Basis expansions, kernels and support vector machines

Hastie, Tibshirani, Friedman Ch 5, 6, 12 Kevin Murphy Ch. 14

> CS 6140 Machine Learning Professor Olga Vitek

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Basis Expansion

HTF Ch5

Move beyond linearity

- Linear regression, logistic regression, LDA
 - Classification by linear hyperplanes
 - Easy to fit and to interpret
- f(Y|X) is typically non-linear and non-additive
 - augment X with transformations of X
 - use as input in linear models
 - Denote mth transformation $h_m(\mathbf{X}): \mathbb{R}^p \to \mathbb{R}^p$

- Model
$$f(\mathbf{X}) = \sum_{m=1}^{M} \beta_m h_m(\mathbf{X})$$

- $f(\mathbf{X})$ is linear in $h_m(\mathbf{X})$
- This is a *linear basis expansion* in ${f X}$

Linear basis expansion

- Linear model
 - $-h_m(\mathbf{X}) = X_m$
- Polynomial terms (Taylor expansion)
 - $h_m(\mathbf{X}) = X_m^2$ or $h_m(\mathbf{X}) = X_i X_j^2$
 - # variables \uparrow exponentially in p
 - tweak one region ↑ flap another
- Functions of a vector

-
$$h_m(\mathbf{X}) = log(X_j), h_m(\mathbf{X}) = \sqrt{X_j}, h_m(\mathbf{X}) = ||\mathbf{X}||$$

- Indicators $h_m(\mathbf{X}) = I(L_m \le X_k < U_m)$
 - Break the range of X_k into regions
 - Piecewise constant model in each region
- Piece-wise polynomials and splines
 - Dictionary \mathcal{D} of basis functions
 - Method for controlling model complexity:
 restriction, selection, regularization

- Restriction:
$$f(\mathbf{X}) = \sum_{j=1}^{p} f_j(X_j) = \sum_{j=1}^{p} \sum_{m=1}^{M_j} \beta_{jm} h_{jm}(X_j)$$

Piecewise fits

- ullet Assume X is one-dimensional (i.e., X)
 - Divide domain(X) into contiguous intervals
 - f(X): a separate polynomial in each interval
- Example: piecewise constant

$$- h_1(X) = I(X < \xi_1), \ h_2(X) = I(\xi_i \le X \le \xi_2),$$
$$h_3 = I(\xi_2 \le X)$$

$$- f(X) = \sum_{m=1}^{3} \beta_m h_m(X)$$

- $\hat{\beta}_m = \bar{Y}_m$, the mean of mth region
- Example: piecewise linear
 - Need extra basis $h_{m+3} = h_m(X) \cdot X$, $m = 1, \dots, 3$
- Example: restricted piecewise linear

$$- f(\xi_1^-) = f(\xi_1^+) \rightarrow \beta_1 + \xi_1 \beta_4 = \beta_2 + \xi_1 \beta_5$$

- 3 intervals: 4 free parameters out of 6
- Alternatively: $h_1(X) = 1$, $h_2(X) = X$, $h_3(X) = (X \xi_1)_+$, $h_4(X) = (X \xi_2)_+$

Piecewise fit

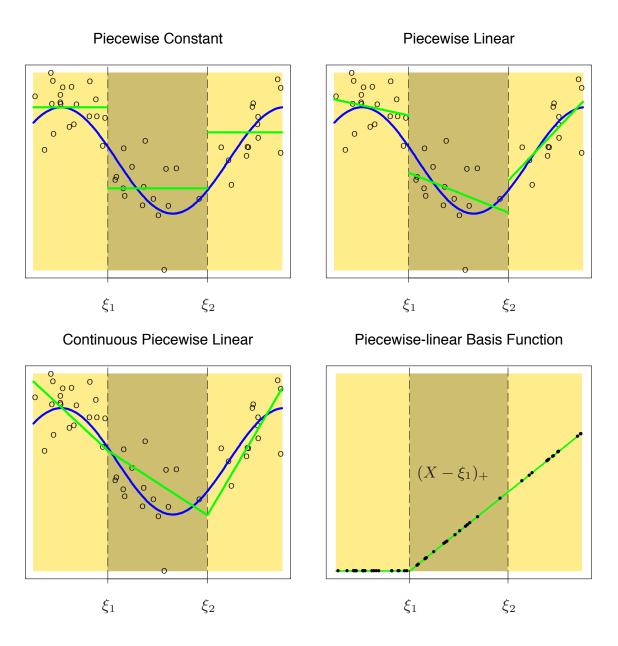
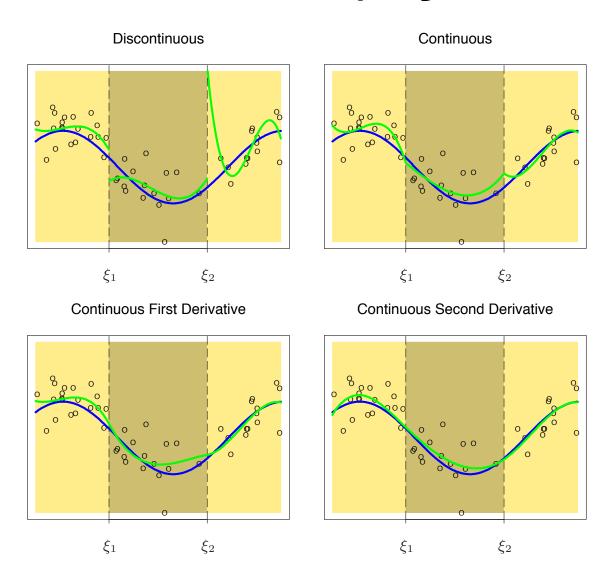


