Nonparametric Methods II

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OVERVIEW

Polynomial regression and Polynomial splines

Natural splines

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Smoothing splines

(effective) degrees of freedom

LOOCV for linear smoothers

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Polynomial regression and Polynomial splines

Polynomial regression

Suppose p = 1.

Replace simple linear regression

$$y_i = \beta_0 + \beta_1 x_i + \epsilon_i$$

with polynomial regression of degree d

$$y_i = \beta_0 + \beta_1 x_i + \beta_2 x_i^2 + \beta_3 x_i^3 + \ldots + \beta_d x_i^d + \epsilon_i$$

- ▶ Higher degree polynomials give more flexible fit.
- ▶ limitations: impose global structure on the non-linear functions of predictors and leads to global fitting.

To overcome the global restriction, one obvious local method would be to fit polynomials piecewise. For example,

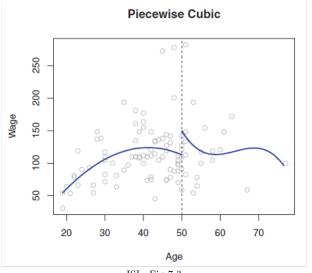
Piecewise cubic (degree=3) with a single cutpoint (knot) at c=50.

$$y_i = \begin{cases} \beta_{01} + \beta_{11}x_i + \beta_{21}x_i^2 + \beta_{31}x_i^3 + \epsilon_i & \text{if } x_i < c \\ \beta_{02} + \beta_{12}x_i + \beta_{22}x_i^2 + \beta_{32}x_i^3 + \epsilon_i & \text{if } x_i \ge c \end{cases}$$

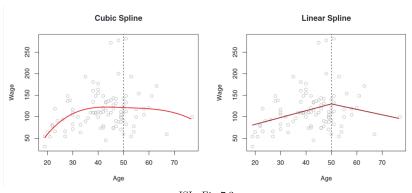
In general, if we place K different knots throughout the range of X, then we will end up fitting K+1 different cubic polynomials.

Polynomial splines

Note the piecewise polynomials lack of continuity at the knots.



ISL: Fig 7.3



ISL: Fig 7.3.

This motivate the polynomial splines:

A d **th-degree spline** f is a piecewise polynomial function of degree d that is continuous and has continuous derivatives of degrees $1, \ldots, d-1$, at its knot points.

- ▶ there are inner nots $t_1 < ... < t_K$
- \triangleright f is a polynomial of degree d on each of the intervals

$$(-\infty,t_1],[t_1,t_2],\ldots,[t_K,\infty)$$

▶ $f^{(j)}$ is continuous at t_1, \ldots, t_K , for each $j = 0, 1, \ldots, d-1$ e.g.: cubic spline, d=3, linear, d=1

A degree-d polynomial spline has the basis given by the "**truncated power basis**":

$$h_j(x) = x^{j-1}, j = 1, \dots, d+1$$

 $h_{d+1+l}(x) = (x-t_l)_+^d, \quad l = 1, \dots, K$

- Unconstrained piecewise polynomial would have d.f. (K+1)(d+1) = Kd + d + K + 1.
- ▶ polynomial splines have d constraints for each of K knots: left with df = d + K + 1.
 - Cubic: d = 3, df = K + 4
 - ▶ Linear: d = 1, df = K + 2

Regression splines use splines as basis functions:

with a set of $d+{\bf K}+1$ basis functions, known as a degree-d spline basis with K knots:

$$b_1(x_i), b_2(x_i), \dots, b_{K+d+1}(x_i)$$

e.g., fit a cubic splines (d = 3):

$$y_i = \beta_1 1 + \beta_2 b_2(x_i) + \beta_3 b_3(x_i) + \dots + \beta_{K+4} b_{K+4}(x_i) + \epsilon_i$$

e.g., with knots $(t_1, t_2), K = 2,$

$$y_i = \beta_1 1 + \beta_2 x_i + \beta_3 x_i^2 + \beta_4 x_i^3 + \beta_5 (x_i - t_1)_+^2 + \beta_6 (x_i - t_2)_+^3 + \epsilon_i$$

Natural splines

Natural splines

Regression splines often give superior results to polynomial regressions.

