Lecture 7: Splines and Generalized Additive Models Computational Statistics

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Lecture 7: Splines and Generalized Additive Models Introduction

Overview

Introduction

Simple approaches Polynomials Step functions

Splines

Regression splines Natural splines Splines for classification

Smoothing splines

- Definition
- Computation
- Nonparametric logistic regression

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Lecture 7: Splines and Generalized Additive Models Introduction

Moving beyond linearity

- Linear models are widely used in econometrics.
- In particular, linear regression, linear discriminant analysis, logistic regression all rely on a linear model.
- It is extremely unlikely that the true function f(X) is actually linear in X. In regression problems, $f(X) = \mathbb{E}(Y|X)$ will typically be nonlinear and nonadditive in X, and representing f(X) by a linear model is usually a convenient, and sometimes a necessary, approximation.
 - Convenient because a linear model is easy to interpret, and is the first-order Taylor approximation to f(X).
 - Sometimes necessary, because with N small and/or p large, a linear model might be all we are able to fit to the data without overfitting.
- Likewise in classification, it is usually assumed that some monotone transformation of $\mathbb{P}(Y = 1|X)$ is linear in X. This is inevitably an approximation.

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Linear basis expansion

- The core idea in this chapter is to augment/replace the vector of inputs X with additional variables, which are transformations of X, and then use linear models in this new space of derived input features.
- Denote by $h_m(X) : \mathbb{R}^p \to \mathbb{R}$ the *m*-th transformation of *X*, m = 1, ..., M. We then model

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

a linear basis expansion in X.

 The beauty of this approach is that once the basis functions h_m have been determined, the models are linear in these new variables, and the fitting proceeds as for linear models.

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Popular choices for basis functions h_m

Some simple and widely used examples of the h_m are the following:

- $h_m(X) = X_m$, m = 1, ..., p recovers the original linear model.
- h_m(X) = X_j² or h_m(X) = X_jX_k allows us to augment the inputs with polynomial terms to achieve higher-order Taylor expansions. Note, however, that the number of variables grows exponentially in the degree of the polynomial. A full quadratic model in p variables requires O(p²) square and cross-product terms, or more generally O(p^d) for a degree-d polynomial.
- h_m(X) = log(X_j), √X_j, ... permits other nonlinear transformations of single inputs. More generally one can use similar functions involving several inputs, such as h_m(X) = ||X||.
- $h_m(X) = I(L_m \le X_k < U_m)$, an indicator for a region of X_k . By breaking the range of X_k up into M_k such nonoverlapping regions results in a model with a piecewise constant contribution for X_k .

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Discussion

- Sometimes the problem at hand will call for particular basis functions h_m , such as logarithms or power functions.
- More often, however, we use the basis expansions as a device to achieve more flexible representations for f(X).
- Polynomials are an example of the latter, although they are limited by their global nature – tweaking the coefficients to achieve a functional form in one region can cause the function to flap about madly in remote regions.



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Polynomials

Fitting polynomials

- In most of this lecture, we assume p = 1.
- Create new variables $h_1(X) = X$, $h_2(X) = X^2$, $h_3(X) = X^3$, etc. and then do multiple linear regression on the transformed variables.
- We either fix the degree *d* at some reasonably low value, else use cross-validation to choose *d*.
- Polynomials have unpredictable tail behavior very bad for extrapolation.



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Example in R

```
x = seq(0, 10, 0.5)
n < -length(x)
y1=x[1:10]+2*cos(x[1:10])+2*rnorm(10)
xtest<- seq(-2,12,0.01)
ftest<- xtest+2*cos(xtest)</pre>
d<-3
plot(x1,y1,xlim=c(-2,12),ylim=c(-5,15),
                   main=paste('degree = ',as.character(d)))
for(i in 1:10){
    y2=x[11:21]+2*cos(x[11:21])+2*rnorm(11)
    points(x[11:21],y2,pch=i+1)
    v < -c(v1, v2)
    reg<-lm(y ~ poly(x,degree=d))</pre>
    ypred<-predict(reg,newdata=data.frame(x=xtest),interval="c'</pre>
    lines(xtest,ypred[,"fit"],lty=1)
}
```

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Simple approaches

Polynomials

Result, d = 2

degree = 3





Simple approaches

Polynomials

Result, d = 3

degree = 5





Simple approaches

Polynomials

Result, d = 4

degree = 9





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Step Functions

 Another way of creating transformations of a variable is to cut the variable into distinct regions.

$$h_1(X) = I(X < \xi_1), h_2(X) = I(\xi_1 \le X < \xi_2), \dots,$$

 $h_M(X) = I(X \ge \xi_{M-1})$

• Since the basis functions are positive over disjoint regions, the least squares estimate of the model $f(X) = \sum_{m=1}^{M} \beta_m h_m(X)$ is $\hat{\beta}_m = \overline{Y}_m$, the mean of Y in the *m*-th region.



