[70240413 Statistical Machine Learning, Spring, 2019]

Probabilistic Methods for Classification

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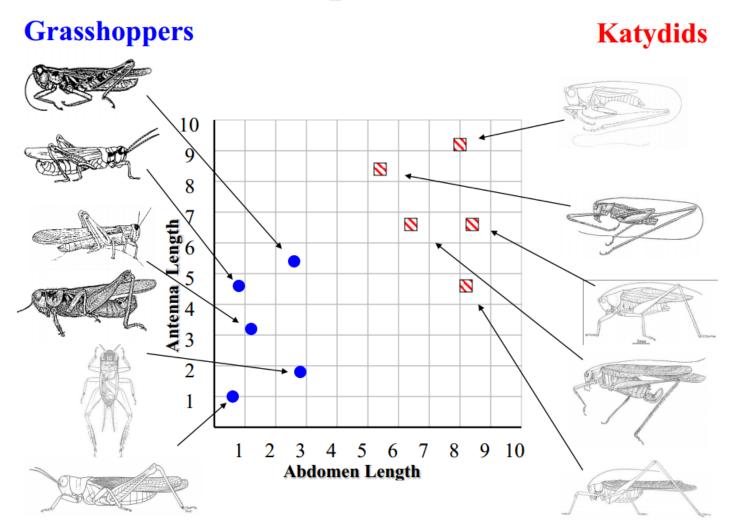
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Outline

- Probabilistic methods for supervised learning
- Naive Bayes classifier
- Logistic regression
- Exponential family distributions
- Generalized linear models

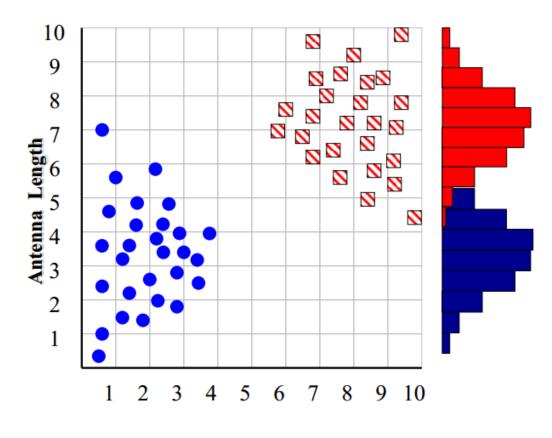
An Intuitive Example



[Courtesy of E. Keogh]

With more data ...

Build a histogram, e.g., for "Antenna length"

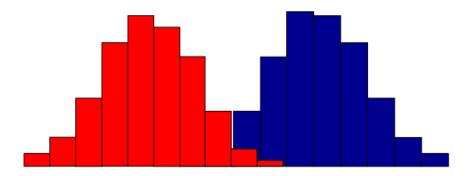


⊠ Katydids

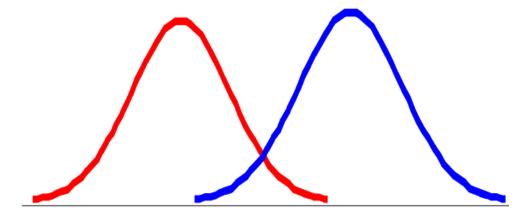
Grasshoppers

Empirical distribution

Histogram (or empirical distribution)

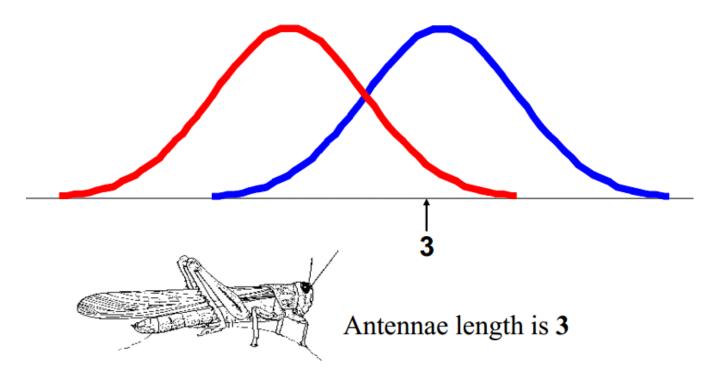


Smooth with kernel density estimation (KDE):



Classification?

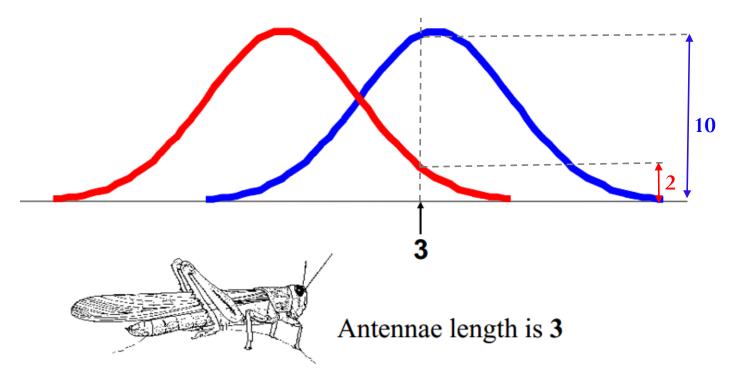
- Classify another insect we find. Its antennae are 3 units long
- Is it more probable that the insect is a Grasshopper or a Katydid?



Classification Probability

$$P(Grasshopper | 3) = 10 / (10 + 2) = 0.833$$

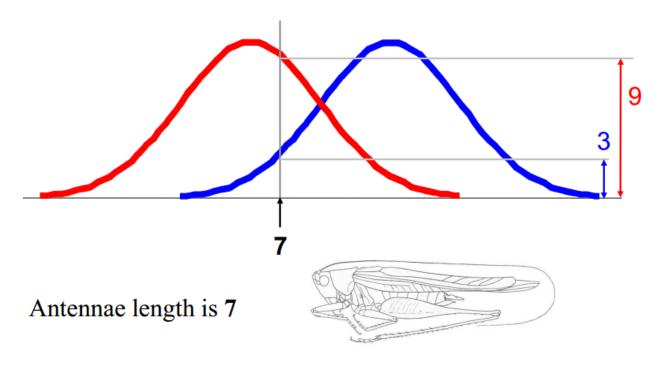
 $P(Katydid | 3) = 2 / (10 + 2) = 0.166$



Classification Probability

$$P(Grasshopper | 7) = 3 / (3 + 9) = 0.250$$

 $P(Katydid | 7) = 9 / (3 + 9) = 0.750$

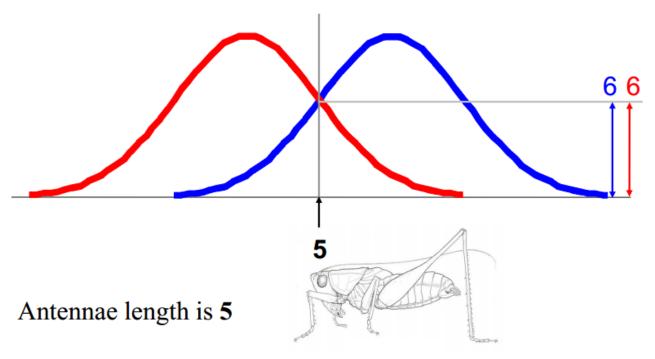


[Courtesy of E. Keogh]