A problem with regression splines is that the estimates tend to display erratic behavior, i.e., they have high variance, at the boundaries of the input domain. This only gets worse as the polynomial order k gets larger.

natural (cubic) splines are regression splines with additional boundary constraints:

▶ the function is required to be linear at the boundary (in the region where X is smaller than the smallest knot, or larger than the largest knots).

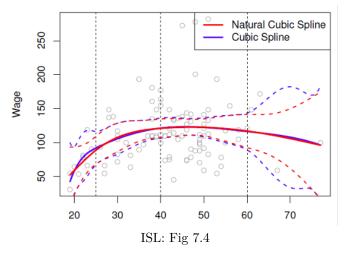
A natural spline of degree d, with knots at $t_1 < \ldots < t_K$, is a piecewise polynomial function f such that

- ▶ f is a polynomial of degree d on each of $[t_1, t_2], \ldots, [t_{K-1}, t_K]$
- ▶ f is a polynomial of degree (d-1)/2 on $(-\infty, t_1]$ and $[t_K, \infty)$
- $ightharpoonup f^{(j)} \ (j=0,\ldots,d-1)$ is continuous at t_1,\ldots,t_K

The dimension of the span of d-th degree natural splines with knots t_1, \ldots, t_K is given by K itself, so independent of the degree.

$$\underbrace{(d+1)(K-1)}_{(1)} + \underbrace{2\left(\frac{(d-1)}{2} + 1\right)}_{(2)} - \underbrace{Kd}_{(3)} = K$$

- (1): the number of free parameters in the interior intervals $[t_1, t_2], \ldots, [t_{K-1}, t_K],$
- (2): the number of free parameters in the exterior intervals $(-\infty, t_1], [t_n, \infty)$,
- ▶ (3): the number of constraints at the knots t_1, \ldots, t_K .



A natural cubic spline with 15 d.f. compared to a degree-15 polynomials (using up to X^{15}).

Better yet, the most popular class of splines are called **B-splines**.

B-Splines

B-Splines

Assume the X is supported on [0,1].

- Let a sequence of knots $\{t_i : 0 = t_0 < \cdots < t_{K+1} = 1\}$ be a partition of the interval [0,1] into K+1 subintervals.
 - ightharpoonup (t_0 , t_{K+1} called boundary knots, the rest K interior knots).
- ► Extended knots are required and the actual values beyond the boundary are arbitrary
 - customary to make them all the same and equal to t_0 and t_{K+1} : $0 = t_{-(q-1)} = \cdots = t_0 < t_1 < \ldots < t_K < t_{K+1} = \cdots = t_{K+q} = 1$
 - ▶ if extended by uniform distance, called "Cardinal B-splines"
- ▶ Denote the B-spline basis functions of order q by $\{B_{1,q}, \ldots, B_{K+q,q}\}$ -defined recursively

For
$$q \ge 2$$
, $B_{i,q}(x) = \frac{x - t_{i-q}}{t_{i-1} - t_{i-q}} B_{i-1,q-1}(x) + \frac{t_i - x}{t_i - t_{i+1-q}} B_{i,q-1}(x)$,
 $i = 1, \dots, q + K$
For $q = 1, B_{i,q}(x) = 1_{[t_{i-1}, t_{i})}(x)$, $i = 1, \dots, K + 1$

$$f(x) = \sum_{j=1}^{J} \beta_j B_{j,q}(x)$$

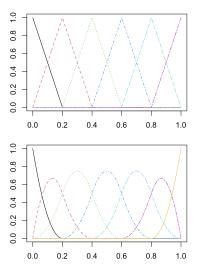
where J = k + q denotes the total number of basis functions.