Example in R

library("ISLR")

reg<-lm(wage ~ cut(age, c(18, 25, 50, 65, 90)),data=Wage)
ypred<-predict(reg,newdata=data.frame(age=18:80),interval="c")</pre>



Simple approaches

Step functions

Result





Step functions - continued

- Easy to work with. Creates a series of dummy variables representing each group.
- Useful way of creating interactions that are easy to interpret. For example, interaction effect of Year and Age:

$$I(Year < 2005) \cdot Age, I(Year \ge 2005) \cdot Age$$

would allow for different linear functions in each age category.

• Choice of cutpoints or knots can be problematic. For creating nonlinearities, smoother alternatives such as splines are available.



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Piecewise Polynomials

 Instead of a single polynomial in X over its whole domain, we can rather use different polynomials in regions defined by knots. E.g. (see figure)

$$y_{i} = \begin{cases} \beta_{01} + \beta_{11}x_{i} + \beta_{21}x_{i}^{2} + \beta_{31}x_{i}^{3} + \epsilon_{i} & \text{if } x_{i} < \xi, \\ \beta_{02} + \beta_{12}x_{i} + \beta_{22}x_{i}^{2} + \beta_{32}x_{i}^{3} + \epsilon_{i} & \text{if } x_{i} \ge \xi, \end{cases}$$

- Better to add constraints to the polynomials, e.g. continuity.
- Splines have the "maximum" amount of continuity.



Splines

Regression splines





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Linear Splines

- A linear spline with knots at ξ_k , k = 1, ..., K is a piecewise linear polynomial continuous at each knot.
- The set of linear splines with fixed knots is a vector space.
- The number of degrees of freedom is 2(K+1) K = K+2. We can thus decompose linear splines on a basis of K + 2 basis functions,

$$y = \sum_{m=1}^{K+2} \beta_m h_m(x) + \epsilon.$$

• The basis functions can be chosen as

$$egin{aligned} h_1(x) &= 1 \ h_2(x) &= x \ h_{k+2}(x) &= (x-\xi_k)_+, \quad k = 1, \dots, K, \end{aligned}$$

where $(\cdot)_+$ denotes the positive part, i.e., $(x - \xi_k)_+ = x - \xi_k$ if $x > \xi_k$ and $(x - \xi_k)_+ = 0$ otherwise.



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Cubic Splines

- A cubic spline with knots at ξ_k, k = 1,..., K is a piecewise cubic polynomial with continuous derivatives up to order 2 at each knot.
- Enforcing one more order of continuity would lead to a global cubic polynomial.
- Again, the set of cubic splines with fixed knots is a vector space, and the number of degrees of freedom is 4(K + 1) 3K = K + 4. We can thus decompose cubic splines on a basis of K + 4 basis functions,

$$y = \sum_{m=1}^{K+4} \beta_m h_m(x) + \epsilon.$$

• We can choose truncated power basis functions,

$$h_k(x) = x^{k-1}, \quad k = 1, \dots, 4,$$

 $h_{k+4}(x) = (x - \xi_k)^3_+, \quad k = 1, \dots, K.$



order-M splines

- More generally, an order-M spline with knots ξ_k , k = 1, ..., K is a piecewise-polynomial of order M 1, which has continuous derivatives up to order M 2.
- A cubic spline has M = 4. A piecewise-constant function is an order-1 spline, while a continuous piecewise linear function is an order-2 spline.
- The general form for the truncated-power basis set is

$$h_k(x) = x^{k-1}, \quad k = 1, \dots, M,$$

 $h_{k+M}(x) = (x - \xi_k)_+^{M-1}, \quad k = 1, \dots, K.$

- It is claimed that cubic splines are the lowest-order spline for which the knot-discontinuity is not visible to the human eye. There is seldom any good reason to go beyond cubic-splines.
- In practice the most widely used orders are M = 1, 2 and 4.



