Logistic Regression

Machine Learning 10-601

Logistic Regression

Idea:

- Naïve Bayes allows computing P(Y|X) by learning P(Y) and P(X|Y)
 - Essentially learns P(Y)P(X|Y) = P(Y,X)
- Why not learn P(Y|X) directly?

- Consider learning f: X → Y, where
 - Problem set-up:
 - X is a vector of real-valued features, < X₁ ... X_n >
 - Y is boolean
 - Naïve Bayes assumption: assume all X_i are conditionally independent given Y
 - model $P(X_i \mid Y = y_k)$ as Gaussian $N(\mu_{ik}, \sigma_i)$
 - model P(Y) as Bernoulli (π)
- What does that imply about the form of P(Y|X)?

$$P(Y = 1|X = \langle X_1, ...X_n \rangle) = \frac{1}{1 + exp(w_0 + \sum_i w_i X_i)}$$

Derive form for P(Y|X) for continuous X_i

$$P(Y=1|X) = \frac{P(Y=1)P(X|Y=1)}{P(Y=1)P(X|Y=1) + P(Y=0)P(X|Y=0)}$$

$$= \frac{1}{1 + \frac{P(Y=0)P(X|Y=0)}{P(Y=1)P(X|Y=1)}}$$

$$= \frac{1}{1 + \exp(\ln\frac{P(Y=0)P(X|Y=0)}{P(Y=1)P(X|Y=1)})}$$

$$= \frac{1}{1 + \exp((\ln\frac{1-\pi}{\pi}) + \sum_{i} \ln\frac{P(X_{i}|Y=0)}{P(X_{i}|Y=1)})}$$

$$\sum_{i} \left(\frac{\mu_{i0} - \mu_{i1}}{\sigma_{i}^{2}} X_{i} + \frac{\mu_{i1}^{2} - \mu_{i0}^{2}}{2\sigma_{i}^{2}}\right)$$

$$P(Y=1|X) = \frac{1}{1 + \exp(w_{0} + \sum_{i=1}^{n} w_{i}X_{i})}$$

Very convenient!

$$P(Y = 1 | X = < X_1, ...X_n >) = \frac{1}{1 + exp(w_0 + \sum_i w_i X_i)}$$
 implies

$$P(Y = 0|X = < X_1, ...X_n >) = \frac{exp(w_0 + \sum_i w_i X_i)}{1 + exp(w_0 + \sum_i w_i X_i)}$$

implies

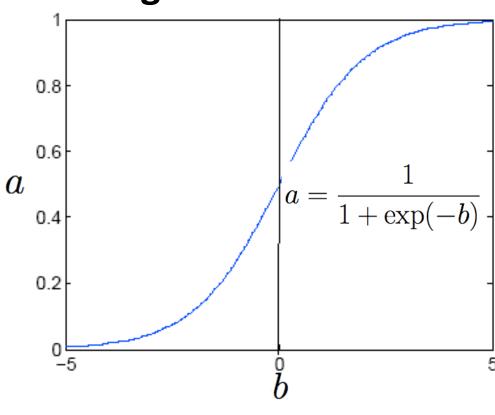
$$\frac{P(Y = 0|X)}{P(Y = 1|X)} = exp(w_0 + \sum_i w_i X_i)$$

∕linear classification rule!

implies

$$\ln \frac{P(Y=0|X)}{P(Y=1|X)} = w_0 + \sum_i w_i X_i$$

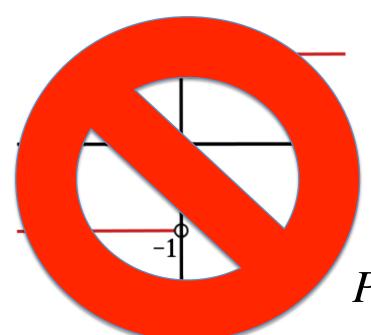
Logistic function



$$P(Y = 1|X) = \frac{1}{1 + \exp(w_0 + \sum_{i=1}^{n} w_i X_i)} -b$$

Logistic function for classifiers

 Replace sign(x•w) with something differentiable: e.g. the logistic(x•w)



logistic(
$$u$$
) $\equiv \frac{1}{1 + e^{-u}}$

$$P(Y = 1 \mid X = \mathbf{x}) \equiv \frac{1}{1 + e^{-\mathbf{x} \cdot \mathbf{w}}}$$