Some key properties:

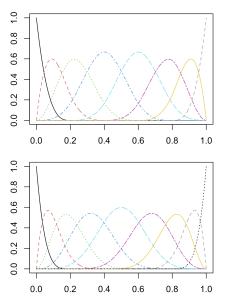
- \triangleright $B_{i,q} \ge 0, i = 1, ..., q + K$; and $B_{i,q} > 0$ on (t_{i-q}, t_i)
- $\sum_{i=1}^{q+K} B_{i,q} = 1$
- At most q adjacent B-splines functions are nonzero at any give $x \in [0, 1]$.

Let
$$b_{J,q}(x) := (B_{1,q}(x), B_{2,q}(x), \dots, B_{J,q}(x))^{\top}$$
 and $\boldsymbol{\beta} := (\beta_1, \dots, \beta_J)^{\top}$, $f(x) = b_{J,q}^{\top}(x)\boldsymbol{\beta}$.

The following are the *B*-splines of order 2 and 3 (K=4 interior knots) even spaced from 0 to 1.



The following are the *B*-splines of order 4 and 5 (K=4 interior knots) even spaced from 0 to 1.



Choosing number and location of knots

Where to place knots?

- ▶ Ideally, place more knots in regions that function varies more, and less in regions of lower variation.
- ▶ In practice, placing knots at uniform quantiles of data.

How many knots?

- number of knots and order of the basis functions count as together
- Try out different number of knots, then use CV to decide order needed
- ▶ Try out different order, then use CV to decide number of knots

Inference about series estimation

Let
$$f(x) = E(Y|X = x) \approx \sum_{j} \beta_{j} B_{j}(x)$$

- ▶ $\{B_j : j = 1, ..., M\}$ are some basis functions (M < n)
- **B** denote the matrix whose (i,j)-th element is given by $B_j(x_i)$.

Then the estimate for β is

$$\hat{\beta} = \left(\mathbf{B}^{\top}\mathbf{B}\right)^{-1}\mathbf{B}^{\top}\mathbf{y}$$

The estimated covariance matrix of $\hat{\beta}$ is

$$\widehat{\operatorname{Var}}(\hat{\beta}) = \left(\mathbf{B}^{\top}\mathbf{B}\right)^{-1}\hat{\sigma}^2$$

$$-\hat{\sigma}^2 = \sum_{i=1}^n (y_i - \hat{f}(x_i))^2 / n.$$

Let $b(x) = (B_1(x), \dots, B_M(x))^{\top}$. The pointwise standard error for f at x is given by

$$\widehat{\operatorname{se}}(\widehat{f}(x)) = \left(b(x)^{\top} \left(\mathbf{B}^{\top} \mathbf{B}\right)^{-1} b(x)\right)^{\frac{1}{2}} \widehat{\sigma}$$

Confidence interval

Note that $\hat{\mathbf{f}} = \mathbf{B} (\mathbf{B}^{\top} \mathbf{B})^{-1} \mathbf{B}^{\top} \mathbf{y} = S_{\mathbf{B}} \mathbf{y}$ (linear smoother).

$$S_{\mathbf{B}} = \mathbf{B} \left(\mathbf{B}^{\top} \mathbf{B} \right)^{-1} \mathbf{B}^{\top}$$

The variance the regression function at the observed data points x_i is given by

$$\widehat{\operatorname{Var}}(\widehat{\mathbf{f}}) = \widehat{\operatorname{Var}}(\mathbf{B}\hat{\beta}) = \hat{\sigma}^2 S_{\mathbf{B}}$$

The pointwise confidence interval at the observed data points can be obtained similarly, using $\widehat{\text{Var}}(\hat{f}_i) = \hat{\sigma}^2[S_{\mathbf{B}}]_{ii}$.

A more accurate estimate $\hat{\sigma}^2 = \frac{1}{n-2\nu_1+\nu_2} \sum_{i=1}^n (y_i - \hat{f}(x_i))^2$, where $\nu_1 = \text{tr}(S), \nu_2 = \text{tr}(SS^\top)$.