Splines in R

```
library('splines')
fit<-lm(wage~bs(age,5),data=Wage)</pre>
```

ypred<-predict(fit,newdata=data.frame(age=18:80),interval="c")</pre>

```
plot(Wage$age,Wage$wage,cex=0.5,xlab="age",ylab="wage")
lines(18:80,ypred[,"fit"],lty=1,col="blue",lwd=2)
lines(18:80,ypred[,"lwr"],lty=2,col="blue",lwd=2)
lines(18:80,ypred[,"upr"],lty=2,col="blue",lwd=2)
```

- By default, degree=3, and the intersect is not included in the basis functions.
- The number of knots is df-degree. If not specified, the knots are placed at quantiles.

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Result





B-spline basis

- Since the space of spline functions of a particular order and knot sequence is a vector space, there are many equivalent bases for representing them (just as there are for ordinary polynomials.)
- While the truncated power basis is conceptually simple, it is not too attractive numerically: powers of large numbers can lead to severe rounding problems.
- In practice, we often use another basis: the B-spline basis, which allows for efficient computations even when the number of knots *K* is large (each basis function has a local support).



B-spline basis Construction

- Before we can get started, we need to augment the knot sequence.
- Let $\xi_0 < \xi_1$ and $\xi_K < \xi_{K+1}$ be two boundary knots, which typically define the domain over which we wish to evaluate our spline. We now define the augmented knot sequence τ such that

•
$$\tau_1 \leq \tau_2 \leq \ldots \leq \tau_M \leq \xi_0$$

•
$$\tau_{j+M} = \xi_j, \ j = 1, \dots, K$$

- $\xi_{K+1} \leq \tau_{K+M+1} \leq \tau_{K+M+2} \leq \ldots \leq \tau_{K+2M}$.
- The actual values of these additional knots beyond the boundary are arbitrary, and it is customary to make them all the same and equal to ξ_0 and ξ_{K+1} , respectively.



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Lecture 7: Splines and Generalized Additive Models Splines Regression splines B-spline basis Construction – Continued

Denote by B_{i,m}(x) the *i*th B-spline basis function of order m for the knot-sequence τ, m ≤ M. They are defined recursively in terms of divided differences as follows:

$$B_{i,1}(x) = egin{cases} 1 & ext{if } au_i \leq x < au_{i+1} \ 0 & ext{otherwise} \end{cases}$$

for $i = 1, \ldots, K + 2M - 1$. (By convention, $B_{i,1} = 0$ if $\tau_i = \tau_{i+1}$).

$$B_{i,m} = \frac{x - \tau_i}{\tau_{i+m-1} - \tau_i} B_{i,m-1}(x) + \frac{\tau_{i+m} - x}{\tau_{i+m} - \tau_{i+1}} B_{i+1,m-1}(x)$$

for
$$i = 1, ..., K + 2M - m$$
.

 Thus with M = 4, B_{i,4}, i = 1,..., K + 4 are the K + 4 cubic B-spline basis functions for the knot sequence ξ.



Properties

- The B-splines span the space of cubic splines for the knot sequence ξ .
- They have local support and they are nonzero on an interval spanned by M + 1 knots (see next slide).



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Sequence of B-splines up to order 4 with 10 knots evenly spaced from 0 to 1





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Variance of splines beyond the boundary knots

- We know that the behavior of polynomials fit to data tends to be erratic near the boundaries, and extrapolation can be dangerous.
- These problems are exacerbated with splines. The polynomials fit beyond the boundary knots behave even more wildly than the corresponding global polynomials in that region.



Splines

Natural splines

Example





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Explanation of the previous figure

- Pointwise variance curves for four different models, with X consisting of 50 points drawn at random from U[0, 1], and an assumed error model with constant variance.
- The linear and cubic polynomial fits have 2 and 4 df, respectively, while the cubic spline and natural cubic spline each have 6 df.
- The cubic spline has two knots at 0.33 and 0.66, while the natural spline has boundary knots at 0.1 and 0.9, and four interior knots uniformly spaced between them.



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Natural cubic spline

- A natural cubic spline adds additional constraints, namely that the function is linear beyond the boundary knots.
- This frees up four degrees of freedom (two constraints each in both boundary regions), which can be spent more profitably by putting more knots in the interior region.
- There will be a price paid in bias near the boundaries, but assuming the function is linear near the boundaries (where we have less information anyway) is often considered reasonable.