Logistic regression more generally

- Logistic regression when Y not boolean (but still discrete-valued).
- Now $y \in \{y_1 \dots y_R\}$: learn R-1 sets of weights

for
$$k < R$$
 $P(Y = y_k | X) = \frac{\exp(w_{k0} + \sum_{i=1}^n w_{ki} X_i)}{1 + \sum_{j=1}^{R-1} \exp(w_{j0} + \sum_{i=1}^n w_{ji} X_i)}$

for
$$k=R$$
 $P(Y = y_R|X) = \frac{1}{1 + \sum_{j=1}^{R-1} \exp(w_{j0} + \sum_{i=1}^{n} w_{ji}X_i)}$

Training Logistic Regression: Maximum Conditional Likelihood Estimation (MCLE)

- we have L training examples: $\{\langle X^1, Y^1 \rangle, \ldots \langle X^L, Y^L \rangle\}$
- maximum likelihood estimate for parameters W

$$W_{MLE} = \arg \max_{W} P(\langle X^{1}, Y^{1} \rangle \dots \langle X^{L}, Y^{L} \rangle | W)$$

= $\arg \max_{W} \prod_{l} P(\langle X^{l}, Y^{l} \rangle | W)$

maximum conditional likelihood estimate

Training Logistic Regression: MCLE

• Choose parameters $W=\langle w_0, ... w_n \rangle$ to <u>maximize</u> conditional likelihood of training data, where

$$P(Y = 0|X, W) = \frac{1}{1 + exp(w_0 + \sum_i w_i X_i)}$$
$$P(Y = 1|X, W) = \frac{exp(w_0 + \sum_i w_i X_i)}{1 + exp(w_0 + \sum_i w_i X_i)}$$

- Training data D = $\{\langle X^1, Y^1 \rangle, \dots \langle X^L, Y^L \rangle\}$
- Data likelihood = $\prod_{l} P(X^{l}, Y^{l}|W)$
- Data <u>conditional</u> likelihood = $\prod_{l} P(Y^{l}|X^{l}, W)$

$$W_{MCLE} = \arg\max_{W} \prod_{l} P(Y^{l}|W, X^{l})$$

Expressing Conditional Log Likelihood

$$l(W) \equiv \ln \prod_{l} P(Y^{l}|X^{l}, W) = \sum_{l} \ln P(Y^{l}|X^{l}, W)$$

$$P(Y = 0|X, W) = \frac{1}{1 + exp(w_{0} + \sum_{i} w_{i}X_{i})}$$

$$P(Y = 1|X, W) = \frac{exp(w_{0} + \sum_{i} w_{i}X_{i})}{1 + exp(w_{0} + \sum_{i} w_{i}X_{i})}$$

$$l(W) = \sum_{l} [Y^l \ln P(Y^l = 1|X^l, W) + (1 - Y^l) \ln P(Y^l = 0|X^l, W)]$$
 For the samples with $Y^l = 0$

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$$= \sum_{l} [Y^{l} \ln \frac{P(Y^{l} = 1 | X^{l}, W)}{P(Y^{l} = 0 | X^{l}, W)} + \ln P(Y^{l} = 0 | X^{l}, W)]$$

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$$= \sum_{l} [Y^{l} \ln \frac{P(Y^{l} = 1 | X^{l}, W)}{P(Y^{l} = 0 | X^{l}, W)} + \ln P(Y^{l} = 0 | X^{l}, W)]$$

$$= \sum_{l} [Y^{l} (w_{0} + \sum_{i}^{n} w_{i} X_{i}^{l}) - \ln(1 + exp(w_{0} + \sum_{i}^{n} w_{i} X_{i}^{l}))]$$