Confidence interval (general)

Suppose that $\hat{f}(x) = \sum_i w_i(x)y_i$. The conditional variance is $\sum_i w_i^2(x)\sigma^2(x)$ which can be estimated by $\sum_i w_i^2(x)\hat{\sigma}^2(x)$.

An asymptotic, (biased) pointwise confidence interval is

$$\widehat{f}(x) \pm z_{\alpha/2} \sqrt{\sum_i w_i^2(x) \widehat{\sigma}^2(x)}.$$

 $\hat{\sigma}^2(x)$ can be estimated as: using the regression $e_i^2:=(y_i-\hat{f}(x_i))^2$ v.s. x_i , then take $\hat{\sigma}^2(x_i)=\hat{e}_i^2$.

To eliminate bias, one may want to **undersmooth** the fit \hat{f} , that is M should increase with n at a sufficiently fast rate.

Smoothing splines

Smoothing splines

With inputs x_1, \ldots, x_n lying in an interval [0, 1], the **smoothing** spline estimate f, of a given odd integer degree $d \ge 0$, is defined as

$$\hat{f} = \underset{f}{\operatorname{argmin}} \sum_{i=1}^{n} (y_i - f(x_i))^2 + \lambda \int_0^1 (f^{(m)}(x))^2 dx$$
, where $m = (d+1)/2$

The cubic smoothing splines are given by

$$\hat{f} = \underset{f}{\operatorname{argmin}} \sum_{i=1}^{n} (y_i - f(x_i))^2 + \lambda \int_0^1 f^{(2)}(x)^2 dx.$$

The larger the λ , the smoother f will be.

The solution to above problem is a shrunken version of a natural cubic spline with knots at the unique values of the data x_1, \ldots, x_n .

The solution is $\hat{f}(x) = \sum_{j=1}^{n} \hat{\beta}_{j} N_{j}(x)$, where N_{j} is the set of d-th degree natural splines with knots at x_{1}, \ldots, x_{n} . The problem is

$$\hat{\boldsymbol{\beta}} = \underset{\boldsymbol{\beta} \in \mathbb{R}^n}{\operatorname{argmin}} \| \boldsymbol{y} - \mathbf{N} \boldsymbol{\beta} \|_2^2 + \lambda \boldsymbol{\beta}^\top \Omega \boldsymbol{\beta}$$

where

$$\mathbf{N}_{ij} = N_j(x_i)$$
 and $\Omega_{ij} = \int_0^1 N_i^{(2)}(x) N_j^{(2)}(x) dx$ for $i, j = 1, \dots, n$

The solution is a ridge-type estimator:

$$\hat{\boldsymbol{\beta}} = \left(\mathbf{N}^{\top} \mathbf{N} + \lambda \Omega\right)^{-1} \mathbf{N}^{\top} \mathbf{y}$$

and therefore the fitted values $\hat{\mathbf{f}} = (\hat{f}(x_1), \dots, \hat{f}(x_n))^{\top}$ are

$$\hat{\mathbf{f}} = \mathbf{N} \left(\mathbf{N}^{\top} \mathbf{N} + \lambda \Omega \right)^{-1} \mathbf{N}^{\top} \mathbf{y}$$

$$= \mathbf{N} \left(\mathbf{N}^{\top} \left(I + \lambda \left(\mathbf{N}^{\top} \right)^{-1} \Omega \mathbf{N}^{-1} \right) \mathbf{N} \right)^{-1} \mathbf{N}^{\top} \mathbf{y}$$

$$= (I + \lambda \mathbf{Q})^{-1} \mathbf{y}$$

where $Q = (\mathbf{N}^{\top})^{-1} \Omega \mathbf{N}^{-1}$.