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Example





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Natural cubic spline basis

- A natural cubic spline with K knots has K degrees of freedom: it can ve represented by K basis functions.
- One can start from a basis for cubic splines, and derive the reduced basis by imposing the boundary constraints. For example, starting from the truncated power series basis,

$$f(X) = \sum_{j=0}^{3} \beta_j X^j + \sum_{k=1}^{K} \theta_k (X - \xi_k)^3_+,$$

the contraints f''(X) = 0 and $f^{(3)}(X) = 0$ for $X < \xi_1$ and $X > \xi_K$ lead to the conditions

$$\beta_2 = \beta_3 = 0, \quad \sum_{k=1}^{K} \theta_k = 0, \quad \sum_{k=1}^{K} \xi_k \theta_k = 0$$

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Natural cubic spline basis - continued

 These conditions are automatically satisfied by choosing the following basis,

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

 $N_{k+2}(X) = d_k(X) - d_{K-1}(X), \quad k = 1, \dots, K-2$

with

$$d_{k} = \frac{(X - \xi_{k})_{+}^{3} - (X - \xi_{K})_{+}^{3}}{\xi_{K} - \xi_{k}}$$



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Example in R

```
fit1<-lm(y ~ ns(x,df=5))
fit2<-lm(y ~ bs(x,df=5))
```

ypred1<-predict(fit1,newdata=data.frame(x=xtest),interval="c")
ypred2<-predict(fit2,newdata=data.frame(x=xtest),interval="c")</pre>

```
plot(x,y,xlim=range(xtest))
lines(xtest,ftest)
lines(xtest,ypred1[,"fit"],lty=1,col="red",lwd=2)
lines(xtest,ypred2[,"fit"],lty=1,col="blue",lwd=2)
```



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Using splines with logistic regression

- Until now, we have discussed regression problems. However, splines can also be used when the response variable is qualitative.
- Consider, for instance, natural splines with K knots. For binary classification, we can fit the logistic regression model,

$$\log \frac{\mathbb{P}(Y=1 \ X=x)}{\mathbb{P}(Y=0 \ X=x)} = f(x)$$

with $f(x) = \sum_{k=1}^{K} \beta_k N_k(x)$.

- Once the basis functions have been defined, we just need to estimate coefficients β_k using a standard logistic regression procedure.
- A smooth estimate of the conditional probability $\mathbb{P}(Y = 1|x)$ can then be used for classification or risk scoring.



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Example in R

class<-glm(I(wage>250) ~ ns(age,3),data=Wage,family='binomial')
proba<-predict(class,newdata=data.frame(age=18:80),type='response')</pre>

plot(18:80,proba,xlab="age",ylab="P(wage>250)",type="1") ii<-which(Wage\$wage>250) points(Wage\$age[ii],rep(max(proba),length(ii)),cex=0.5) points(Wage\$age[-ii],rep(0,nrow(Wage)-length(ii)),cex=0.5)



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Smoothing splines Problem formulation

- Here we discuss a spline basis method that avoids the knot selection problem completely by using a maximal set of knots. The complexity of the fit is controlled by regularization.
- Problem: among all functions f(x) with two continuous derivatives, find one that minimizes the penalized residual sum of squares

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

where λ is a fixed smoothing parameter.

The first term measures closeness to the data, while the second term penalizes curvature in the function, and λ establishes a tradeoff between the two. Special cases: λ = 0 (no constraint on f) and λ = ∞ (f has to be linear).

Smoothing splines Solution

- It can be shown that this problem has an explicit, finite-dimensional, unique minimizer which is a natural cubic spline with knots at the unique values of the x_i , i = 1, ..., N.
- At face value it seems that the family is still over-parametrized, since there are as many as *N* knots, which implies *N* degrees of freedom. However, the penalty term translates to a penalty on the spline coefficients, which are shrunk some of the way toward the linear fit.
- The solution is thus of the form

$$f(x) = \sum_{j=1}^{N} N_j(x)\theta_j,$$

where the $N_j(x)$ are an *N*-dimensional set of basis functions for representing this family of natural splines.



