problem on  $\mathbf{w}$  since  $\ln Z(\mathbf{w}, \mathbf{x})$  is a convex function of  $\mathbf{w}$ . The solution can be obtained by gradient descent since  $\ln Z(\mathbf{w}, \mathbf{x})$  is also differentiable. We have

$$\nabla_{\mathbf{w}} R(\mathbf{w}, \mathbf{X}, \mathbf{Y}) = \frac{\mathbf{w}}{\sigma^2} - \sum_{i=1}^m \left( \Phi(\mathbf{x}_i, \mathbf{y}_i) - \nabla_{\mathbf{w}} \ln(Z(\mathbf{w}, \mathbf{x}_i)) \right).$$

It follows from direct computation that

$$abla_{\mathbf{w}} \ln(Z(\mathbf{w}, \mathbf{x})) = \mathbb{E}_{\mathbf{y} \sim \mathbf{P}(\mathbf{y} \,|\, \mathbf{x}; \mathbf{w})} [\Phi(\mathbf{x}, \mathbf{y})].$$

Since  $\Phi(\mathbf{x}, \mathbf{y})$  are decomposed over nodes and edges, it is straightforward to show that the expectation also decomposes into expectations on nodes  $\mathcal{V}$  and edges  $\mathcal{E}$ 

$$\begin{split} & \mathbb{E}_{\mathbf{y} \sim \mathbf{P}(\mathbf{y} \mid \mathbf{x}; \mathbf{w})} [\Phi(\mathbf{x}, \mathbf{y})] = \\ & \sum_{i \in \mathcal{V}} \mathbb{E}_{y^{(i)} \sim \mathbf{P}(y^{(i)} \mid \mathbf{x}; \mathbf{w})} [\Phi_i(y^{(i)}, \mathbf{x})] \\ & + \sum_{(ij) \in \mathcal{E}} \mathbb{E}_{y^{(i)}, y^{(j)} \sim \mathbf{P}(y^{(i)}, y^{(j)} \mid \mathbf{x}; \mathbf{w})} [\Phi_{i,j}(y^{(i)}, y^{(j)}, \mathbf{x})], \end{split}$$

where the node and edge expectations can be computed given  $\mathbf{P}(y^{(i)}|\mathbf{x};\mathbf{w})$  and  $\mathbf{P}(y^{(i)},y^{(j)}|\mathbf{x};\mathbf{w})$ , which can be computed by Marginal inference methods such as variable elimination, junction tree, e.g. (loopy) belief propagation, or being circumvented through sampling.

Parameters for MRFs Max Margin Approaches (e.g. Structured SVMs) Probabilistic Approaches (e.g. CRFs)

### Break

Take a break ...