- $ightharpoonup S_{\lambda} = (I + \lambda \mathbf{Q})^{-1} \text{ (called Reinsch form)}$
- ▶ The eigen-decomposition $Q = UDU^{\top}$, $D = diag(d_1, ..., d_n)$
- ▶ U is orthogonal whose columns are basis $\{u\}_{i=1}^n$ (Memmler-Reinsch basis)
- $0 = d_1 = d_2 \le d_3 \le \ldots \le d_n$
 - \triangleright corresponding to the increasingly complex u_i
 - $ightharpoonup d_1 = d_2$ are always equal to 0: linear functions are not penalized

The eigendecomposition of S_{λ} is

$$S_{\lambda} = (I + \lambda Q)^{-1}$$
$$= (I + \lambda U D U^{-1})^{-1}$$
$$= \sum_{j=1}^{n} \rho_{j}(\lambda) u_{j} u_{j}^{\top}$$

The eigenvalues of S_{λ} are given by $\{\rho_j(\lambda) := \frac{1}{1+\lambda d_j} : j=1,\ldots,n\}$.

- $ightharpoonup 0 = d_1 = d_2 \le d_3 \le \ldots \le d_n$
- $ightharpoonup u_1, u_2, \ldots, u_n$ is ordered by increasing complexity

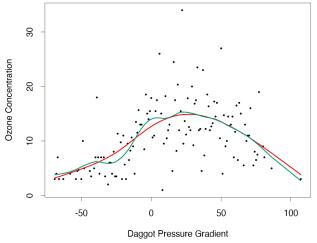
$$\hat{\mathbf{f}} = S_{\lambda} \mathbf{y} = \sum_{j=1}^{n} \frac{u_{j}^{\top} \mathbf{y}}{1 + \lambda d_{j}} u_{j}$$

The smoothing spline operates by decomposing y w.r.t. the orthonormal basis $\{u\}_{i=1}^n$ and differentially shrinking the contributions using $1/(1+\lambda d_i)$.

- The sequence of vectors $u_j \in \mathbb{R}^n$, ordered by decreasing $1/(1+\lambda d_j)$ (as d_j in increasing order), appear to increase in complexity
 - ightharpoonup i.e., u_1, u_2, \ldots, u_n is ordered by increasing complexity

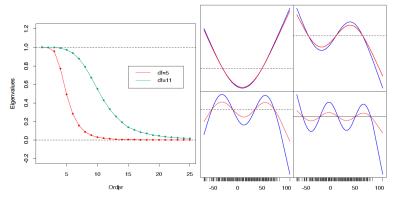
$$S_{\lambda}u_j = \rho_j(\lambda)u_j$$

- Each of the eigenvectors themselves are shrunk by the smoothing spline: the higher the complexity (higher d_j), the more they are shrunk
- ▶ Increasing λ in the smoothing spline estimator tunes out the more wiggly components.



The two fits correspond to different values of the smoothing parameter, chosen to achieve five and eleven effective degrees of freedom.

ESL: Figure 5.7



(left) eigenvalues $\rho_k(\lambda)$: red = large λ , green = small λ . First 25 largest eigenvalues for the two smoothing-spline matrices. The first two are exactly 1, and all are ≥ 0 .

(right) eigenvectors $\rho_k(\lambda)u_k$: red = $\lambda > 0$, blue = $\lambda = 0$. Third to sixth eigenvectors of the spline smoother matrices. In each case, u_k is plotted against x, and as such is viewed as a function of x.

ESL: Figure 5.7.

A comparison between smoothing spline and the regression splines

For regression splines,

$$\hat{\mathbf{f}}_P = B_P (B_P^\top B_P)^{-1} B_P^\top \mathbf{y}$$

- ▶ B_P is $n \times M$ matrix of basic functions evaluated at data points. B_P is full column rank.
- ▶ $H_P = B_P (B_P^\top B_P)^{-1} B_P^\top$, which is symmetric, p.s.d., idempotent and rank (H_P) =M.
- ▶ SVD: $B_P = \tilde{U}\tilde{D}\tilde{V}^{\top}$, and $H_P = \tilde{U}\tilde{U}^{\top}$, where \tilde{U} is $n \times M$.