Classification Probability

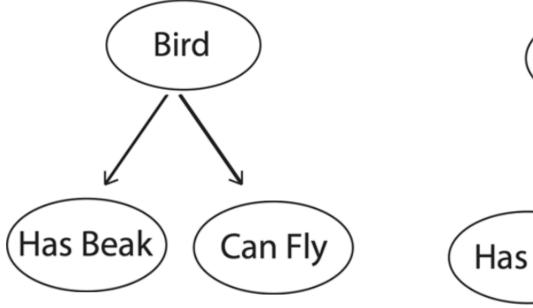
$$P(Grasshopper | 5) = 6 / (6 + 6) = 0.500$$

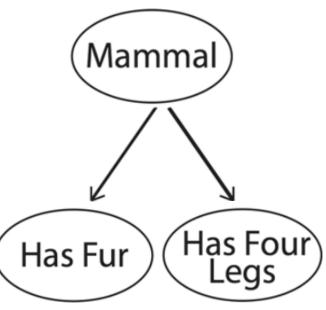
 $P(Katydid | 5) = 6 / (6 + 6) = 0.500$



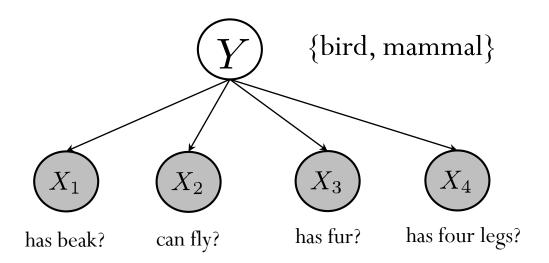
[Courtesy of E. Keogh]

- The simplest "category-feature" generative model:
 - Category: "bird", "Mammal"
 - □ **Features**: "has beak", "can fly" ...





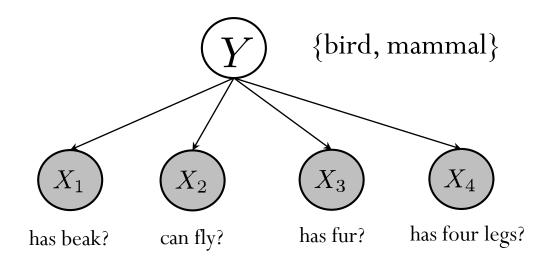
- A mathematic model:
 - **Naive Bayes assumption**: features X_1, \ldots, X_d are conditionally independent given the class label Y



A joint distribution:

tion: prior likelihood
$$p(\mathrm{x},y) = p(y)p(\mathrm{x}|y)$$

♦ A mathematic model:



Inference via Bayes rule:

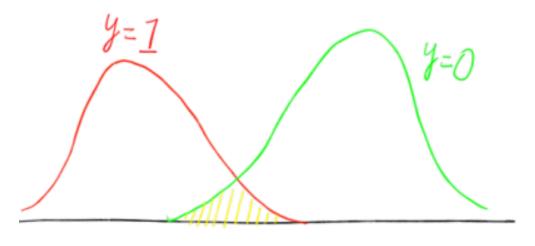
$$p(y|\mathbf{x}) = \frac{p(\mathbf{x}, y)}{p(\mathbf{x})} = \frac{p(y)p(\mathbf{x}|y)}{p(\mathbf{x})}$$

Bayes' decision rule:

$$y^* = \arg\max_{y \in \mathcal{Y}} p(y|\mathbf{x})$$

Bayes Error

Theorem: Bayes classifier is optimal!



$$p(error|\mathbf{x}) = \begin{cases} p(y=1|\mathbf{x}) & \text{if we decide } y=0\\ p(y=0|\mathbf{x}) & \text{if we decide } y=1 \end{cases}$$

$$p(error) = \int_{-\infty}^{\infty} p(error|\mathbf{x})p(\mathbf{x})d\mathbf{x}$$

• However, the true distribution is unknown.

Learning!

□ We need to estimate it!

- How to learn model parameters?
 - Assume *X* are *d* binary features, *Y* has 2 possible labels

$$p(y|\pi) = \left\{ \begin{array}{ll} \pi & \text{if } y = 1 \; (i.e., \; \text{bird}) \\ 1 - \pi & \text{otherwise} \end{array} \right. \quad \left\{ \begin{array}{ll} \text{bird, mammal} \right\} \\ X_1 & X_2 & X_3 & X_4 \\ \text{has beak?} & \text{can fly?} & \text{has fur?} & \text{has four legs?} \end{array} \right.$$

$$p(x_j|y=0,q) = \begin{cases} q_{0j} & \text{if } x_j = 1\\ 1 - q_{0j} & \text{otherwise} \end{cases}$$
 $p(x_j|y=1,q) = \begin{cases} q_{1j} & \text{if } x_j = 1\\ 1 - q_{1j} & \text{otherwise} \end{cases}$

■ How many parameters to estimate?

- How to learn model parameters?
- A set of training data:
 - \circ (1, 1, 0, 0; 1)
 - (1,0,0,0;1)
 - (0, 1, 1, 0; 0)
 - (0, 0, 1, 1; 0)
- **♦ Maximum likelihood estimation** (*N*: # of training data)

$$p(\{\mathbf{x}_i, y_i | \pi, q\}) = \prod_{i=1}^{N} p(\mathbf{x}_i, y_i | \pi, q)$$

♦ Maximum likelihood estimation (*N*: # of training data)

$$(\hat{\pi}, \hat{q}) = \arg \max_{\pi, q} p(\{\mathbf{x}_i, y_i\} | \pi, q)$$

$$(\hat{\pi}, \hat{q}) = \arg\max_{\pi, q} \log p(\{\mathbf{x}_i, y_i\} | \pi, q)$$

Results (count frequency! Exercise?):

$$\hat{\pi} = \frac{N_1}{N}$$
 $\hat{q}_{0j} = \frac{N_0^j}{N_0}$ $\hat{q}_{1j} = \frac{N_1^j}{N_1}$

$$N_k = \sum_{i=1}^{N} \mathbf{I}(y_i = k)$$
: # of data in category k

$$N_k^j = \sum_{i=1}^{N} \mathbf{I}(y_i = k, x_{ij} = 1)$$
: # of data in category k that has feature j

Data scarcity issue (zero-counts problem):

$$\hat{\pi} = \frac{N_1}{N}$$
 $\hat{q}_{0j} = \frac{N_0^j}{N_0}$ $\hat{q}_{1j} = \frac{N_1^j}{N_1}$

- How about if some features do not appear?
- Laplace smoothing (Additive smoothing):

$$\hat{q}_{0j} = \frac{N_0^j + \alpha}{N_0 + 2\alpha}$$

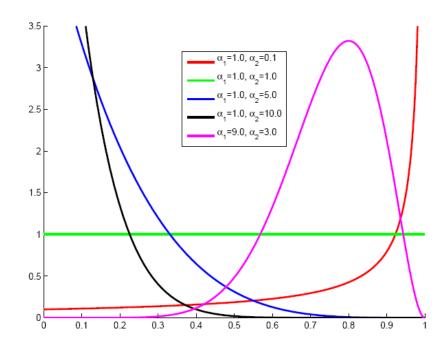
$$\alpha > 0$$

$$\hat{q}_{1j} = \frac{N_1^j + \alpha}{N_1 + 2\alpha}$$

A Bayesian Treatment

Put a prior on the parameters

$$p_0(q_{0j}|\alpha_1,\alpha_2) = \text{Beta}(\alpha_1,\alpha_2) = \frac{\Gamma(\alpha_1 + \alpha_2)}{\Gamma(\alpha_1)\Gamma(\alpha_2)} q_{0j}^{\alpha_1 - 1} (1 - q_{0j})^{\alpha_2 - 1}$$