Fig. 5.1. Hastie, Tibshirani, Friedman
The Elements of Statistical Learning 2008

Splines

- Smoother functions
 - Increase the order of polynomials
 - Cubic splines: continuous to second derivative
 - $h_1(X) = 1, h_s(X) = X, h_3(X) = X^2, h_4(X) = X^3$ $h_5(X) = (X \xi_1)_+^3, \ h_6(X) = (X \xi_2)_+^3$
- Example on the next page
 - 6 basis functions
 - 6-dimensional linear space of functions
 - (3 regions) \times (4 parameters per region)
 - (2 knots) \times (3 constrains per knot) = 6
- Order-M spline with knots ξ_j , $j=1,\ldots,K$
 - Piecewise polynomial up to order ${\cal M}$
 - Has continuous derivatives up to order M-2.
 - Piecewise constant fit is order-1 spline.
 Piecewise continuous fit is order-2 spline.
 Cubic spline is order-4 spline.
 - Basis set: $h_j(X) = X^{j-1}, \ j = 1, ..., M$ and $h_{M+l}(X) = (X \xi_l)_+^{M-1}, \ l = 1, ..., K$

Piecewise cubic polynomials



Cubic spline is the lowest-order spline for which the knot-discontinuity is not visible to human eye

Fig. 5.2. Hastie, Tibshirani, Friedman

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Natural cubic splines

- Need for extra stability
 - Polynomials and splines have erratic behavior near the boundaries
 - Variance explodes
 - Extrapolation is problematic
- Natural cubic splines: extra constraints
 - Linear functions beyond boundary knots
 - K knots = K basis functions
 - Basis for cubic splines → impose constraints
 - Start from basis set, impose boundary constraint, derive reduced basis:

$$N_1(X) = 1, \ N_2(X) = X,$$

$$N_{k+2}(X) = d_k(X) - d_{K-1}(X)$$
 where
$$d_k(X) = \left\{ (X - \xi_k)_+^3 - (X - \xi_K)_+^3 \right\} / \left\{ \xi_K - \xi_k \right\}$$

- Second and third derivatives are 0 for $X \geq \xi_K$

Pointwise variance curve

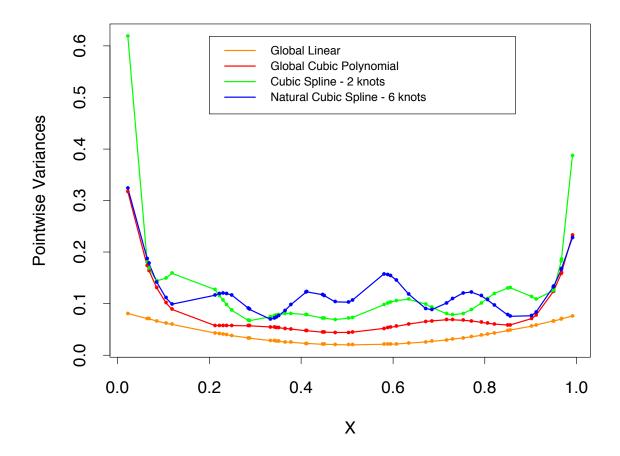


Fig. 5.3. Hastie, Tibshirani, Friedman *The Elements of Statistical Learning* 2008

Example: South African Heart Disease

Logistic regression

$$\frac{\Pr\{\mathsf{chd} = 1|X\}}{1 - \Pr\{\mathsf{chd} = 1|X\}} = \theta_0 + h_1(X_1)'\theta_1 + \ldots + h_p(X_p)'\theta_p$$

- θ_j are vectors of coefficients multiplying the vector of natural spline basis functions h_j
- Use 4 natural spline basis functions for each term in the model
- Knots chosen at uniform quantiles of X_i :
 - 3 internal knots +
 - 2 boundary knots at the extremes of X_i
- Binary predictor has a single coefficient
- More compactly: combine p vectors of basis functions + constant term in a big vector h(X), and model $h(X)'\theta$
- Backward stepwise selection + AIC to drop terms
- Plot prediction $\pm 2 \cdot SE$