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Computation

• The criterion can be written as

$$RSS(\theta, \lambda) = (\boldsymbol{y} - \boldsymbol{N}\theta)^{T} (\boldsymbol{y} - \boldsymbol{N}\theta) + \lambda \theta^{T} \boldsymbol{\Omega}_{N} \theta,$$

where $\{\boldsymbol{N}\}_{ij} = N_j(x_i)$ and $\{\boldsymbol{\Omega}_N\}_{jk} = \int N_j''(t)N_k''(t)dt$.

The solution is

$$\widehat{\theta} = (\boldsymbol{N}^{\mathsf{T}} \boldsymbol{N} + \lambda \boldsymbol{\Omega}_{N})^{-1} \boldsymbol{N}^{\mathsf{T}} \boldsymbol{y},$$

a generalized ridge regression.

• The fitted smoothing spline is given by

$$\widehat{f}(x) = \sum_{j=1}^{N} N_j(x) \widehat{\theta}_j.$$

In practice, when N is large, we can use only a subset of the N interior knots (rule of thumb: number of knots proportional to log N).



Degrees of freedom

Denote by *f* the *N*-vector of fitted values *f*(*x_i*) at the training predictors *x_i*. Then,

$$\widehat{\boldsymbol{f}} = \boldsymbol{N}\widehat{ heta} = (\boldsymbol{N}^{\mathsf{T}}\boldsymbol{N} + \lambda \boldsymbol{\Omega}_N)^{-1} \boldsymbol{N}^{\mathsf{T}} \boldsymbol{y} = \boldsymbol{S}_\lambda \boldsymbol{y}$$

- As matrix *S_λ* does not depend on *y*, the smoothing spline is a linear smoother.
- In the case of cubic spline with knot sequence ξ and, we have

$$\widehat{\boldsymbol{f}} = \boldsymbol{B}_{\xi}\widehat{\theta} = (\boldsymbol{B}_{\xi}^{\mathsf{T}}\boldsymbol{B}_{\xi})^{-1}\boldsymbol{B}_{\xi}^{\mathsf{T}}\boldsymbol{y} = \boldsymbol{H}_{\xi}\boldsymbol{y},$$

where \boldsymbol{B}_{ξ} is the $N \times M$ matrix of basis functions. The degrees of freedom is $M = \text{trace}(\boldsymbol{H}_{\xi})$.

 By analogy, the effective degrees of freedom of a smoothing spline is defined as

$$\mathsf{df}_{\lambda} = \mathsf{trace}(\boldsymbol{S}_{\lambda})$$



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Computation

Selection of smoothing parameters

- As $\lambda \to 0$, df_{λ} $\to N$ and $S_{\lambda} \to I$. As $\lambda \to \infty$, df_{λ} $\to 2$ and $S_{\lambda} \to H$, the hat matrix for linear regression on x.
- Since df_λ is monotone in λ, we can invert the relationship and specify λ by fixing df_λ (this can be achieved by simple numerical methods). Using df in this way provides a uniform approach to compare many different smoothing methods.
- The leave-one-out (LOO) cross-validated error is given by

$$RSS_{cv}(\lambda) = \sum_{i=1}^{N} (y_i - \widehat{f}_{\lambda}^{(-i)}(x_i))^2 = \sum_{i=1}^{N} \left[\frac{y_i - \widehat{f}_{\lambda}(x_i)}{1 - \{\boldsymbol{S}_{\lambda}\}_{ii}} \right]^2$$



Smoothing splines in R

```
ss1<-smooth.spline(x,y,df=3)
ss2<-smooth.spline(x,y,df=15)
ss<-smooth.spline(x,y)</pre>
```

```
plot(x,y)
lines(x,ss1$y,col="blue",lwd=2)
lines(x,ss2$y,col="blue",lwd=2,lty=2)
lines(x,ss$y,col="red",lwd=2)
```

> ss\$df 7.459728



Lecture 7: Splines and Generalized Additive Models

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Application to logistic regression

- The smoothing spline problem has been posed in a regression setting. It is typically straightforward to transfer this technology to other domains.
- Here we consider logistic regression with a single quantitative input X. The model is

$$\log \frac{\mathbb{P}(Y=1|X=x)}{\mathbb{P}(Y=0|X=x)} = f(x),$$

which implies

$$\mathbb{P}(Y = 1 | X = x) = \frac{e^{f(x)}}{1 + e^{f(x)}} = p(x).$$



Penalized log-likelihood

• We construct the penalized log-likelihood criterion

$$\ell(f;\lambda) = \sum_{i=1}^{N} [y_i \log p(x_i) + (1 - y_i) \log(1 - p(x_i))] - \frac{1}{2}\lambda \int \{f''(t)\}^2 dt$$
$$= \sum_{i=1}^{N} [y_i f(x_i) - \log(1 + e^{f(x)})] - \frac{1}{2}\lambda \int \{f''(t)\}^2 dt$$