A Bayesian Treatment

Maximum a Posterior Estimate (MAP):

$$\hat{q} = \arg \max_{q} \log p(q | \{\mathbf{x}_i, y_i\})$$

$$= \arg \max_{q} \log p(q) + \log p(\{\mathbf{x}_i, y_i\} | q)$$

Results (Exercise?):

$$\hat{q}_{0j} = \frac{N_0^j + \alpha_1 - 1}{N_0 + \alpha_1 + \alpha_2 - 2}$$

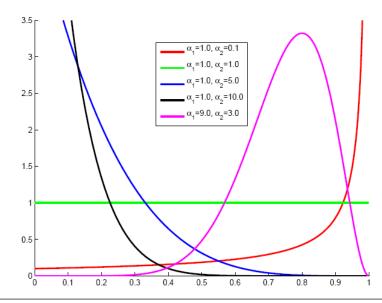
$$\hat{q}_{1j} = \frac{N_1^j + \alpha_1 - 1}{N_1 + \alpha_1 + \alpha_2 - 2}$$

A Bayesian Treatment

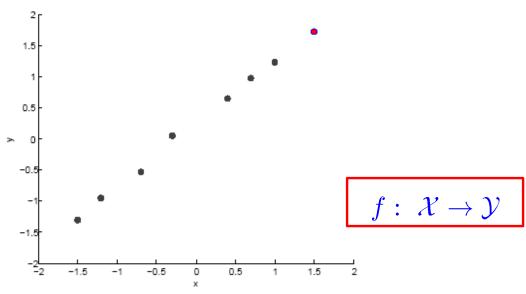
Maximum a Posterior Estimate (MAP):

$$\hat{q}_{0j} = \frac{N_0^j + \alpha_1 - 1}{N_0 + \alpha_1 + \alpha_2 - 2}$$

- If $\alpha_1 = \alpha_2 = 1$ (non-informative prior), no effect
 - MLE is a special case of Bayesian estimate
- \bullet Increase α_1, α_2 , lead to heavier influence from prior



Bayesian Regression



• Goal: learn a function from noisy observed data

Linear
$$\mathcal{F}_{linear} = \{f: f = wx + b, w, b \in \mathbb{R}\}$$

$$\begin{array}{ll} & \text{Polynomial} & \mathcal{F}_{polynomial} = \{f: \ f = \sum_k w_k x^k, \ w_k \in \mathbb{R}\} \\ & \dots \end{array}$$

•

Bayesian Regression

Noisy observations

$$y = f(\mathbf{x}) + \epsilon$$
, where $\epsilon \sim \mathcal{N}(0, \sigma_n^2)$

lacktriangle Gaussian likelihood function for linear regression $f(\mathbf{x}_i) = \mathbf{w}^{\top} \mathbf{x}_i$

$$p(\mathbf{y}|X, \mathbf{w}) = \prod_{i=1}^{N} p(y_i|\mathbf{x}_i, \mathbf{w}) = \mathcal{N}(X^{\top}\mathbf{w}, \sigma_n^2 I)$$

Gaussian prior (Conjugate)

$$\mathbf{w} \sim \mathcal{N}(0, \Sigma_d)$$

- Inference with Bayes' rule
 - Posterior $p(\mathbf{w}|X,\mathbf{y}) = \mathcal{N}(\frac{1}{\sigma_n^2}A^{-1}X\mathbf{y}, A^{-1}), \text{ where } A = \sigma_n^{-2}XX^\top + \Sigma_d^{-1}$
 - Marginal likelihood
 - Prediction $p(\mathbf{y}|X) = \int p(\mathbf{y}|X, \mathbf{w}) p(\mathbf{w}) d\mathbf{w}$

$$p(f_*|\mathbf{x}_*, X, \mathbf{y}) = \int p(f_*|\mathbf{x}_*, \mathbf{w}) p(\mathbf{w}|X, \mathbf{y}) d\mathbf{w} = \mathcal{N}\left(\frac{1}{\sigma_n^2} \mathbf{x}_*^\top A^{-1} X \mathbf{y}, \mathbf{x}_*^\top A^{-1} \mathbf{x}_*\right)$$

Extensions of NB

- We covered the case with binary features and binary class labels
- NB is applicable to the cases:
 - Discrete features + discrete class labels
 - Continuous features + discrete class labels
 - **.** . . .
- More dependency between features can be considered
 - Tree augmented NB
 - **...**

Gaussian Naive Bayes (GNB)

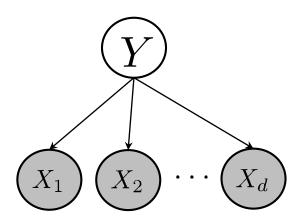
- \bullet E.g.: character recognition: feature X_i is intensity at pixel i:
- The generative process is

$$Y \sim \text{Bernoulli}(\pi)$$

$$P(X_i|Y=y) = \mathcal{N}(\mu_{iy}, \sigma_{iy}^2)$$



- Different mean and variance for each class k and each feature i
- Sometimes assume variance is:
 - independent of Y (i.e., σ_i)
 - ullet or independent of X (i.e., σ_y)
 - $lue{}$ or both (i.e., σ)



Estimating Parameters & Prediction

MLE estimates

$$\hat{\mu}_{ik} = \frac{1}{\sum_{n} \mathbb{I}(y_n = k)} \sum_{n} x_{ni} \mathbb{I}(y_n = k)$$
pixel i in training image n

$$\hat{\sigma}_{ik}^{2} = \frac{1}{\sum_{n} \mathbb{I}(y_{n} = k)} \sum_{n} (x_{ni} - \hat{\mu}_{ik})^{2} \mathbb{I}(y_{n} = k)$$

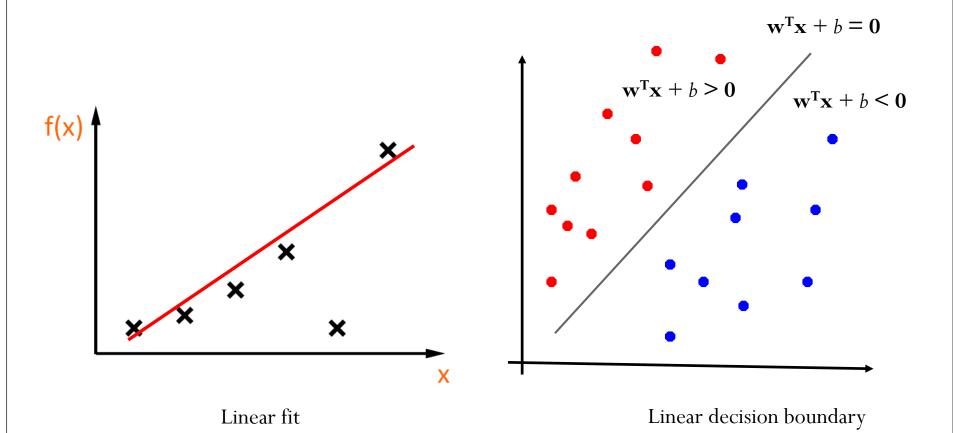
Prediction:

$$h(\mathbf{x}) = \underset{y}{\operatorname{argmax}} P(y) \prod_{i} P(x_i|y)$$

What you need to know about NB classifier

- What's the assumption
- Why we use it
- How do we learn it
- Why is Bayesian estimation (MAP) important

Linear regression and linear classification



What's the decision boundary of NB?