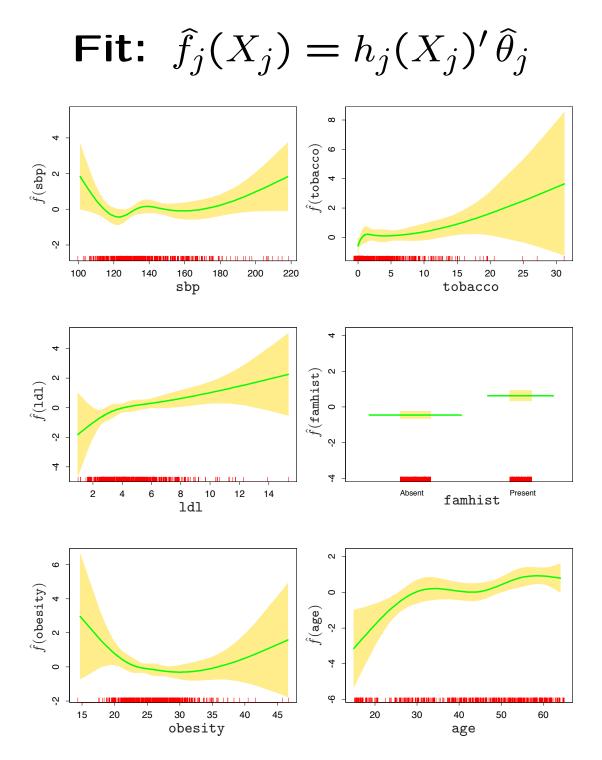
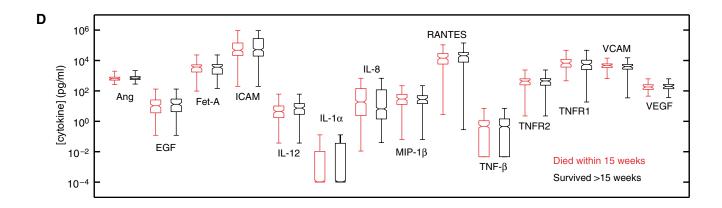


Fig. 5.4. Hastie, Tibshirani, Friedman *The Elements of Statistical Learning* 2008

Example: disease classification

Knickerbocker *et al.* "An integrated approach to prognosis using protein microarrays and nonparametric methods". *Molecular Systems Biology*, 2007

- Goal: predict mortality in patients w/kidney dialysis
 - 468 patients, 208 died within 15 weeks of diagnoses
 - 14 proteins ("messenger" molecule that allows cells to communicate and alter function)
 - 11 clinical characteristics (age, race, bmi...)
 - Proteins were uninformative in isolation



Approach

- Separate predictive model for clinical and molecular measurements
 - Logistic regression (Y = death within 15 weeks)

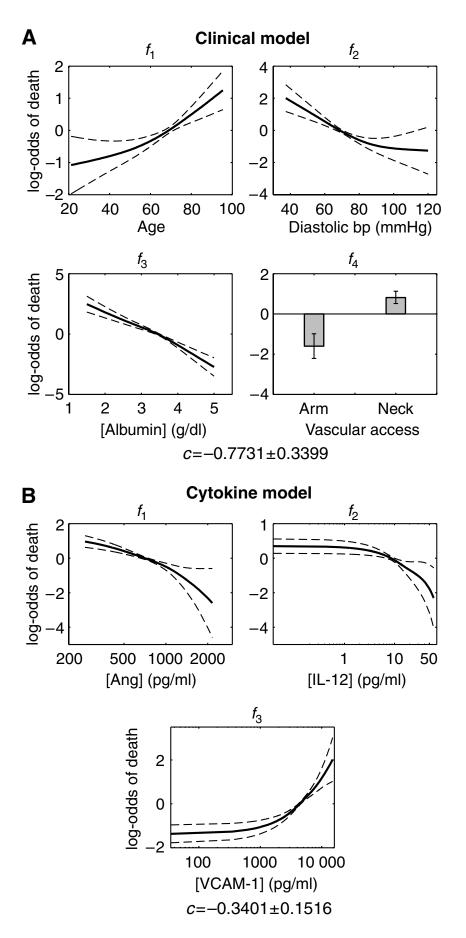
$$log - odds of death = log(P_{sample}(death)/P_{sample}(survival))$$