Maximizing Conditional Log Likelihood

$$P(Y = 0|X, W) = \frac{1}{1 + exp(w_0 + \sum_i w_i X_i)}$$
$$P(Y = 1|X, W) = \frac{exp(w_0 + \sum_i w_i X_i)}{1 + exp(w_0 + \sum_i w_i X_i)}$$

$$l(W) \equiv \ln \prod_{l} P(Y^{l}|X^{l}, W)$$

$$= \sum_{l} [Y^{l}(w_{0} + \sum_{i}^{n} w_{i}X_{i}^{l}) - \ln(1 + exp(w_{0} + \sum_{i}^{n} w_{i}X_{i}^{l}))]$$

Good news: l(W) is concave function of W

Bad news: no closed-form solution to maximize l(W)

Learning Logistic Regression with Gradient Descent

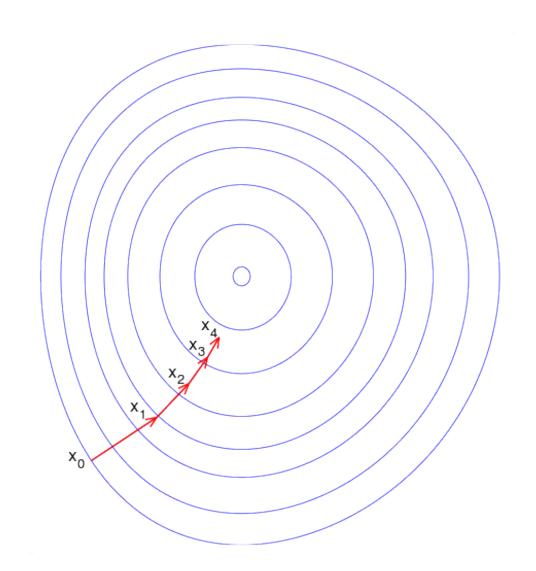
Learning as optimization: general procedure

- Goal: Learn the parameter w of ...
- Dataset: $D=\{(x_1,y_1),...,(x_n,y_n)\}$
- Write down a loss function
 - $Loss_D(\mathbf{w}) = \dots$
- Set w to minimize Loss
 - Usually we use numeric methods to find the optimum
 - i.e., gradient descent: repeatedly take a small step in the direction of the gradient

Gradient descent

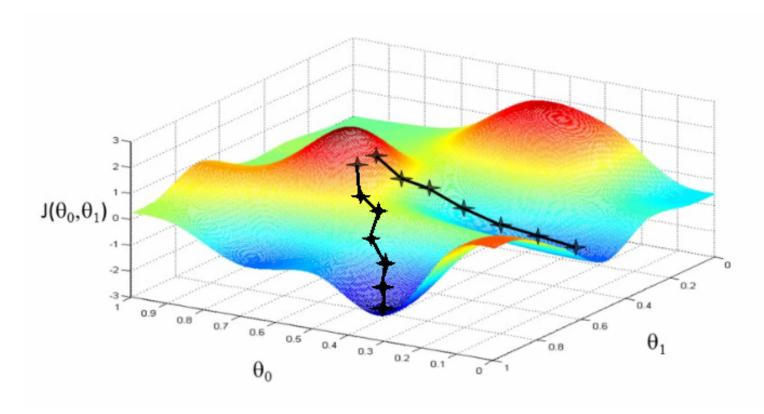
To find $\operatorname{argmin}_{\mathbf{x}} f(\mathbf{x})$:

- Start with \mathbf{x}_0
- For t=1....
 - $\mathbf{x}_{t+1} = \mathbf{x}_t \lambda f'(\mathbf{x}_t)$ where λ is small