- Is it linear or non-linear?
- There are several distributions that lead to a linear decision boundary, e.g., GNB with equal variance

$$P(X_i|Y=y) = \mathcal{N}(\mu_{iy}, \sigma_i^2)$$

Decision boundary (??):

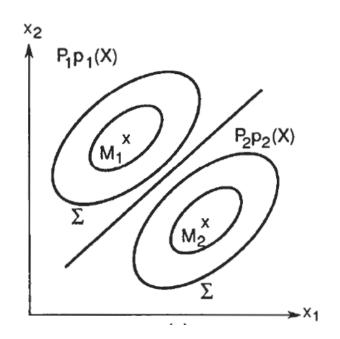
$$\log \frac{\prod_{i=1}^{a} P(X_i|Y=0)P(Y=0)}{\prod_{i=1}^{d} P(X_i|Y=1)P(Y=1)} = 0$$

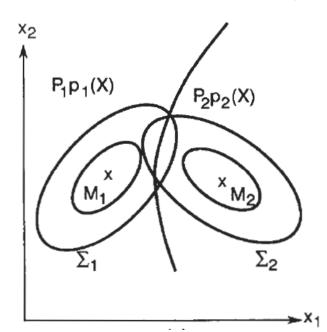
$$\Rightarrow \log \frac{1-\pi}{\pi} + \sum_{i} \frac{\mu_{i1}^2 - \mu_{i0}^2}{2\sigma_i^2} + \sum_{i} \frac{\mu_{i0} - \mu_{i1}}{\sigma_i^2} x_i = 0$$

$$\Rightarrow w_0 + \sum_{i} w_i x_i = 0$$

Gaussian Naive Bayes (GNB)

Decision boundary (the general multivariate Gaussian case):





$$P_1 = P(Y = 0), \quad P_2 = P(Y = 1)$$

 $p_1(X) = p(X|Y = 0) = \mathcal{N}(M_1, \Sigma_1)$
 $p_2(X) = p(X|Y = 1) = \mathcal{N}(M_2, \Sigma_2)$

The predictive distribution of GNB

• Understanding the predictive distribution

$$p(y = 1 | \mathbf{x}, \mu, \Sigma, \pi) = \frac{p(y = 1, \mathbf{x} | \mu, \Sigma, \pi)}{p(\mathbf{x} | \mu, \Sigma, \pi)}$$

• Under naive Bayes assumption:

$$p(y = 1 | \mathbf{x}, \mu, \Sigma, \pi) = \frac{1}{1 + \frac{p(y = 0, \mathbf{x} | \mu, \Sigma, \pi)}{p(y = 1, \mathbf{x} | \mu, \Sigma, \pi)}}$$

$$= \frac{1}{1 + \frac{(1 - \pi) \prod_{i} \mathcal{N}(x_{i} | \mu_{i0}, \sigma_{i}^{2})}{\pi \prod_{i} \mathcal{N}(x_{i} | \mu_{i1}, \sigma_{i}^{2})}}$$

$$= \frac{1}{1 + \exp(-\mathbf{w}^{\top}\mathbf{x} - w_{0})}$$

♦ Note: For multi-class, the predictive distribution is softmax!

Generative vs. Discriminative Classifiers

- ♦ Generative classifiers (e.g., Naive Bayes)
 - Assume some functional form for P(X,Y) (or P(Y) and P(X|Y))
 - Estimate parameters of P(X,Y) directly from training data
 - Make prediction

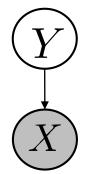
$$\hat{y} = \operatorname*{argmax}_{y} P(\mathbf{x}, Y = y)$$

But, we note that

$$\hat{y} = \operatorname*{argmax}_{y} P(Y = y | \mathbf{x})$$



- Discriminative classifiers (e.g., Logistic regression)
 - $lue{}$ Assume some functional form for P(Y | X)
 - ullet Estimate parameters of P(Y | X) directly from training data



Logistic Regression

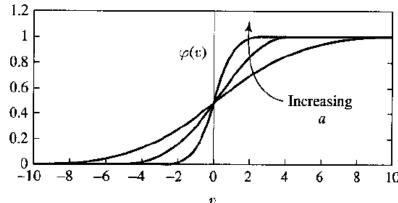
- Recall the predictive distribution of GNB!
- \diamond Assume the following functional form for P(Y | X)

$$P(y = 1|\mathbf{x}) = \frac{1}{1 + \exp(-(w_0 + \mathbf{w}^{\mathsf{T}}\mathbf{x}))}$$

Logistic function (or Sigmoid) applied to a linear function of the data (for $\alpha = 1$):

$$\psi_{\alpha}(v) = \frac{1}{1 + \exp(-\alpha v)}$$

 $a \to \infty$: step function



use a large α can be good for some neural networks

Logistic Regression

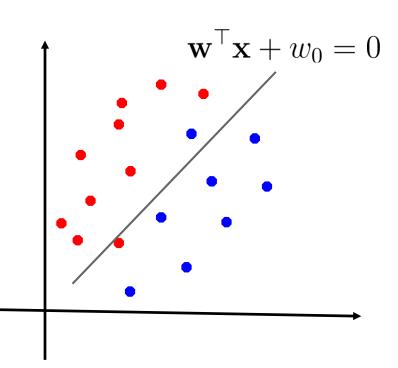
• What's the decision boundary of logistic regression? (linear or nonlinear?)

$$P(y = 1|\mathbf{x}) = \frac{1}{1 + \exp(-(w_0 + \mathbf{w}^{\top}\mathbf{x}))}$$

$$\log \frac{P(Y=1|\mathbf{x})}{P(y=0|\mathbf{x})} = 0$$

$$\mathbf{w}^{\top}\mathbf{x} + w_0 = 0$$

Logistic regression is a linear classifier!



Representation

Logistic regression

$$P(y = 1|\mathbf{x}) = \frac{1}{1 + \exp(-(w_0 + \mathbf{w}^{\top}\mathbf{x}))}$$

For notation simplicity, we use the augmented vector:

input features:
$$\begin{pmatrix} 1 \\ \mathbf{x} \end{pmatrix}$$
 model weights: $\begin{pmatrix} w_0 \\ \mathbf{w} \end{pmatrix}$

□ Then, we have

$$P(y = 1|\mathbf{x}) = \frac{1}{1 + \exp(-\mathbf{w}^{\top}\mathbf{x})}$$

Multiclass Logistic Regression

 \bullet For more than 2 classes, where $y \in \{1, \dots, K\}$, logistic regression classifier is defined as

$$\forall k < K : P(Y = k | \mathbf{x}) = \frac{\exp(\mathbf{w}_k^{\top} \mathbf{x})}{1 + \sum_{j=1}^{K-1} \exp(\mathbf{w}_j^{\top} \mathbf{x})}$$
$$P(Y = K | \mathbf{x}) = \frac{1}{1 + \sum_{j=1}^{K-1} \exp(\mathbf{w}_j^{\top} \mathbf{x})}$$

- Well normalized distribution! No weights for class K!
- Is the decision boundary still linear?

Training Logistic Regression

We consider the binary classification

$$P(y = 1|\mathbf{x}) = \frac{1}{1 + \exp(-\mathbf{w}^{\top}\mathbf{x})}$$

- Training data $D = \{(\mathbf{x}_i, y_i)\}_{i=1}^N$
- How to learn the parameters?
- Can we do MLE?

$$\hat{\mathbf{w}}_{MLE} = \underset{\mathbf{w}}{\operatorname{argmax}} \prod_{i=1}^{N} P(\mathbf{x}_i, y_i | \mathbf{w})$$

- No! Don't have a model for P(X) or P(X | Y)
- Can we do large-margin learning?

