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CS 584 - Summer 2020.

1,

- ~~Gene~~ In generative learning, we model the feature distribution and class priors \rightarrow allow generating new examples and also do classification.
- In discriminative learning, we model $P(Y|X) \rightarrow$ strictly doing prediction.

- Given nD feature vector:

$$g(x) = \theta^T x = \theta_0 + \theta_1 x_1 + \dots + \theta_n x_n.$$

- Parameters $\theta_1, \dots, \theta_n$ can be thought of as the components of the normal vector to the decision boundary.
 θ_0 can be thought of as the negative distance from decision boundary to origin.

2, We have 2 options

- One vs all ^{boundaries} where we model the decision boundary that separates each class against ~~the~~ all remaining classes
 \hookrightarrow This will produce k discriminant functions (one for each class)
- One vs each other where we separates each class against each of the remaining classes \rightarrow this will produce $\frac{k(k-1)}{2}$ discriminant functions, (one for each pair of classes)

→ One vs each other produces more discriminant functions and thus more parameters

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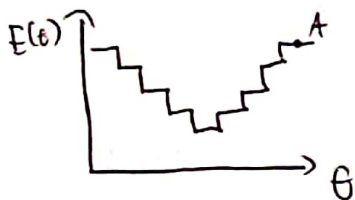
- Empirical Error:

$$E(\theta) = \#x^* + \#x^{**}$$

$$\text{where } x^* = \{x^{(i)} \mid 1(y^{(i)}=0) \wedge \theta^T x > 0\}$$

$$x^{**} = \{x^{(i)} \mid 1(y^{(i)}=1) \wedge \theta^T x < 0\}.$$

↳ The problem is that $E(\theta)$ is a piecewise function:



↳ We can't compute the gradient since we can easily stop at point A (gradient = 0).

→ Can't optimize using GD.

4,

$$PC: E(\theta) = \sum_{x^{(i)} \in x^*} \theta^T x^{(i)} - \sum_{x^{(i)} \in x^{**}} \theta^T x^{(i)}$$

$$\rightarrow \nabla E(\theta) = \sum_{x^{(i)} \in x^*} x^{(i)} - \sum_{x^{(i)} \in x^{**}} x^{(i)}$$

↳ Update:

$$\theta \leftarrow \theta - \eta \left(\sum_{x^{(i)} \in x^*} x^{(i)} - \sum_{x^{(i)} \in x^{**}} x^{(i)} \right)$$

5,

- In order to arrive at the logistic function as the hypothesis, we assume that the log of probabilities ratio can be modeled as a linear function:

$$\log \frac{P(y=1|x)}{P(y=0|x)} = \theta^T x.$$

- We simply equate $P(y=1|x) = \text{Sigmoid} = \frac{1}{1 + e^{-\theta^T x}}$

\Rightarrow the result is in the range of $(0, 1)$ where 0.5 represents complete uncertainty.

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$$\begin{aligned} l(\theta) &= \log \left(\prod_{x^{(i)} \in C_1} P(y=1|x^{(i)}) \prod_{x^{(i)} \in C_0} P(y=0|x^{(i)}) \right) \\ &= \log \prod_{i=1}^m P(y=1|x^{(i)})^{y^{(i)}} P(y=0|x^{(i)})^{1-y^{(i)}} \\ &= \sum_{i=1}^m y^{(i)} \log P(y=1|x^{(i)}) + (1-y^{(i)}) \log P(y=0|x^{(i)}) \\ &\quad \underbrace{\log P(y=0|x^{(i)})}_{\log(1 - P(y=1|x^{(i)}))} \end{aligned}$$

where $h_{\theta}(x) = P(y=1|x)$

$$\rightarrow l(\theta) = \sum_{i=1}^m y^{(i)} \log h_{\theta}(x^{(i)}) + (1-y^{(i)}) \log (1 - h_{\theta}(x^{(i)}))$$

$$\begin{aligned} - \frac{d}{d\theta} l(\theta) &= \frac{d}{d\theta} \sum_{i=1}^m y^{(i)} \log h_{\theta}(x^{(i)}) + (1-y^{(i)}) \log (1 - h_{\theta}(x^{(i)})) \\ &= \sum_{i=1}^m y^{(i)} (1 - h_{\theta}(x^{(i)})) x^{(i)} + (1-y^{(i)}) (-h_{\theta}(x^{(i)})) x^{(i)} \\ &= \sum_{i=1}^m x^{(i)} (y^{(i)} - y^{(i)} h_{\theta}(x^{(i)}) - h_{\theta}(x^{(i)}) + y^{(i)} h_{\theta}(x^{(i)})) \end{aligned}$$