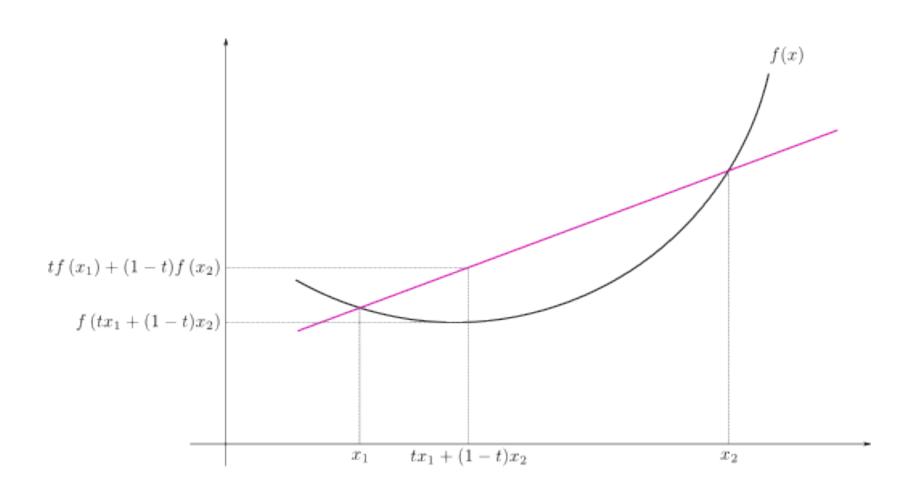
Pros and cons of gradient descent

- Simple and often quite effective on ML tasks
- Only applies to smooth functions (differentiable)
- Might find a local minimum, rather than a global one



Pros and cons of gradient descent

There is only one local optimum if the function is *convex*



Gradient Descent:

Batch gradient: use error $E_D(\mathbf{w})$ over entire training set D

Do until satisfied:

- 1. Compute the gradient $\nabla E_D(\mathbf{w}) = \left[\frac{\partial E_D(\mathbf{w})}{\partial w_0} \dots \frac{\partial E_D(\mathbf{w})}{\partial w_n} \right]$
- 2. Update the vector of parameters: $\mathbf{w} \leftarrow \mathbf{w} \eta \nabla E_D(\mathbf{w})$

Stochastic gradient: use error $E_d(\mathbf{w})$ over single examples $d \in D$

Do until satisfied:

- 1. Choose (with replacement) a random training example $d \in D$
- 2. Compute the gradient just for $d: \nabla E_d(\mathbf{w}) = \left[\frac{\partial E_d(\mathbf{w})}{\partial w_0} \dots \frac{\partial E_d(\mathbf{w})}{\partial w_n}\right]$
- 3. Update the vector of parameters: $\mathbf{w} \leftarrow \mathbf{w} \eta \nabla E_d(\mathbf{w})$

Stochastic approximates Batch arbitrarily closely as $\eta o 0$

Stochastic can be much faster when D is very large

Intermediate approach: use error over subsets of D

Maximize Conditional Log Likelihood: Gradient Ascent

$$l(W) \equiv \ln \prod_{l} P(Y^{l}|X^{l}, W)$$

$$= \sum_{l} Y^{l}(w_{0} + \sum_{i}^{n} w_{i}X_{i}^{l}) - \ln(1 + exp(w_{0} + \sum_{i}^{n} w_{i}X_{i}^{l}))$$

$$\frac{\partial l(W)}{\partial w_i} = \sum_{l} X_i^l (Y^l - \hat{P}(Y^l = 1|X^l, W))$$

Maximize Conditional Log Likelihood: Gradient Ascent

$$l(W) \equiv \ln \prod_{l} P(Y^{l}|X^{l}, W)$$

$$= \sum_{l} Y^{l}(w_{0} + \sum_{i}^{n} w_{i}X_{i}^{l}) - \ln(1 + exp(w_{0} + \sum_{i}^{n} w_{i}X_{i}^{l}))$$

$$\frac{\partial l(W)}{\partial w_i} = \sum_{l} X_i^l (Y^l - \hat{P}(Y^l = 1|X^l, W))$$

$$(\log f)' = \frac{1}{f}f'$$

$$(e^f)' = e^f f'$$

Gradient ascent algorithm: iterate until change $< \varepsilon$ For all i, repeat

$$w_i \leftarrow w_i + \eta \sum_{l} X_i^l (Y^l - \hat{P}(Y^l = 1 | X^l, W))$$

MAP Estimation with Regularization

That's all for M(C)LE. How about MAP?