$$= c + \sum_{p=1}^{M} b_p x_p$$

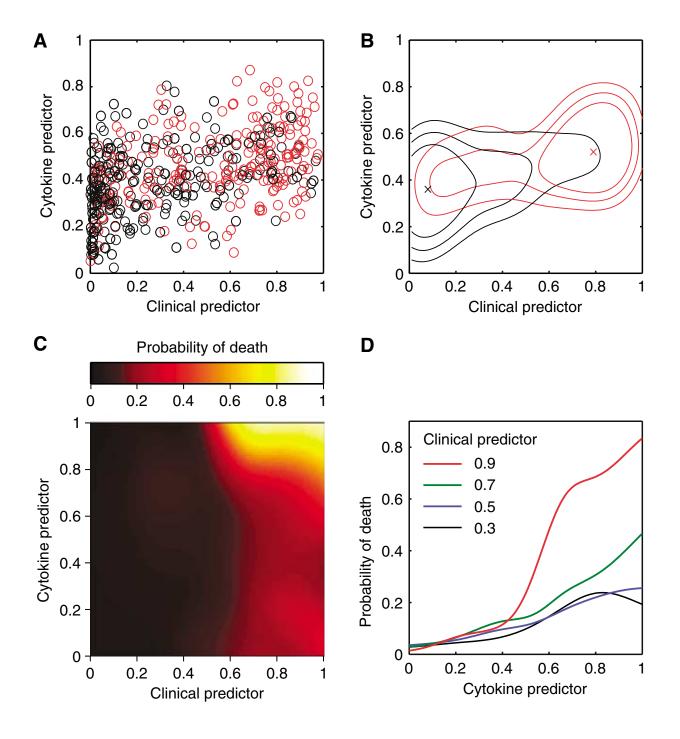
- Additive model to reduce feature space
- Exhaustive search for best M-variable model
- Non-parametric version of the best model: splines

$$log - odds of death = log(P_{sample}(death)/P_{sample}(survival))$$

$$= c + \sum_{p=1}^{M} f_p(x_p)$$



Combined prediction



Controlling model complexity (fixed knots)

Restriction

Limit the class of functions

$$f(\mathbf{X}) = \sum_{j=1}^{p} f_j(X_j) = \sum_{j=1}^{p} \sum_{m=1}^{M_j} \beta_{jm} h_{jm}(X_j)$$

- Model complexity \sim number of basis functions

Selection

- AIC, BIC, significance testing
- Adaptively include basis functions that contribute to prediction (include CART, boosting...)

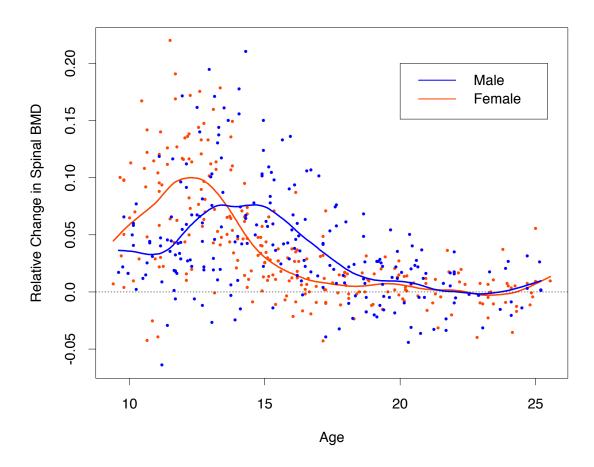
Regularization

 Include the entire dictionary of basis functions, but restrict the coefficients

$$RSS(f,\lambda) = \sum_{i=1}^{N} \{y_i - f(x_i)\}^2 + \lambda \int \{f''(t)\}^3 dt$$

- λ is a smoothing parameter ($\lambda = 0$: any function; $\lambda = \infty$: linear fit)
- Maximized with a natural cubic spline
- Equivalent to generalized ridge regression

Example



Rel. change in bone mineral density \sim age, $\lambda = 0.00022$

$$\begin{aligned} \mathsf{EPE}(\widehat{f}_{\lambda}) &= E\left(Y - \widehat{f}_{\lambda}\right)^{2} \\ &= \mathsf{Var}(Y) + E\left[\mathsf{Bias}^{2}(\widehat{f}_{\lambda}) + \mathsf{Var}(\widehat{f}_{\lambda})\right] \\ &= \sigma^{2} + \mathsf{MSE}(\widehat{f}_{\lambda}) \end{aligned}$$

Choose λ by cross-validation

Fig. 5.6. Hastie, Tibshirani, Friedman

The Elements of Statistical Learning 2008

Fitting multivariate models

Example: linear additive model

$$Y = \alpha + \sum_{j=1}^{p} f_j(X_j) + \varepsilon$$

Algorithm 9.1 The Backfitting Algorithm for Additive Models.