• As before, the optimal *f* is a finite-dimensional natural spline with knots at the unique values of x. We can represent *f* as

$$f(x) = \sum_{j=1}^{N} N_j(x)\theta_j.$$



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Optimization

• We compute the first and second derivatives

$$\frac{\partial \ell(\theta)}{\partial \theta} = \mathbf{N}^{\mathsf{T}} (\mathbf{y} - \mathbf{p}) - \lambda \Omega \theta$$
$$\frac{\partial^2 \ell(\theta)}{\partial \theta \partial \theta^{\mathsf{T}}} = -\mathbf{N}^{\mathsf{T}} \mathbf{W} \mathbf{N} - \lambda \Omega,$$

where **p** is the *N*-vector with elements p(xi), and **W** is a diagonal matrix of weights $p(x_i)(1 - p(x_i))$.

• Parameters θ_j can be estimated using the Newton method,

$$\theta^{new} = \theta^{old} - \left(\frac{\partial^2 \ell(\theta^{old})}{\partial \theta \partial \theta^{\mathsf{T}}}\right)^{-1} \frac{\partial \ell(\theta^{old})}{\partial \theta}$$



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Nonparametric logistic regression in R

library(gam)
class<-gam(I(wage>250) ~ s(age,df=3),data=Wage,family='binomial')
proba<-predict(class,newdata=data.frame(age=18:80),type='response')</pre>

```
plot(18:80,proba,xlab="age",ylab="P(wage>250)",type="l")
ii<-which(Wage$wage>250)
points(Wage$age[ii],rep(max(proba),length(ii)),cex=0.5)
points(Wage$age[-ii],rep(0,nrow(Wage)-length(ii)),cex=0.5)
```



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Motivation

- Regression models play an important role in many data analyses, providing prediction and classification rules, and data analytic tools for understanding the importance of different inputs.
- Although attractively simple, the traditional linear model often fails in these situations: in real life, effects are often not linear.
- Here, we describe more automatic flexible statistical methods that may be used to identify and characterize nonlinear regression effects. These methods are called generalized additive models (GAMs).



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GAM for regression

• In the regression setting, a generalized additive model has the form

$$\mathbb{E}(Y|X_1,X_2,\ldots,X_p) = \alpha + f_1(X_1) + f_2(X_2) + \ldots + f_p(X_p)$$

- As usual X_1, X_2, \ldots, X_p represent predictors and Y is the outcome.
- The f_j's are unspecified smooth (nonparametric) functions.



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GAM for binary classification

 For two-class classification, recall the logistic regression model for binary data discussed previously. We relate the mean of the binary response μ(X) = P(Y = 1|X) to the predictors via a linear regression model and the logit link function:

$$\log \frac{\mu(X)}{1-\mu(X)} = \alpha + \beta_1 X_1 + \ldots + \beta_p X_p$$

• The additive logistic regression model replaces each linear term by a more general functional form

$$\log \frac{\mu(X)}{1-\mu(X)} = \alpha + f_1(X_1) + \ldots + f_p(X_p)$$

where again each f_j is an unspecified smooth function.

• While the nonparametric form for the functions f_j makes the model more flexible, the additivity is retained and allows us to interpret the model in much the same way as before.

GAM: general form

 In general, the conditional mean μ(X) of a response Y is related to an additive function of the predictors via a link function g:

$$g[\mu(X)] = \alpha + f_1(X_1) + \ldots + f_p(X_p)$$

- Examples of classical link functions are the following:
 - $g(\mu) = \mu$ is the identity link, used for linear and additive models for Gaussian response data.
 - g(μ) = logit(μ) as above, or g(μ) = probit(μ), the probit link function, for modeling binomial probabilities. The probit function is the inverse Gaussian cumulative distribution function: probit(μ) = Φ⁻¹(μ).
 - g(µ) = log(µ) for log-linear or log-additive models for Poisson count data.