Maximum Conditional Likelihood Estimate

• We learn the parameters by solving

$$\hat{\mathbf{w}} = \underset{\mathbf{w}}{\operatorname{argmax}} \prod_{i=1}^{N} P(y_i | \mathbf{x}_i, \mathbf{w})$$

Discriminative philosophy – don't waste effort on learning P(X), focus on P(Y | X) – that's all that matters for classification!

Maximum Conditional Likelihood Estimate

$$\hat{\mathbf{w}} = \underset{\mathbf{w}}{\operatorname{argmax}} \prod_{i=1}^{N} P(y_i | \mathbf{x}_i, \mathbf{w})$$

$$P(y = 1|\mathbf{x}) = \frac{1}{1 + \exp(-\mathbf{w}^{\top}\mathbf{x})}$$

• We have:

$$\mathcal{L}(\mathbf{w}) = \log \prod_{i=1}^{N} P(y_i | \mathbf{x}_i, \mathbf{w})$$
$$= \sum_{i} \left[y_i \mathbf{w}^{\top} \mathbf{x}_i - \log(1 + \exp(\mathbf{w}^{\top} \mathbf{x}_i)) \right]$$

Maximum Conditional Likelihood Estimate

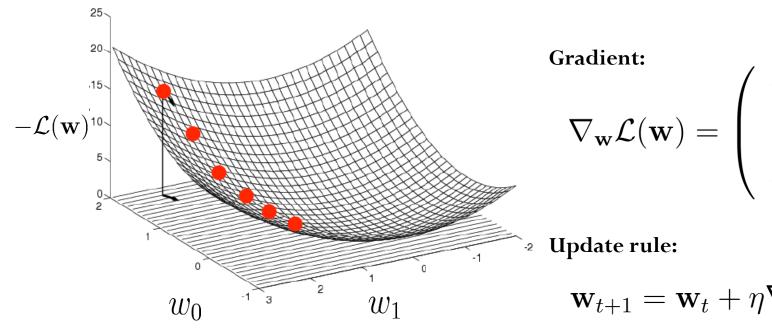
$$\hat{\mathbf{w}} = \operatorname*{argmax}_{\mathbf{w}} \mathcal{L}(\mathbf{w})$$

$$\mathbf{\mathcal{L}}(\mathbf{w}) = \sum_{i} \left[y_{i} \mathbf{w}^{\top} \mathbf{x}_{i} - \log(1 + \exp(\mathbf{w}^{\top} \mathbf{x}_{i})) \right]$$

- Bad news: no closed-form solution!
- \bullet Good news: $\mathcal{L}(\mathbf{w})$ is a concave function of w!
 - Is the original logistic function concave?

Optimizing concave/convex function

- Conditional likelihood for logistic regression is concave
- Maximum of a concave function = minimum of a convex function
 - Gradient ascent (concave) / Gradient descent (convex)



$$\nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w}) = \begin{pmatrix} \frac{\partial \mathcal{L}(\mathbf{w})}{\partial w_0} \\ \vdots \\ \frac{\partial \mathcal{L}(\mathbf{w})}{\partial w_d} \end{pmatrix}$$

$$\mathbf{w}_{t+1} = \mathbf{w}_t + \eta \nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w})|_{\mathbf{w}_t}$$

Gradient Ascent for Logistic Regression

Property of sigmoid function



$$\Rightarrow \nabla_{\nu}\psi = \psi(1-\psi)$$

Gradient ascent algorithm iteratively does:

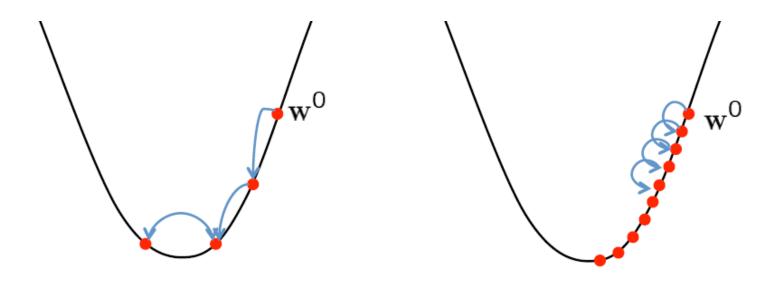
$$\mathbf{w}_{t+1} \leftarrow \mathbf{w}_t + \eta \sum_{i=1}^{N} \mathbf{x}_i \left(y_i - \mu_i^t \right)$$

- where $\mu_i^t = P(y = 1 | \mathbf{x}_i, \mathbf{w}_t)$ is the prediction made by the current model
- Until the change (of objective or gradient) falls below some threshold

Issues

- Gradient descent is the simplest optimization methods, faster convergence can be obtained by using
 - E.g., Newton method, conjugate gradient ascent, IRLS (iterative reweighted least squares)
- The vanilla logistic regression often over-fits; using a regularization can help a lot!

Effects of step-size

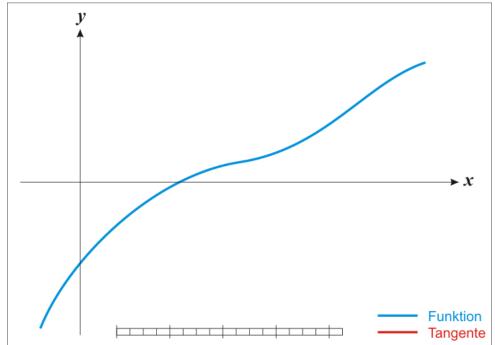


- \bullet Large $\eta =>$ fast convergence but larger residual error; Also possible oscillations
- \bullet Small η => slow convergence but small residual error

The Newton's Method

- AKA: Newton-Raphson method
- A method that finds the root of: f(x) = 0

$$x_{t+1} = x_t - \frac{f(x_t)}{f'(x_t)}$$



For Wikipedia

The Newton's Method

To maximize the conditional likelihood

$$\mathcal{L}(\mathbf{w}) = \sum_{i} \left[y_i \mathbf{w}^{\top} \mathbf{x}_i - \log(1 + \exp(\mathbf{w}^{\top} \mathbf{x}_i)) \right]$$

We need to find w* such that

$$\nabla \mathcal{L}(\mathbf{w}^*) = 0$$

So we can perform the following iteration:

$$\mathbf{w}_{t+1} \leftarrow \mathbf{w}_t - H^{-1} \nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w})|_{\mathbf{w}_t}$$

• where *H* is known as the Hessian matrix:

$$H = \nabla_{\mathbf{w}}^2 \mathcal{L}(\mathbf{w})|_{\mathbf{w}_t}$$

Newton's Method for LR

The update equation

$$\mathbf{w}_{t+1} \leftarrow \mathbf{w}_t - H^{-1} \nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w})|_{\mathbf{w}_t}$$

• where the gradient is:

$$\nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w})|_{\mathbf{w}_t} = \sum_i (y_i - \mu_i) \mathbf{x}_i = X(\mathbf{y} - \boldsymbol{\mu})$$
$$\mu_i = \psi(\mathbf{w}_t^{\top} \mathbf{x}_i)$$

The Hessian matrix is:

$$H = \nabla_{\mathbf{w}}^{2} \mathcal{L}(\mathbf{w})|_{\mathbf{w}_{t}} = -\sum_{i} \mu_{i} (1 - \mu_{i}) \mathbf{x}_{i} \mathbf{x}_{i}^{\top} = -XRX^{\top}$$
where $R_{ii} = \mu_{i} (1 - \mu_{i})$