- 1. Initialize: $\hat{\alpha} = \frac{1}{N} \sum_{1}^{N} y_i, \ \hat{f}_j \equiv 0, \forall i, j.$
- 2. Cycle: $j = 1, 2, \dots, p, \dots, 1, 2, \dots, p, \dots$

$$\hat{f}_j \leftarrow \mathcal{S}_j \left[\{ y_i - \hat{\alpha} - \sum_{k \neq j} \hat{f}_k(x_{ik}) \}_1^N \right],$$

$$\hat{f}_j \leftarrow \hat{f}_j - \frac{1}{N} \sum_{i=1}^N \hat{f}_j(x_{ij}).$$

until the functions \hat{f}_j change less than a prespecified threshold.

 S_j is a cubic spline. Iteratively smooth the residual fit for one predictor at a time, until convergence.

See Algorithm 9.2 for logistic regression

Hastie, Tibshirani, Friedman
The Elements of Statistical Learning 2008

Kernel Methods

HTF Ch6

One-dimensional kernel smoothers

- Different model at each point x_0
 - Only use observations close to target x_0
 - Weigh the neighbors x_1 with a kernel $K_{\lambda}(x_0, x_i)$
 - Weight based on distance from x_0
 - $-\lambda$ is a parameter
 - The resulting function is smooth
- K nearest neighbors
 - $-\hat{f}(x_0) = \text{Ave}(y_i|x_i \in N_k(x_0))$ (discontinuous in x)
- Nadaraya-Watson kernel-weighted average

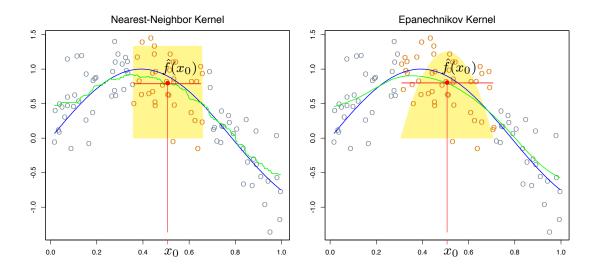
$$-\widehat{f}(x_0) = \frac{\sum_{i=1}^{N} K_{\lambda}(x_0, x_i) y_i}{\sum_{i=1}^{N} K_{\lambda}(x_0, x_i)}, \text{ where}$$

$$K_{\lambda}(x_0, x) = D\left(\frac{|x - x_0|}{\lambda}\right) \text{ and}$$

$$D(t) = \begin{cases} \frac{3}{4}(1 - t^2) & \text{if} \quad |t| \leq 1\\ 0 & \text{otherwise} \end{cases}$$

— Points near the boundary have weight ~ 0 \rightarrow smoothing

Example



- 100 pairs (x_i, y_i)
- Green: Left: 30-NN running mean. Right: Kernel-weighted average, $\lambda = 0.2$
- Orange: observations contributing to the fit

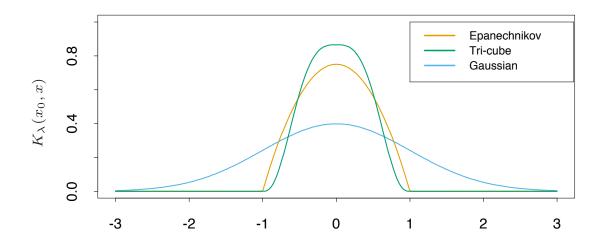
Fig. 6.1. Hastie, Tibshirani, Friedman

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Adaptive width

- Define $h_{\lambda}(x_0)$ a width function that determines the neighborhood of x_0
- Define $K_{\lambda} = D\left(\frac{|x-x_0|}{h_{\lambda}(x_0)}\right)$
- Concerns:
 - Determine λ . Larger $\lambda \to \text{lower variance but}$ higher bias
 - $h_{\lambda}(x)$ constant \rightarrow variance inversely proportional to density of points
 - Nearest neighbor \rightarrow bias inversely proportional to density of points

Kernel examples



• Epanechnikov quadratic kernel $K_{\lambda}(x_0,x) = D\left(rac{|x-x_0|}{\lambda}
ight)$

$$D(t) = \begin{cases} \frac{3}{4}(1-t^2) & if \quad |t| \le 1\\ 0 & \text{otherwise} \end{cases}$$

• Tri-cube function

$$D(t) = \begin{cases} (1-t^3)^3 & if & |t| \le 1\\ 0 & \text{otherwise} \end{cases}$$

Fig. 6.2. Hastie, Tibshirani, Friedman

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Local linear regression

- Smoothly varying locally weighted average
 - Problem: asymmetry of the kernel on the boundary of the domain \rightarrow bias
 - Solution: fit straight lines rather than constants
 - Also helps if values of x are unequally spaced
- Separate weighted least squares problem