$$= \sum_{i=1}^m (y^{(i)} - h_{\theta}(x^{(i)})) x^{(i)}$$

→ ~~negative log like~~ gradient of negative log likelihood:

$$\frac{d}{d\theta} (-\ell(\theta)) = \sum_{i=1}^m (h_{\theta}(x^{(i)}) - y^{(i)}) x^{(i)}$$

- SGD update: For batch X_k :

$$\theta \leftarrow \theta - \eta \sum_{i \in X_k} (h_{\theta}(x^{(i)}) - y^{(i)}) x^{(i)}$$

- Should initialize parameters randomly ~~& close to~~ close to 0

Since as $\theta \rightarrow 0$, $h_{\theta}(x) \rightarrow 0.5 \rightarrow$ represents ~~great~~ greatest uncertainty

7,

- For k classes, we use the softmax function since it outputs a distribution over all the ~~classes~~ classes and adds up to 1.

- Softmax is essentially a "one vs all" version of ~~logis~~ sigmoid.

- Update equation:

$$\theta_j \leftarrow \theta_j - \eta (h_{\theta_j}(x) - 1(y=j)) x^{(i)}$$

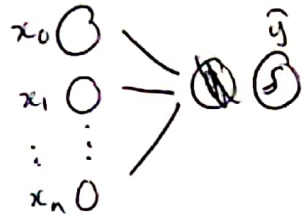
where $j=1 \dots k$ and θ_j = parameter vector for class j

⑤ Neural Net

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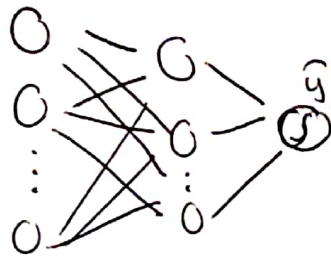
- Basic Structure:

A logistic unit:



→ We can stack multiple logistic unit on top of each other to form a neural net.

Ex:



- Activation function is important as it introduces non-linearity into the neural net model → give it ability to learn complex patterns.
- Several types of activation function: sigmoid, softmax, ReLU, leaky ReLU, linear (no activation), etc

2,

- In ~~FF~~ Feedforward net pushes the input ~~atth~~ forward all the way to output, Feedback net also does the forward push, but the output loops back to the input layer

- In the case of single output, we have exactly 1 parameter vector θ to plug into activation function

In K -class case, we have K parameter vectors $\theta_1 \dots \theta_K$

In terms of classification, we use sigmoid for single output and softmax for K -class outputs.

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⑥

- number of hidden units $<$ number of ~~inp~~ inputs :

→ We are basically doing dimensionality reduction

Since we are condensing a larger number of features into a smaller dimensional space

→ Extracting the more important features

- number of hidden units $>$ no of inputs :

→ We are doing non-linear mapping of features to higher dimensional space

→ Allows for more complex models.

4,

- Difficulty is that the update equations require knowing if $z^{(i)} = j$ or not where $z_j^{(i)}$ is the output of ~~hid~~ unit j in hidden layer.

- Chain rule is used to derive update equations by moving backward one layer at a time, calculating gradient at each step. Since

Since neural net is essentially a chain of functions, Consider

parameter v ~~into~~ one output unit \hat{y} with error E

$$\rightarrow \frac{\partial E}{\partial v} = \frac{\partial E}{\partial \hat{y}} \cdot \frac{\partial \hat{y}}{\partial v}$$

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5, Assume 2-layer

- Single output regression:

$$V = \begin{bmatrix} v_0 \\ \vdots \\ v_n \end{bmatrix} = \text{parameter vector of output layer}$$

$$W_j = \begin{bmatrix} w_{j0} \\ \vdots \\ w_{jn} \end{bmatrix} = \text{parameter vector of hidden unit } z_j.$$

$$E = \text{loss} = \frac{1}{2} \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)})^2$$

$$\begin{aligned} \hookrightarrow \frac{\partial E}{\partial v} &= \frac{\partial E}{\partial \hat{y}} \cdot \frac{\partial \hat{y}}{\partial v} = \frac{1}{2} \cdot 2 \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)}) \cdot z_j^{(i)} \\ &= \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)}) z_j^{(i)} \end{aligned}$$