Columns in \tilde{U} are eigenvectors of H_P corresponding to eigenvalue 1.

$$\hat{\mathbf{f}}_P = H_P y = \sum_{i=1}^M \tilde{u}_i \left\langle \tilde{u}_j, \mathbf{y} \right\rangle$$

Regression Splines :
$$\hat{\mathbf{f}}_P = H_P \mathbf{y} = \sum_{j=1}^M \tilde{u}_j \langle \tilde{u}_j, \mathbf{y} \rangle$$

Smoothing Splines :
$$\hat{\mathbf{f}} = S_{\lambda} \mathbf{y} = \sum_{j=1}^{n} \frac{1}{1 + \lambda d_j} u_j \langle u_j, \mathbf{y} \rangle$$

- regression splines smoother are called **projection smoothers**
 - using projection matrix formed by a subset of bases
- smoothing splines are called shrinking smoothers
 - using a complete basis along with shrinkage

Recall: $\mathbf{X}: n \times p$

OLS:
$$\hat{\mathbf{f}}^{\text{ols}} = P_X \mathbf{y} = \sum_{j=1}^{p} \tilde{u}_j \langle \tilde{u}_j, \mathbf{y} \rangle$$

$$\text{Ridge }: \hat{\mathbf{f}}^{\text{ridge}} = \sum_{j=1}^p \tilde{u}_j \, \frac{\tilde{d}_j^2}{\tilde{d}_j^2 + \lambda} \, \langle \tilde{u}_j, \mathbf{y} \rangle$$

where \tilde{u}_j in the above expression is the j-th column of \tilde{U} , $\mathbf{X} = \tilde{U} \tilde{D} \tilde{V}^{\top}$ (reduced form SVD), $\tilde{d}_1^2 \geq \tilde{d}_2^2 \geq \dots$ Ridge shrinks the most for \tilde{u}_j with small \tilde{d}_j^2 .

(effective) degrees of freedom

(effective) degrees of freedom

More generally, if $y = f(x) + \epsilon$ where $var(\epsilon) = \sigma^2$, the effective d.f. for a linear smoother \hat{f} is

$$\mathrm{df}(\hat{f}) = \frac{\sum_{i=1}^{n} \mathrm{cov}\left(\hat{y}_{i}, y_{i}\right)}{\sigma^{2}} = \frac{\mathrm{trace}(\mathrm{Cov}(\hat{\mathbf{y}}, \mathbf{y}))}{\sigma^{2}}, \text{ conditional on } x_{i}'s$$

Note: the covariance treats only $\{y_i\}_i$ as random ($\{x_i\}_i$ fixed). It measures how strongly the predicted outcomes are associated the actual outcomes.

This definition holds for models that are adaptively fitted to the data (i.e., with tuning parameter chosen adaptively).

For linear smoother $f(x) = \sum_{i=1}^{n} w(x, x_i) y_i$, in the matrix form for the vector of $\mathbf{f} = (f(x_1), \dots, f(x_n))$ to be $\mathbf{f} = W\mathbf{y}$, where W depends on the training data \mathbf{X} , possibly some tuning parameter, but not \mathbf{y} .

The effective degree of freedom for \hat{f} is equal to the trace (W):

$$\mathrm{df}\left(\hat{f}\right) = \sum_{i=1}^{n} w\left(x_{i}, x_{i}\right).$$

- ▶ For the projection linear smoother, $\hat{\mathbf{f}}_p = H_P \mathbf{y}$, $M = \text{trace}(H_P)$ gives the dimension of the projection space.
- ► The (cubic) smoothing spline: $df_{\lambda} = trace(S_{\lambda}) = \sum_{k=1}^{n} \frac{1}{1+\lambda d_{k}}$
 - ▶ as $\lambda \to 0$, $df_{\lambda} \to n$ and $S_{\lambda} \to I$
 - ▶ as $\lambda \to \infty$, $df_{\lambda} \to 2$ and $S_{\lambda} \to P_{\mathbf{X}}$ the projection matrix for linear regression on X.