Iterative reweighted least squares (IRLS)

In least square estimate of linear regression, we have

$$\mathbf{w} = (XX^{\top})^{-1}X\mathbf{y}$$

Now, for logistic regression

$$\mathbf{w}_{t+1} = \mathbf{w}_t - H^{-1} \nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w}_t)$$

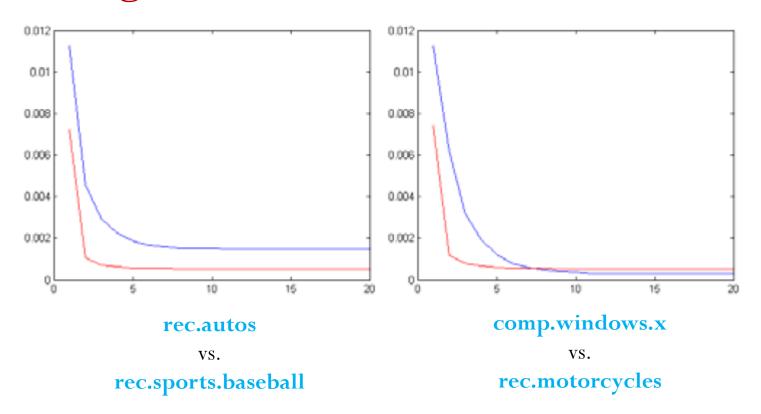
$$= \mathbf{w}_t - (XRX^\top)^{-1} X(\boldsymbol{\mu} - \mathbf{y})$$

$$= (XRX^\top)^{-1} \left\{ XRX^\top \mathbf{w}_t - X(\boldsymbol{\mu} - \mathbf{y}) \right\}$$

$$= (XRX^\top)^{-1} XR\mathbf{z}$$

where
$$\mathbf{z} = X^{\top} \mathbf{w}_t - R^{-1} (\boldsymbol{\mu} - \mathbf{y})$$

Convergence curves



- □ Legend: X-axis: Iteration #; Y-axis: classification error
- □ In each figure, red for IRLS and blue for gradient descent

LR: Practical Issues

- ♦ IRLS takes $O(N + d^3)$ per iteration, where N is # training points and d is feature dimension, but converges in fewer iterations
- Quasi-Newton methods, that approximate the Hessian, work faster
- lacktriangle Conjugate gradient takes O(Nd) per iteration, and usually works best in practice
- ♦ Stochastic gradient descent can also be used if N is large c.f. perceptron rule

Gaussian NB vs. Logistic Regression

<u>GNB</u> Gaussian parameters

VS

Regression parameters

- Representation equivalence
 - But only in some special case! (GNB with class independent variances)
- What's the differences?
 - □ LR makes no assumption about P(X | Y) in learning
 - They optimize different functions, obtain different solutions

Generative vs. Discriminative

- Given infinite data (asymptotically)
 - (1) If conditional independence assumption holds, discriminative and generative NB perform similar

$$\epsilon_{\mathrm{Dis},\infty} \sim \epsilon_{\mathrm{Gen},\infty}$$

 (2) If conditional independence assumption does NOT hold, discriminative outperform generative NB

$$\epsilon_{\mathrm{Dis},\infty} < \epsilon_{\mathrm{Gen},\infty}$$

Generative vs. Discriminative

Given finite data (N data points, d features)

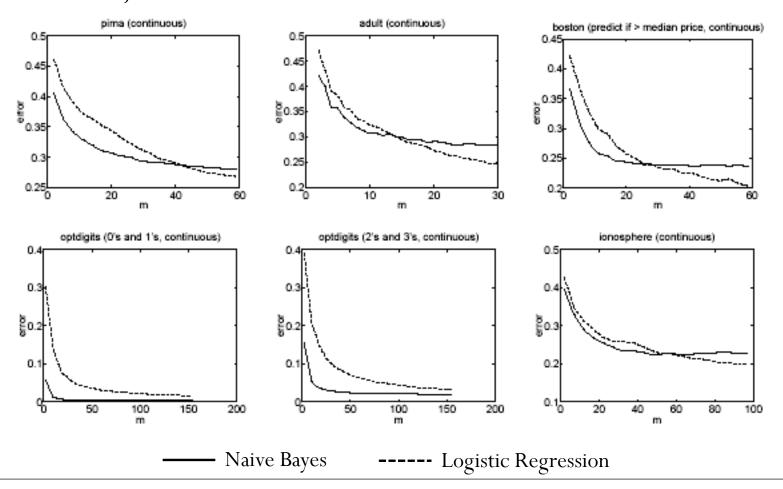
$$\epsilon_{\mathrm{Dis},n} \le \epsilon_{\mathrm{Dis},\infty} + O\left(\sqrt{\frac{d}{N}}\right)$$

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

- lacktriangledown Naive Bayes (generative) requires $N=O(\log d)$ to converge to its asymptotic error, whereas logistic regression (discriminative) requires N=O(d) .
- Why?
 - "Independent class conditional densities" parameter estimates are not coupled, each parameter is learnt independently, not jointly, from training data

Experimental Comparison

• UCI Machine Learning Repository 15 datasets, 8 continuous features, 7 discrete features



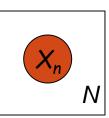
What you need to know

- LR is a linear classifier
 - Decision boundary is a hyperplane
- LR is learnt by maximizing conditional likelihood
 - No closed-form solution
 - Concave! Global optimum by gradient ascent methods
- GNB with class-independent variances representationally equivalent to LR
 - Solutions differ because of objective (loss) functions
- ♦ In general, NB and LR make different assumptions
 - ullet NB: features independent given class, assumption on $P(X \mid Y)$
 - \blacksquare LR: functional form of P(Y | X), no assumption on P(X | Y)
- Convergence rates:
 - GNB (usually) needs less data
 - □ LR (usually) gets to better solutions in the limit

Exponential family

 \diamond For a numeric random variable X

$$p(\mathbf{x}|\boldsymbol{\eta}) = h(\mathbf{x}) \exp\left(\boldsymbol{\eta}^{\top} T(\mathbf{x}) - A(\boldsymbol{\eta})\right)$$
$$= \frac{1}{Z(\boldsymbol{\eta})} h(\mathbf{x}) \exp\left(\boldsymbol{\eta}^{\top} T(\mathbf{x})\right)$$



is an **exponential family distribution** with natural (canonical) parameter η

- \diamond Function T(x) is a sufficient statistic.
- Function $A(\eta) = \log Z(\eta)$ is the log normalizer.
- Examples: Bernoulli, multinomial, Gaussian, Poisson, gamma,...

Recall Linear Regression

• Let us assume that the target variable and the inputs are related by the equation:

$$y_i = \boldsymbol{\theta}^{\top} \mathbf{x}_i + \epsilon_i$$

where \mathcal{E} is an error term of unmodeled effects or random noise

 \bullet Now assume that ε follows a Gaussian $N(0,\sigma)$, then we have:

$$p(y_i|\mathbf{x}_i,\boldsymbol{\theta}) = \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{(y_i - \boldsymbol{\theta}^{\top}\mathbf{x}_i)^2}{2\sigma^2}\right)$$

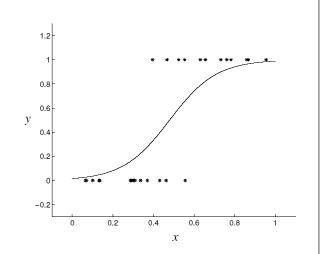
Recall: Logistic Regression (sigmoid classifier)

The condition distribution: a Bernoulli

$$p(y|\mathbf{x}) = \mu(\mathbf{x})^y (1 - \mu(\mathbf{x}))^{1-y}$$

where μ is a logistic function

$$\mu(\mathbf{x}) = \frac{1}{1 + e^{-\boldsymbol{\theta}^{\top} \mathbf{x}}}$$



We can use the brute-force gradient method as in LR

 \diamond But we can also apply generic laws by observing the p(y|x) is an exponential family function, more specifically, a generalized linear model!