+, Update for v :

$$v \leftarrow v - \eta \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)}) z_j^{(i)}$$

$$\begin{aligned} \frac{\partial E}{\partial w_j} &= \frac{\partial E}{\partial \hat{y}} \frac{\partial \hat{y}}{\partial z_j} \frac{\partial z_j}{\partial w_j} \\ &= \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)}) v_j z_j^{(i)} (1 - z_j^{(i)}) x^{(i)} \end{aligned}$$

+, Update for w_j :

$$w_j \leftarrow w_j - \eta \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)}) v_j z_j^{(i)} (1 - z_j^{(i)}) x^{(i)}$$

- Multiple outputs: assume k outputs

→ v_j = parameter vector of \hat{y}_j , $j = 1 \dots k$.

w_j = same definition as before.

$$E = \text{loss} = \frac{1}{2} \sum_{i=1}^m \sum_{j=1}^k (\hat{y}_j^{(i)} - y_j^{(i)})^2$$

↳ sum over all k classes.

$$+, \frac{\partial E}{\partial v_j} = \frac{\partial E}{\partial \hat{y}_j} \cdot \frac{\partial \hat{y}_j}{\partial v_j}$$

$$= \sum_{i=1}^m (\hat{y}_j^{(i)} - y_j^{(i)}) z^{(i)}$$

→ Update for v_j where $j = 1 \dots k$:

$$v_j \leftarrow v_j - \eta \sum_{i=1}^m (\hat{y}_j^{(i)} - y_j^{(i)}) z^{(i)}$$

$$+, \frac{\partial E}{\partial w_j} = \sum_{l=1}^k \frac{\partial E}{\partial \hat{y}_l} \cdot \frac{\partial \hat{y}_l}{\partial z_j} \cdot \frac{\partial z_j}{\partial w_j}$$

↳ summation because w_j leads to z_j which is input to all of k classes
Output units.

$$= \sum_{i=1}^m \sum_{l=1}^k (\hat{y}_l^{(i)} - y_l^{(i)}) v_{lj} z_j^{(i)} (1 - z_j^{(i)}) x^{(i)}$$

→ Update for w_{ji}

$$w_{ji} \leftarrow w_{ji} - \eta \sum_{i=1}^m \sum_{l=1}^k (\hat{y}_l^{(i)} - y_l^{(i)}) v_{lj} z_j^{(i)} (1 - z_j^{(i)}) x^{(i)}$$

- In case of regression, objective function is MSE, for multiclass, it's MSE over all classes.

7, Assume 2-layer:

- Single output classification:

Using the sigmoid derivative rules

$$\text{and } \frac{\partial E}{\partial v} = \frac{\partial E}{\partial \hat{y}} \cdot \frac{\partial \hat{y}}{\partial v}$$

$$\frac{\partial E}{\partial w_j} = \frac{\partial E}{\partial \hat{y}} \cdot \frac{\partial \hat{y}}{\partial z_j} \cdot \frac{\partial z_j}{\partial w_j}$$

$$\text{where } E = \sum_{i=1}^m y^{(i)} \log \hat{y}^{(i)} + (1 - y^{(i)}) \log (1 - \hat{y}^{(i)})$$

↳ We observe that $\frac{\partial E}{\partial v}$ and $\frac{\partial E}{\partial w_j}$ has same value as their regression counterpart.

$$\rightarrow v \leftarrow v - \eta \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)}) z^{(i)}$$

$$w_j \leftarrow w_j - \eta \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)}) v_j z_j^{(i)} (1 - z_j^{(i)}) x^{(i)}$$

- ~~the~~ K-class ~~the~~ classification:

Similarly, we also ~~ob~~ obtain same gradient and update equations as the regression ~~counter~~ counterpart ^{the} in case of multiple outputs.

- Objective function used is the cross-entropy / negative log likelihood obtained from negating the ~~likelihood~~ likelihood.