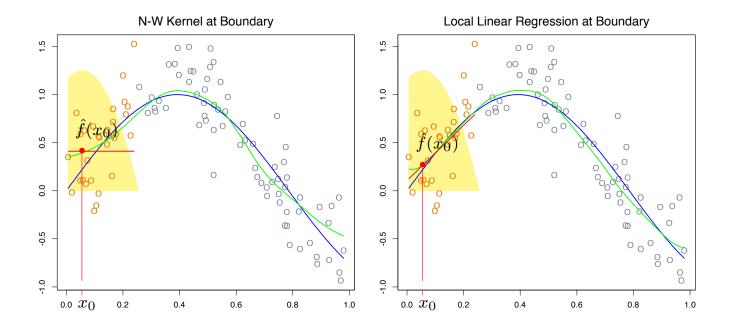
- At each
$$x_0$$
: $\min_{\alpha(x_0),\beta(x_0)} \sum_{i=1}^N K_{\lambda}(x_0,x_i) \left[y_i - \alpha(x_0) - \beta(x_0) x_i \right]^2$

- Define b(x)' = (1, x)
 - B the $X \times 2$ regression matrix with *i*th row b(x)'
 - $\mathbf{W}(x_0)$ the $N \times N$ diagonal matrix diag $\{K_{\lambda}(x_0, x_i)\}$.

$$\widehat{f}(x_0) = b(x_0)' \left(\mathbf{B}' \mathbf{W}(x_0) \mathbf{B} \right)^{-1} \mathbf{B}' \mathbf{W}(x_0) \mathbf{y}$$
$$= \sum_{i=1}^{N} l_i(x_0) y_i, \quad l_i(x_0) \text{ is a local function}$$

- Prediction performed at a single point x_0 : $\hat{f}(x_0) = \hat{\alpha}(x_0) + \hat{\beta}(x_0)x_0$

Example



- True function is linear. Most observations in the neighborhood exceed the target point.
- Left: locally weighted average: bias near the boundaries of *X*
- Right: locally weighted linear regression removes the bias to first order

Fig. 6.3. Hastie, Tibshirani, Friedman

The Elements of Statistical Learning 2008

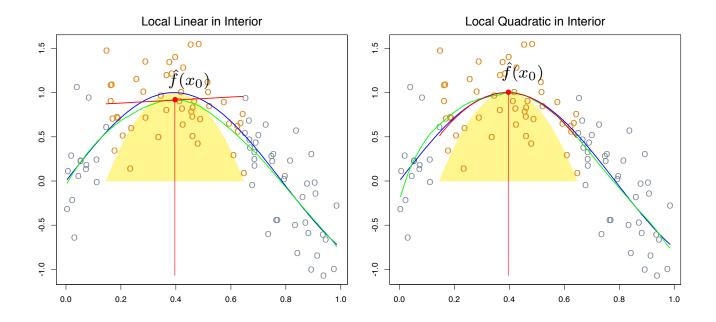
Extensions / comments

- Local polynomial regression
 - Do not have to stop at linear terms
 - Minimize

$$\min_{lpha(x_0),eta_j(x_0),j=1,...,d} \sum_{i=1}^N K_\lambda(x_0,x_i) \left[y_i - lpha(x_0) - \sum_{j=1}^d eta_j(x_0) x_i^j
ight]^2$$

- Solution $\widehat{f}(x_0) = \widehat{\alpha}(x_0) + \sum_{j=1}^d \widehat{\beta}_j(x_0) x_0^j$
- Avoid "trimming the hills and filling the valleys"
- Bias-variance tradeoff
 - Smaller window → lower bias, higher variance
- Default hoice of λ
 - Epanechnikov/tri-cube: radius of support reg.
 - Gaussian: standard deviation
 - Better option: cross-validation
- Model complexity
 - Estimated degrees of freedom $df = trace(S_{\lambda})$, where $\{S_{\lambda}\}_{ij} = l_i(x_j)$

Example: polynomial regression



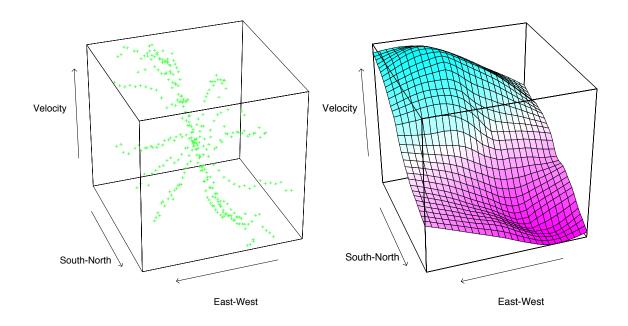
If
$$Y = f(X) + \varepsilon$$
, $\varepsilon \stackrel{iid}{\sim} (0, \sigma^2)$, then

- $Var\{\hat{f}(x_0)\} = \sigma^2 ||l(x_0)||$, where $l(x_0)$ is vector of kernel weights at x_0
- Can show that $l(x_0)$ increases with d
- Implies greater bias-variance trade-off