MAP estimate

$$W \leftarrow \arg\max_{W} \ln P(W) \prod_{l} P(Y^{l}|X^{l}, W)$$

- One common approach is to define priors on W
 - Normal distribution, zero mean, identity covariance
- Helps avoid very large weights and overfitting

• let's assume Gaussian prior: W ~ N(0, σ^2 I) = 1/Z (w^j)⁻² (where Z is a constant)

MLE vs MAP

Maximum conditional likelihood estimate

$$W \leftarrow \arg\max_{W} \ \ln\prod_{l} P(Y^{l}|X^{l},W)$$

$$w_{i} \leftarrow w_{i} + \eta \sum_{l} X_{i}^{l} (Y^{l} - \widehat{P}(Y^{l} = 1|X^{l},W))$$

• Maximum a posteriori estimate with prior $W^{\sim}N(0,\sigma^{2}I)$

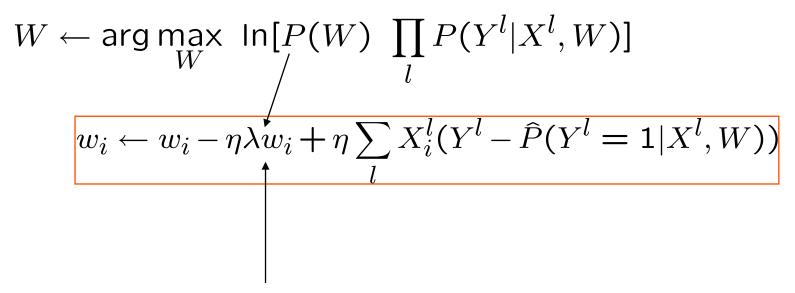
$$W \leftarrow \arg\max_{W} \ \ln[P(W) \ \prod_{l} P(Y^l|X^l,W)]$$

$$w_i \leftarrow w_i - \eta \lambda w_i + \eta \sum_{l} X_i^l (Y^l - \widehat{P}(Y^l = 1|X^l,W))$$

$$\lambda = 1/(2\sigma^2)$$

MAP Estimates and Regularization

• Maximum a posteriori estimate with prior $W^{N}(0,\sigma^{2}I)$



called a "regularization" term

- helps reduce overfitting, especially when training data is sparse
- keep weights nearer to zero (if P(W) is zero mean Gaussian prior), or whatever the prior suggests
- used very frequently in Logistic Regression

The Bottom Line

- Consider learning f: $X \rightarrow Y$, where
 - X is a vector of real-valued features, $< X_1 ... X_n >$
 - Y is boolean
- assume all X_i are conditionally independent given Y
 - model $P(X_i \mid Y = y_k)$ as Gaussian $N(\mu_{ik}, \sigma_i)$
 - model P(Y) as Bernoulli (π)
- Then P(Y|X) is of this form, and we can directly estimate W

$$P(Y = 1|X = \langle X_1, ...X_n \rangle) = \frac{1}{1 + exp(w_0 + \sum_i w_i X_i)}$$



Generative vs. Discriminative Classifiers

Training classifiers involves estimating f: $X \rightarrow Y$, or P(Y|X)

Generative classifiers (e.g., Naïve Bayes)

- Assume some functional form for P(X|Y), P(X) (i.e., P(X,Y))
- Estimate parameters of P(X|Y), P(X) directly from training data
- Use Bayes rule to calculate P(Y|X=x_i)
 - Find θ = argmax $_{\mathbf{w}} \Pi_{\mathbf{i}} \Pr(y_{\mathbf{i}}, x_{\mathbf{i}} | \theta)$
 - Different assumptions about *generative process* for the data: Pr(X,Y), priors on θ ,...