Example: Multivariate Gaussian Distribution

 \diamond For a continuous vector random variable $\mathbf{x} \in \mathbb{R}^d$:

$$p(\mathbf{x}|\boldsymbol{\mu}, \boldsymbol{\Sigma}) = \frac{1}{(2\pi)^{d/2} |\boldsymbol{\Sigma}|^{1/2}} \exp\left(-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^{\top} \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu})\right)$$

$$= \frac{1}{(2\pi)^{d/2}} \exp\left(-\frac{1}{2} \mathrm{tr}(\boldsymbol{\Sigma}^{-1} \mathbf{x} \mathbf{x}^{\top}) + \boldsymbol{\mu}^{\top} \boldsymbol{\Sigma}^{-1} \mathbf{x} - \frac{1}{2} \boldsymbol{\mu}^{\top} \boldsymbol{\Sigma}^{-1} \boldsymbol{\mu} - \log |\boldsymbol{\Sigma}|\right)$$

Exponential family representation

$$\boldsymbol{\eta} = \left[\Sigma^{-1} \boldsymbol{\mu}; -\frac{1}{2} \text{vec}(\Sigma^{-1}) \right] = \left[\boldsymbol{\eta}_1; \text{vec}(\boldsymbol{\eta}_2) \right], \ \boldsymbol{\eta}_1 = \Sigma^{-1} \boldsymbol{\mu} \text{ and } \boldsymbol{\eta}_2 = -\frac{1}{2} \Sigma^{-1}$$

Natural parameter

$$T(\mathbf{x}) = [\mathbf{x}; \text{vec}(\mathbf{x}\mathbf{x}^{\top})]$$

$$A(\boldsymbol{\eta}) = \frac{1}{2} \boldsymbol{\mu}^{\top} \Sigma^{-1} \boldsymbol{\mu} + \log |\Sigma| = -\frac{1}{2} \operatorname{tr}(\boldsymbol{\eta}_2 \boldsymbol{\eta}_1 \boldsymbol{\eta}_1^{\top}) - \frac{1}{2} \log(-2|\boldsymbol{\eta}_2|)$$

$$h(\mathbf{x}) = (2\pi)^{-d/2}$$

Note: a d-dimensional Gaussian is a $(d+d^2)$ -parameter distribution with a $(d+d^2)$ -element vector of sufficient statistics (but because of symmetry and positivity, parameters are constrained and have lower degree of freedom)

Example: Multinomial distribution

• For a binary vector random variable $\mathbf{x} \sim \text{multi}(\mathbf{x}|\pi)$:

$$p(\mathbf{x}|\pi) = \prod_{i=1}^{d} \pi_i^{x_i} = \exp\left(\sum_i x_i \ln \pi_i\right)$$

$$= \exp\left(\sum_{i=1}^{d-1} x_i \ln \pi_i + \left(1 - \sum_{i=1}^{d-1} x_i\right) \ln \left(1 - \sum_{i=1}^{d-1} \pi_i\right)\right)$$

$$= \exp\left(\sum_{i=1}^{d-1} x_i \ln \frac{\pi_i}{1 - \sum_{i=1}^{d-1} \pi_i} + \ln \left(1 - \sum_{i=1}^{d-1} \pi_i\right)\right)$$

Exponential family representation

$$\eta = \left[\ln(\pi_i/\pi_d); 0\right]$$

$$T(\mathbf{x}) = \mathbf{x}$$

$$A(\eta) = -\ln\left(1 - \sum_{i=1}^{d-1} \pi_i\right) = \ln\left(\sum_{i=1}^{d} e^{\eta_i}\right)$$

$$h(\mathbf{x}) = 1$$

Why exponential family?

Moment generating property (proof?)

$$\nabla_{\boldsymbol{\eta}} A(\boldsymbol{\eta}) = \nabla_{\boldsymbol{\eta}} \log Z(\boldsymbol{\eta}) = \cdots = \mathbb{E}_{p(\mathbf{x}|\boldsymbol{\eta})}[T(\mathbf{x})]$$

$$\nabla_{\boldsymbol{\eta}}^2 A(\boldsymbol{\eta}) = \dots = \operatorname{Var}[T(\mathbf{x})]$$

Moment estimation

- We can easily compute moments of any exponential family distribution by taking the derivatives of the log normalizer $A(\eta)$.
- \bullet The q^{th} derivative gives the q^{th} centered moment.

$$\nabla_{\boldsymbol{\eta}} A(\boldsymbol{\eta}) = \text{mean}$$

$$\nabla_{\boldsymbol{\eta}}^2 A(\boldsymbol{\eta}) = \text{variance}$$

:

Moment vs canonical parameters

The moment parameter μ can be derived from the natural (canonical) parameter

$$\nabla_{\boldsymbol{\eta}} A(\boldsymbol{\eta}) = \mathbb{E}_{p(\mathbf{x}|\boldsymbol{\eta})}[T(\mathbf{x})] \triangleq \boldsymbol{\mu}$$

 $A(\eta)$ is convex since

$$\nabla_{\boldsymbol{\eta}}^2 A(\boldsymbol{\eta}) = \text{Var}[T(\mathbf{x})] > 0$$

♦ Hence we can invert the relationship and infer the canonical parameter from the moment parameter (1-to-1):

$$\eta \triangleq \psi(\mu)$$

A distribution in the exponential family can be parameterized not only by η – the canonical parameterization, but also by μ – the moment parameterization.

IID Sampling for Exponential Family

♦ For exponential family distribution, we can obtain the sufficient statistics by inspection once represented in the standard form

$$p(\mathbf{x}|\boldsymbol{\eta}) = h(\mathbf{x}) \exp(\boldsymbol{\eta}^{\top} T(\mathbf{x}) - A(\boldsymbol{\eta}))$$

Sufficient statistics:

$$T(\mathbf{x})$$

For IID sampling, the joint distribution is also an exponential family

$$p(D|\boldsymbol{\eta}) = \prod h(\mathbf{x}_i) \exp \left(\boldsymbol{\eta}^{\top} T(\mathbf{x}_i) - A(\boldsymbol{\eta})\right)$$

$$= \left(\prod_i h(\mathbf{x}_i)\right) \exp\left(\boldsymbol{\eta}^\top \sum_i T(\mathbf{x}_i) - NA(\boldsymbol{\eta})\right)$$

Sufficient statistics:

$$\sum_{i} T(\mathbf{x}_i)$$

MLE for Exponential Family

♦ For *iid* data, the log-likelihood is

$$\mathcal{L}(\boldsymbol{\eta}; D) = \sum_{n} \log h(\mathbf{x}_n) + \left(\boldsymbol{\eta}^{\top} \sum_{n} T(\mathbf{x}_n)\right) - NA(\boldsymbol{\eta})$$