Fig. 6.5. Hastie, Tibshirani, Friedman

The Elements of Statistical Learning 2008

Local regression in \mathbb{R}^p



- ullet p-dimensional Kernel; polynomial fit of degree d
- Define b(X)':

-
$$d = 1$$
, $p = 2$: $(1, X_1, X_2)$
- $d = 2$, $p = 2$: $(1, X_1, X_2, X_1^2, X_2^2, X_1X_2)$

• At each x_0 , solve

$$\min_{eta(x_0)} \sum_{i=1}^N K_\lambda(x_0,x_i) \left(y_i - b(x_i)'eta(x_0)
ight)^2$$
, where $K_\lambda(x_0,x) = d\left(rac{||x-x_0||}{\lambda}
ight)$

Fig. 6.8. Hastie, Tibshirani, Friedman
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Extensions

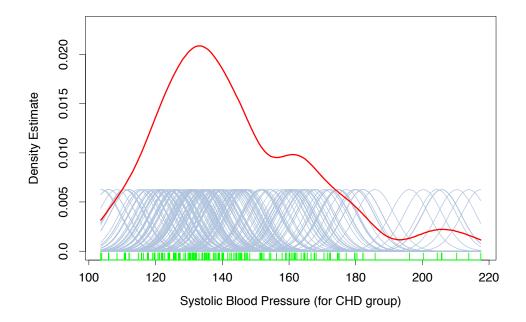
- Structured Kernels
 - As dimensions ↑, # neighbor points ↓
 - Need additional constraints on local regression, to counter curse of dimensionality
 - Option: modify the kernel

$$K_{\lambda,A}(x_0,x) = d\left(\frac{(x-x_0)'\mathbf{A}(x-x_0)}{\lambda}\right)$$

- Restrict A to downgrade or omit directions
- Local likelihood: local fit of any other model
 - E.g., logistic regression

$$l(\beta(x_0) = \sum_{i=1}^{N} K_{\lambda}(x_0, x_i) \, l(y_i, x_i' \, \beta(x_0))$$

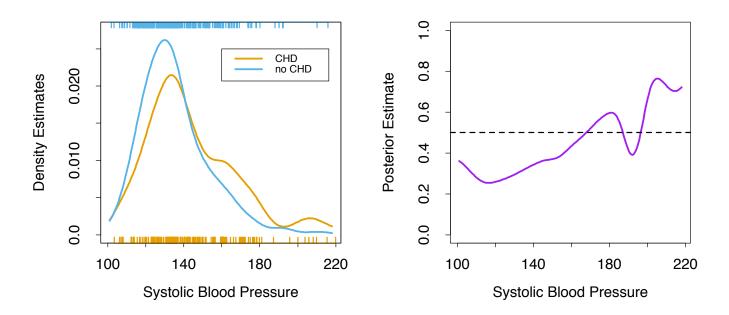
Kernel density estimation



- Natural (bumpy) estimate $\hat{f}_X(x_0) = \frac{\#x_i \in \mathcal{N}(x_0)}{N\lambda}$ where $\mathcal{N}(x_0)$ is a neighborhood of x_0 of width λ
- Smoothed estimate $\hat{f}_X(x_0) = \frac{1}{N\lambda} \sum_{i=1}^N K_\lambda(x_0, x_i)$ E.g., Gaussian kernel $K_\lambda(x_0, x_i) = \phi(|x - x_0|/\lambda)$ where ϕ is pdf of $\mathcal{N}(0, 1)$
- Equivalent to local weighted average

Fig. 6.13. Hastie, Tibshirani, Friedman
The Elements of Statistical Learning 2008

Kernel density classification



Nonparametric classification

$$\widehat{Pr}\{G = j | X = x_0\} = \frac{\widehat{\pi}_j \widehat{f}_j(x_0)}{\sum\limits_{k=1}^{J} \widehat{\pi}_k \widehat{f}_k(x_0)}$$

• Naïve Bayes $\widehat{Pr}\{G = k | X = x_0\} = \frac{\widehat{\pi}_k \widehat{f}_k(x_0)}{\sum\limits_{j=1}^J \widehat{\pi}_j \widehat{f}_j(x_0)} = \frac{\widehat{\pi}_k \prod_{l=1}^p \widehat{f}_{kl}(x_0)}{\sum\limits_{j=1}^J \widehat{\pi}_j \prod_{l=1}^p \widehat{f}_{jl}(x_0)}$

Fig. 6.14. Hastie, Tibshirani, Friedman The Elements of Statistical Learning 2008