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- ~~Overfitt~~ Prevent overfitting by:
 - + Early stopping
 - + L1, L2 regularization ~~and~~
 - + Dropout.
- We can add regularization term to loss function:

Ex of 12:

$$E = \frac{1}{2} \sum_{i=1}^m (\hat{y}^{(i)} - y^{(i)})^2 + \lambda \| \theta \|^2$$

↳ update equation:

$$w_j \leftarrow w_j - \eta \frac{\partial E}{\partial w_j} - \lambda \| w_j \|^2$$

↳ Basically weight decay.

- Scaling input is important ~~some~~ since some input may have way bigger values than the other, which will have more influence on output.
↳ Even though ~~that may not~~ they may not be important features.
- One possible approach to adapt learning rate is learning rate decay, for ex:

$$\eta^{(i)} = \frac{\eta^{(i-1)}}{2}$$

↳ Useful since as we approach the minima, we ~~may~~ need smaller and smaller learning rate to avoid overshooting the minima.

- ~~Moment~~ Update equation with momentum added:

$$w_j^{(t+1)} \leftarrow w_j^{(t)} - \eta \frac{\partial E}{\partial w_j} + \beta (w_j^{(t)} - w_j^{(t-1)})$$

↳ useful to help us ~~avoid~~ overcome local minimum.

- Dropout is performed by randomly "shutdown" units in hidden layers during training

→ Prevent ~~co-adaptation~~ co-adaptation

9,

CNN is used to process image

The idea is that we arrange the parameters in an $n \times n$ matrix
(dot product)
called filter and we perform convolution between ~~the~~ the filter
and each $n \times n$ region of the image

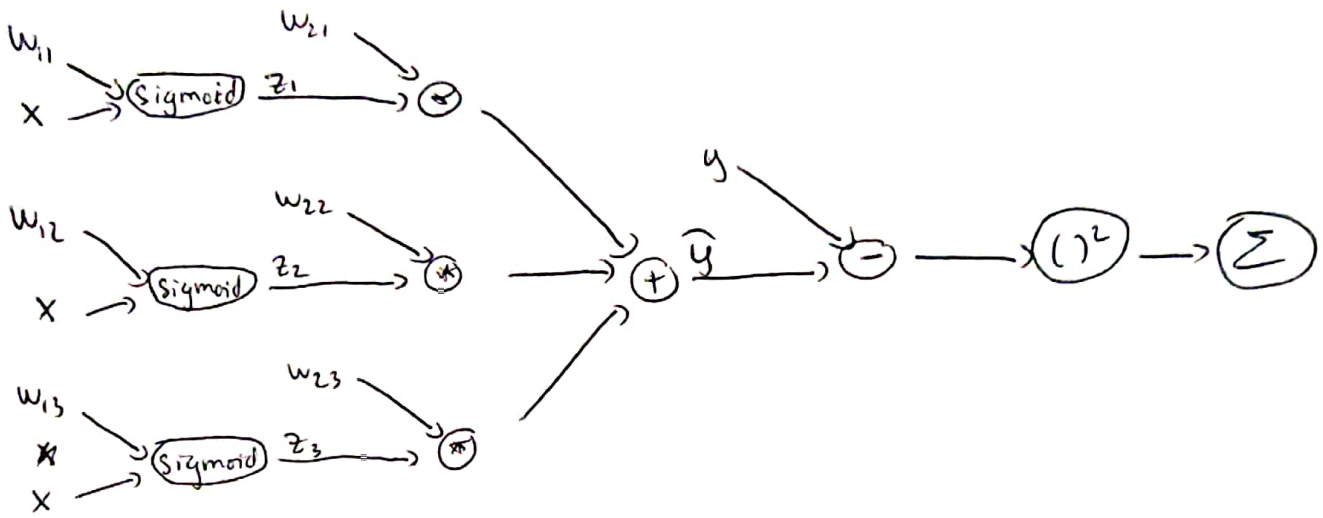
↳ ~~Advantages~~ Advantages of doing this include ~~we get~~:

+ We can keep the ~~image~~ input image in its original form
and not have to vectorize the image

+ There are ~~much~~ fewer parameters compared to vectorizing
the image and feed into MLP.

Deep learning:

1,



~~For x, y
 $w_{11} x = \text{Sigmoid}(w_{11}, x) z_1$~~