◆ Take derivatives and set to zero:

$$\nabla_{\boldsymbol{\eta}} \mathcal{L}(\boldsymbol{\eta}; D) = \sum_{n} T(\mathbf{x}_{n}) - N \nabla_{\boldsymbol{\eta}} A(\boldsymbol{\eta}) = 0$$
$$\nabla_{\boldsymbol{\eta}} A(\boldsymbol{\eta}) = \frac{1}{N} \sum_{n} T(\mathbf{x}_{n})$$

- ightharpoonup This amounts $\hat{\mu}_{MLE} = \frac{1}{N} \sum_{n} T(\mathbf{x}_n)_{n}$ only involve sufficient stiatistics!
- \diamond We can infer the canonical parameters using

$$\hat{\boldsymbol{\eta}}_{MLE} = \psi(\hat{\boldsymbol{\mu}}_{MLE})$$

Examples

• Gaussian:
$$\eta = \left[\Sigma^{-1} \mu; -\frac{1}{2} \text{vec}(\Sigma^{-1}) \right]$$

 $h(\mathbf{x}) = (2\pi)^{-d/2}$

$$T(\mathbf{x}) = [\mathbf{x}; \text{vec}(\mathbf{x}\mathbf{x}^{\top})]$$
$$A(\boldsymbol{\eta}) = \frac{1}{2}\boldsymbol{\mu}^{\top} \boldsymbol{\Sigma}^{-1} \boldsymbol{\mu} + \log |\boldsymbol{\Sigma}|$$

$$oldsymbol{\eta} = [\ln(\pi_i/\pi_d); 0]$$

$$T(\mathbf{x}) = \mathbf{x}$$

$$h(\mathbf{x}) = 1$$

$$\bullet$$
 Poisson: $\eta = \log \lambda$

Poisson:
$$\eta = \log \lambda$$

$$T(x) = x$$

$$A(\eta) = \lambda = e^{\eta}$$

$$h(x) = \frac{1}{x!}$$

 $T(\mathbf{x}) = [\mathbf{x}; \text{vec}(\mathbf{x}\mathbf{x}^{\top})]$ $\Rightarrow \hat{\boldsymbol{\mu}}_{MLE} = \frac{1}{N} \sum T_1(\mathbf{x}_n) = \frac{1}{N} \sum \mathbf{x}_n$

$$\eta = [\ln(\pi_i/\pi_d); 0]$$

$$T(\mathbf{x}) = \mathbf{x}$$

$$A(\eta) = -\ln\left(1 - \sum_{i=1}^{d-1} \pi_i\right) \qquad \Rightarrow \hat{\boldsymbol{\mu}}_{MLE} = \frac{1}{N} \sum_{n} \mathbf{x}_n$$

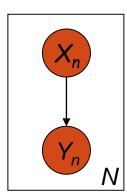
$$\Rightarrow \hat{\mu}_{MLE} = \frac{1}{N} \sum_{n} x_n$$

Generalized Linear Models (GLIMs)

- The graphical model
 - Linear regression
 - Discriminative linear classification
 - Commonality:

model
$$\mathbb{E}_p[y] = \mu = f(\boldsymbol{\theta}^\top \mathbf{x})$$

- What is p()? the cond. dist. of Y.
- What is f()? the response function.



GLIM

- The observed input \boldsymbol{x} is assumed to enter into the model via a linear combination of its elements $\boldsymbol{\xi} = \boldsymbol{\theta}^{\mathsf{T}} \boldsymbol{x}$
- The conditional mean μ is represented as a function $f(\xi)$ of ξ , where f is known as the response function
- The observed output \mathbf{y} is assumed to be characterized by an exponential family distribution with conditional mean μ .

GLIM, cont.

$$\theta \longrightarrow \xi \qquad f \qquad \psi \qquad p(y|\eta) = h(y) \exp\left\{\eta^{T}(x)y - A(\eta)\right\}$$

$$\Rightarrow p(y|\eta, \phi) = h(y, \phi) \exp\left\{\frac{1}{\phi}\left(\eta^{T}(x)y - A(\eta)\right)\right\}$$

- lacktriangle The choice of exp family is constrained by the nature of the data $oldsymbol{Y}$
 - □ Example: y is a continuous vector → multivariate Gaussian
 y is a class label → Bernoulli or multinomial
- ♦ The choice of the response function
 - □ Following some mild constrains, e.g., [0,1]. Positivity ...
 - Canonical response function:
 - In this case $\theta^{\mathsf{T}} \mathbf{X}$ directly corresponds to canonical parameter η .

$$f = \psi^{-1}(\cdot)$$

MLE for GLIMs

Log-likelihood

$$\mathcal{L}(\boldsymbol{\theta}; D) = \sum_{n} \log h(y_n) + \sum_{n} (\eta_n y_n - A(\eta_n))$$
where $\eta_n = \psi(\mu_n), \ \mu_n = f(\xi_n) \text{ and } \xi_n = \boldsymbol{\theta}^\top \mathbf{x}_n$

Derivative of Log-likelihood

$$\nabla_{\boldsymbol{\theta}} \mathcal{L} = \sum_{n} \left(y_n \nabla_{\boldsymbol{\theta}} \eta_n - \frac{dA(\eta_n)}{d\eta_n} \nabla_{\boldsymbol{\theta}} \eta_n \right)$$

 $= \sum_{n} (y_n - \mu_n) \nabla_{\theta} \eta_n$ This is a fixed point function because μ is a function of θ

MLE for GLIMs with canonical response

Log-likelihood

$$\mathcal{L}(\boldsymbol{\theta}; D) = \sum_{n} \log h(y_n) + \sum_{n} (\boldsymbol{\theta}^{\top} \mathbf{x}_n y_n - A(\eta_n))$$

Derivative of Log-likelihood

$$\nabla_{\boldsymbol{\theta}} \mathcal{L} = \sum_{n} \left(\mathbf{x}_{n} y_{n} - \frac{dA(\eta_{n})}{d\eta_{n}} \nabla_{\boldsymbol{\theta}} \eta_{n} \right)$$
$$= \sum_{n} (y_{n} - \mu_{n}) \mathbf{x}_{n}$$
This is a

$$=X(\mathbf{y}-\boldsymbol{\mu})$$

This is a fixed point function because μ is a function of θ

- Online learning for canonical GLIMs
 - Stochastic gradient ascent = least mean squares (LMS) algorithm: $\boldsymbol{\theta}_{t+1} = \boldsymbol{\theta}_t + \rho(y_n \mu_n^t)\mathbf{x}_n$

where
$$\mu_n^t = f(\boldsymbol{\theta}_t^{\top} \mathbf{x}_n)$$
 and ρ is a step size

MLE for GLIMs with canonical response

Log-likelihood

$$\mathcal{L}(\boldsymbol{\theta}; D) = \sum_{n} \log h(y_n) + \sum_{n} (\boldsymbol{\theta}^{\top} \mathbf{x}_n y_n - A(\eta_n))$$

Derivative of Log-likelihood

$$\nabla_{\boldsymbol{\theta}} \mathcal{L} = \sum_{n} \left(\mathbf{x}_{n} y_{n} - \frac{dA(\eta_{n})}{d\eta_{n}} \nabla_{\boldsymbol{\theta}} \eta_{n} \right)$$
$$= \sum_{n} (y_{n} - \mu_{n}) \mathbf{x}_{n}$$
This is a

 $=X(\mathbf{y}-\boldsymbol{\mu})$

This is a fixed point function because μ is a function of θ

- Batch learning applies
 - E.g., the Newton's method leads to an Iteratively Reweighted Least Square (IRLS) algorithm

What you need to know

- Exponential family distribution
- Moment estimation
- Generalized linear models
- Parameter estimation of GLIMs