Discriminative classifiers (e.g., Logistic regression)

- Assume some functional form for P(Y|X)
- Estimate parameters of P(Y|X) directly from training data
 - Find θ = argmax $_{\mathbf{w}} \Pi_{\mathbf{i}} \Pr(y_{\mathbf{i}} | x_{\mathbf{i}}, \theta)$
 - Different assumptions about conditional probability: Pr(Y|X), priors on θ , ...

Use Naïve Bayes or Logisitic Regression?

Consider

- Restrictiveness of modeling assumptions
- Rate of convergence (in amount of training data) toward asymptotic hypothesis

Consider Y boolean, X_i continuous, $X = \langle X_1 ... X_n \rangle$

Number of parameters:

• NB: 4n + 1 (3n + 1 if we assume $\sigma_{ik} = \sigma_i$)

$$P(Y = 0|X, W) = \frac{1}{1 + exp(w_0 + \sum_{i} w_i X_i)}$$

$$P(Y = 1|X, W) = \frac{exp(w_0 + \sum_i w_i X_i)}{1 + exp(w_0 + \sum_i w_i X_i)}$$

Estimation method:

- NB parameter estimates are uncoupled
- LR parameter estimates are coupled

[Ng & Jordan, 2002]

Recall two assumptions deriving form of LR from GNBayes:

- 1. X_i conditionally independent of X_k given Y
- 2. $P(X_i \mid Y = y_k) = N(\mu_{ik}, \sigma_i), \leftarrow \text{not } N(\mu_{ik}, \sigma_{ik})$

Consider three learning methods:

- •GNB (assumption 1 only) -- decision surface can be non-linear
- •GNB2 (assumption 1 and 2) decision surface linear
- •LR -- decision surface linear, trained differently

- •Both (1) and (2) are satisfied:
- •Neither (1) nor (2) is satisfied:
- •(1) is satisfied, but not (2):

[Ng & Jordan, 2002]

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- •Neither (1) nor (2) is satisfied: LR > GNB2, GNB>GNB2
- •(1) is satisfied, but not (2):

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- •Both (1) and (2) are satisfied: LR = GNB2 = GNB
- •Neither (1) nor (2) is satisfied: LR > GNB2, GNB>GNB2
- •(1) is satisfied, but not (2): GNB > LR

[Ng & Jordan, 2002]

What if we have only finite training data?

They converge at different rates to their asymptotic (∞ data) error

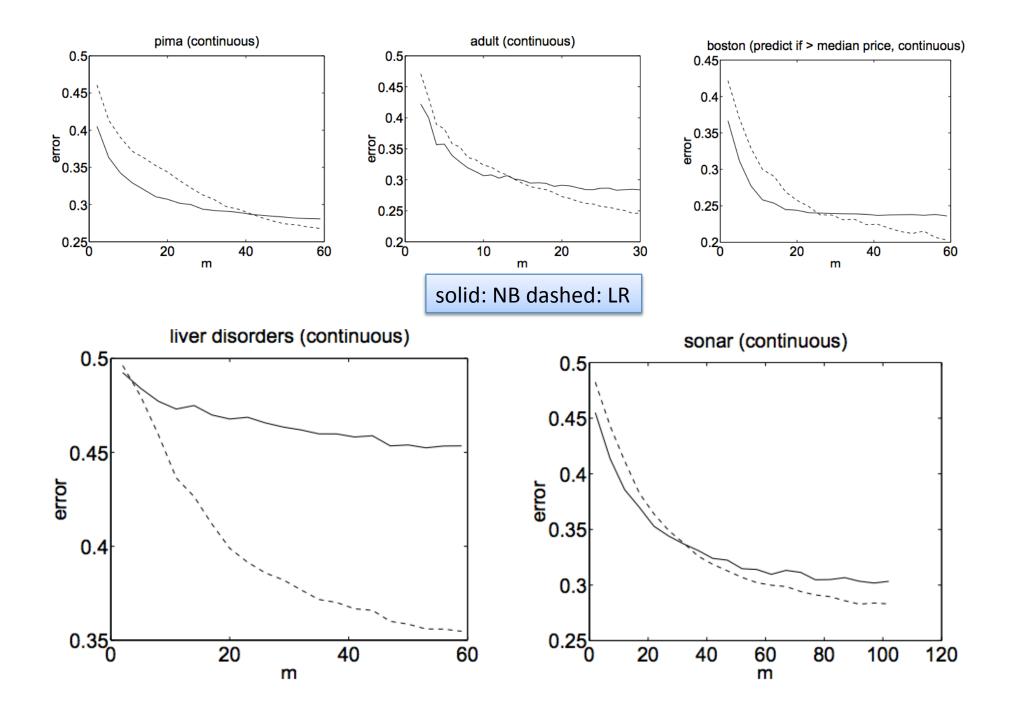
Let $\epsilon_{A,n}$ refer to expected error of learning algorithm A after n training examples