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Mixing linear and nonlinear effects, interactions

- We can easily mix in linear and other parametric forms with the nonlinear terms, a necessity when some of the inputs are qualitative variables (factors).
- The nonlinear terms are not restricted to main effects either; we can have nonlinear components in two or more variables, or separate curves in X_i for each level of the factor X_k, e.g.,
 - $g(\mu) = X^T \beta + \sum_k \alpha_k I(V = k) + f(Z)$ a semiparametric model, where X is a vector of predictors to be modeled linearly, α_k the effect for the kth level of a qualitative input V, and the effect of predictor Z is modeled nonparametrically.
 - g(µ) = f(X) + ∑_k g_k(Z)I(V = k) again k indexes the levels of a qualitative input V, and thus creates an interaction term for the effect of V and Z,
 - etc...



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GAMs with natural splines

- If we model each function f_j as a natural spline, then we can fit the resulting model using simple least square (regression) or likelihood maximization algorithm (classification).
- For instance, with natural cubic splines, we have the following GAM:

$$g(\mu) = \sum_{j=1}^{p} \sum_{k=1}^{\mathcal{K}(j)} \beta_{jk} N_k(X_j) + \epsilon,$$

where K(j) is the number of knots for variable j.



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Example in R

```
library("ISLR") # For the Wage data
library("splines")
fit1<-lm(wage ~ ns(year,df=5)+ns(age,df=5)+education,data=Wage)</pre>
```

```
library("gam")
fit2<-gam(wage ~ ns(year,df=5)+ns(age,df=5)+education,data=Wage)
plot(fit2,se=TRUE)</pre>
```



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Result





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GAMs with smoothing splines

• Consider an additive model of the form

$$Y = \alpha + f_1(X_1) + f_2(X_2) + \ldots + f_p(X_p) + \epsilon,$$

where the error term ϵ has mean zero.

• We can specify a penalized sum of squares for this problem,

$$SS(\alpha, f_1, \ldots, f_p) = \sum_{i=1}^N \left(y_i - \alpha - \sum_{j=1}^p f_j(x_{ij}) \right)^2 + \sum_{j=1}^p \lambda_j \int f_j''(t_j)^2 dt_j,$$

where the $\lambda_j \geq 0$ are tuning parameters.

It can be shown that the minimizer of SS is an additive cubic spline model; each of the functions f_j is a cubic spline in the component with knots at each of the unique values of x_{ij}, i = 1, ..., N.

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Unicity of the solution

- Without further restrictions on the model, the solution is not unique.
- The constant α is not identifiable, since we can add or subtract any constants to each of the functions f_j , and adjust α accordingly.
- The standard convention is to assume that $\sum_{i=1}^{N} f_j(x_{ij}) = 0$ for all j the functions average zero over the data.
- It is easily seen that $\widehat{\alpha} = \operatorname{ave}(y_i)$ in this case.
- If in addition to this restriction, the matrix of input values (having ijth entry xij) has full column rank, then SS is a strictly convex criterion and the minimizer is unique.



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Backfitting algorithm

- A simple iterative procedure exists for finding the solution.
- We set $\widehat{\alpha} = \operatorname{ave}(y_i)$, and it never changes.
- We apply a cubic smoothing spline S_j to the targets $\{y_i \hat{\alpha} \hat{f}(x_{ik})\}_{i=1}^N$, as a function of x_{ij} to obtain a new estimate $\hat{f_j}$.
- This is done for each predictor in turn, using the current estimates of the other functions f
 k when computing y − α̂ − Σ{k≠i} f
 k(x{ik}).
- The process is continued until the estimates f_j stabilize.
- This procedure (known as **backfitting**) is grouped cyclic coordinate descent algorithm.



Backfitting algorithm

1 Initialize:
$$\widehat{\alpha} = \operatorname{ave}(y_i), \ \widehat{f_j} = 0, \ \forall i, j.$$

2 Cycle:
$$j = 1, 2, ..., p, 1, 2, ..., p, ...,$$

$$\widehat{f_j} \leftarrow S_j \left[\{ y_i - \widehat{\alpha} - \sum_{k \neq j} \widehat{f_k}(x_{ik}) \}_{i=1}^N \right]$$
$$\widehat{f_j} \leftarrow \widehat{f_j} - \frac{1}{N} \sum_{i=1}^N \widehat{f_j}(x_{ij})$$

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



Example in R

```
library("gam")
fit3<-gam(wage ~ s(year,df=5)+s(age,df=5)+education,data=Wage)
plot(fit3,se=TRUE)</pre>
```



Result







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