More on naïve Bayes

- Often works well
 - Biased class densities, but less variance
 - Bias may be small near the decision boundary
- Connection to GAMs

$$\begin{aligned} \log \mathrm{it} \frac{Pr\{G = k | X\}}{Pr\{G = J | X\}} &= \log \frac{\pi_k \prod_{l=1}^p f_{kl}(X)}{\pi_J \prod_{l=1}^p f_{Jl}(X)} \\ &= \log \frac{\pi_k}{\pi_J} + \sum_{l=1}^p \log \frac{f_{kl}(X)}{f_{Jl}(X)} \\ &= \beta_{0k} + \sum_{l=1}^p g_l(X) \end{aligned}$$

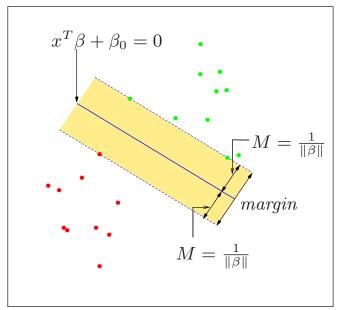
- Similar to GAM
- Difference: GAMs do not specify probability distribution on X, while Naïve Bayes does
- Same distinction as logistic regression vs LDA

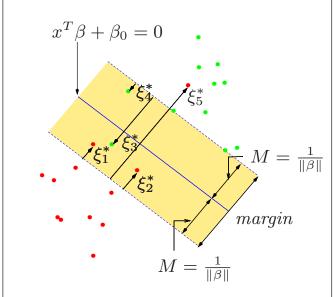
Support Vector Machines

HTF 12.1-12.3

KM Ch14.1-14.7

Support vector classifiers





- Kernel trick
 - Defines high-dimensional feature vector
 - Prevents underfitting
- Sparsity and large margin principle
 - Ensure that we do not use all dimensions
 - Prevent overfitting

Fig. 12.1. Hastie, Tibshirani, Friedman *The Elements of Statistical Learning* 2008

Separable: HTF 12.2

- Data: $\{(x_i, y_i)\}_{i=1,...,N}$, $x_i \in \mathbb{R}^p$, $y_i \in \{-1, 1\}$
- Define a hyperplane
 - $\{x : f(x) = x'\beta + \beta_0 = 0\}, ||\beta|| = 1$
 - Distance from a point x to hyperplane: $f(x) = x'\beta + \beta_0$
 - Classification rule $G(x) = sign(x'\beta + \beta_0)$
- Since the data are separable:
 - Can find β such that $y_i f(x_i) > 0$ for all i
 - Can find β to maximize the margin between training points with class 1 and -1
 - Solve optimization problem $\max_{\beta,\beta_0,||\beta||=1}M \text{ s.t. } y_i(x_i'\beta+\beta_0)\geq M \text{, } i=1,\dots,N$
 - Can show that, equivalently, $\min_{\substack{\beta,\beta_0\\ \text{(drop constraint on } ||\beta||; \ M=1/||\beta||)}} \text{s.t. } y_i(x_i'\beta+\beta_0) \geq 1, \ i=1,\ldots,N$

Non-separable: HTF 12.2

- Classes overlap in feature space
- Goal: maximize ||M||
 - Some points can be on the wrong side of margin
- Solution: slack variables $\xi = (\xi_1, \dots, \xi_N)$
 - Modify previous optimization problem: $\max_{\beta,\beta_0,||\beta||=1} M \text{ s.t. } \xi_i \geq 0, \ \sum_{i=1}^N \xi_i \leq \text{constant}$ $y_i(x_i'\beta+\beta_0) \geq M-\xi_i \text{ or } y_i(x_i'\beta+\beta_0) \geq M(1-\xi_i)$
 - Lead to different solutions; second is "standard"
 - In second formulation, ξ_i is the proportional amount by which prediction $f(x_i) = x_i'\beta + \beta_0$ is on the wrong side of the margin
 - Bound $\sum \xi \le \text{constant} \rightarrow \text{bound this proportion}$
 - Can show that, equivalently, $\min_{\beta,\beta_0} ||\beta|| \text{ s.t. } y_i(x_i'\beta+\beta_0) \geq 1, i=1\dots,N$ $\xi_i \geq 0, \sum \xi_i \leq \text{constant}$

Computation

- Convex optimization problem
 - Quad. optimization; linear inequality constraints
 - Solution: quadratic programming
- Equivalent form

$$\min_{\beta,\beta_0} \frac{1}{2} ||\beta||^2 + C \sum_{i=1}^{N} \xi_i \text{ s.t. } \xi_i \ge 0,
y_i(x_i'\beta x_i + \beta_0) \ge 1 - \xi_i, \ i = 1, \dots, N$$

- C: cost parameter replacing the bound constant
- Separable case $C = \infty$
- Solution

$$\widehat{\beta} = \sum_{i=1}^{N} \widehat{\alpha}_i y_i x_i$$

 $-\alpha_i \neq 0$ only for support vectors (i.e. observations on or on the wrong side of the boundary)