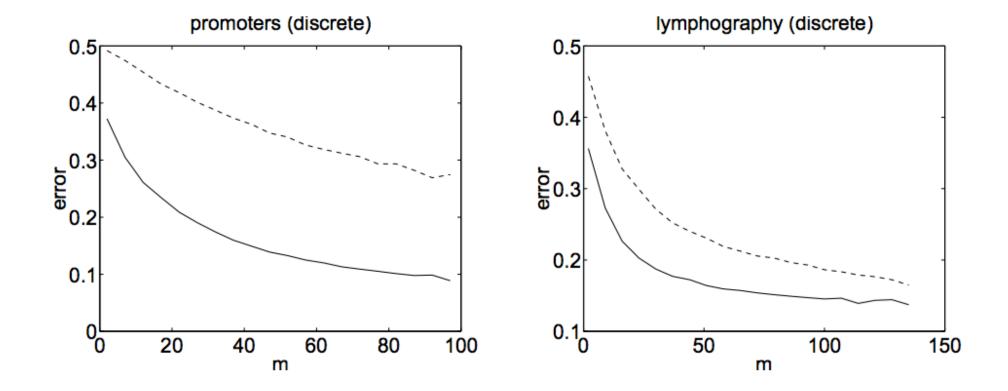
Let d be the number of features: $\langle X_1 ... X_d \rangle$

$$\epsilon_{LR,n} \le \epsilon_{LR,\infty} + O\left(\sqrt{\frac{d}{n}}\right)$$

$$\epsilon_{GNB,n} \le \epsilon_{GNB,\infty} + O\left(\sqrt{\frac{\log d}{n}}\right)$$

So, GNB requires $n = O(\log d)$ to converge, but LR requires n = O(d)





Naïve Bayes makes stronger assumptions about the data but needs fewer examples to estimate the parameters

"On Discriminative vs Generative Classifiers:" Andrew Ng and Michael Jordan, NIPS 2001.

The bottom line:

GNB2 and LR both use linear decision surfaces, GNB need not

Given infinite data, LR is better or equal to GNB2 because *training procedure* does not make assumptions 1 or 2 (though our derivation of the form of P(Y|X) did).

But GNB2 converges more quickly to its perhaps-less-accurate asymptotic error

And GNB is both more biased (assumption1) and less (no assumption 2) than LR, so either might beat the other

Measuring Accuracy of Classifier

- Precision = #(classified as positive AND positive in data)
 #(classified as positive)
 e.g., how many of the emails classified as "spam" are in fact truly "spam"?
- Recall = #(classified as positive AND positive in data)
 #(positive in data)

e.g., how many of the "spam" emails were classified as "spam"?

What you should know:

- Logistic regression
 - Functional form follows from Naïve Bayes assumptions
 - For Gaussian Naïve Bayes assuming variance $\sigma_{i,k} = \sigma_i$
 - For discrete-valued Naïve Bayes too
 - But training procedure picks parameters without making conditional independence assumption
 - MLE training: pick W to maximize P(Y | X, W)
 - MAP training: pick W to maximize P(W | X,Y)
 - 'regularization'
 - helps reduce overfitting
- Gradient ascent/descent
 - General approach when closed-form solutions unavailable
- Generative vs. Discriminative classifiers