The effective d.f. of a linear smoother:

$$d.f.(\hat{f}) = \frac{\sum_{i=1}^{n} cov(\hat{y}_i, y_i)}{\sigma^2} = tr(S)$$

examples:

- ridge regression or smoothing splines, $d.f.(\hat{f}) = tr(S_{\lambda})$.
- ightharpoonup k-nearest-neighbor average, $\hat{\mathbf{y}} = S\mathbf{y}$ where $S_{i,j} = w(x_i, x_j)$ where $w(x_i, x_i) = K_k(x_i, x_i) / \sum_{i=1}^n K_k(x_i, x_i)$, and $K_k(x_0, x) = 1(||x - x_0|| \le ||x_{(k)} - x_0||)$, where x_k is the training observation ranked k-th in distance from x_0 . $tr(S) = \sum_{i=1}^{n} w(x_i, x_i) = n/k.$

$$tr(S) = \sum_{i=1}^{n} w(x_i, x_i) = n/k.$$

We can use bootstrap to estimate degrees of freedom (later).

Effective d.f. and the test error

With additive error, and squared error loss:

$$E_{\mathbf{y}}(Err_{\boldsymbol{\tau},in}) = E_{\mathbf{y}}(\overline{err}) + \frac{2}{n} \sum_{i=1}^{n} Cov(\hat{y}_i, y_i)$$

Note the expectation and covariance is w.r.t \mathbf{y} , with $\{x_i\}$ being conditioned upon.

in-sample test error:
$$\operatorname{Err}_{\boldsymbol{\tau}, \text{in}} = \frac{1}{n} \sum_{i=1}^{n} \operatorname{E}_{Y_0} \left(L\left(Y_{0,i}, \hat{f}\left(x_i\right)\right) \mid \boldsymbol{\tau} \right)$$

here $Y_{0,i}$ is new response at x_i ,

training error:
$$\overline{err}_{\tau} = \frac{1}{n} \sum_{i=1}^{n} L\left(y_{i}, \hat{f}\left(x_{i}\right)\right)$$

Although ${\rm Err}_{\tau,~{\rm in}}~$ is not often the direct interest and is different from EPE. However, the comparison of in-sample error is convenient and often leads to efficient model selection.

To estimate in-sample test error, we can use

$$\frac{1}{n}\sum_{i=1}^{n}(y_i-\hat{y}_i)^2+\frac{2\hat{\sigma}^2}{n}\hat{\mathrm{df}}(\hat{f})$$

- Estimate for df from bootstrap or analytic form.
- \triangleright Estimate for σ^2 is needed.

Suppose our linear smoother of interest depends on a tuning parameter α , and express this as $\hat{f}_{\alpha} = S_{\alpha}y$. Then we could choose the tuning parameter α to minimize the estimated test error, as in

$$\hat{\alpha} = \underset{\alpha}{\operatorname{argmin}} \frac{1}{n} \|y - S_{\alpha}y\|_{2}^{2} + \frac{2\sigma^{2}}{n} \operatorname{tr}(S_{\alpha})$$

 \triangleright variants include AIC, BIC, and Mallow's C_p ...

LOOCV for linear smoothers

LOOCV for linear smoothers

CV estimates are roughly unbiased for Err (expected prediction error):

$$E(CV(\hat{f})) \approx Err.$$

Leave-one-out CV

$$CV(\hat{f}) = \frac{1}{n} \sum_{i=1}^{n} (y_i - \hat{f}^{-(i)}(x_i))^2$$

where $\hat{f}^{-(i)}$ is fitted using all training data except the *i*-th observation.

For some linear smoothers $\hat{\mathbf{f}} = S\mathbf{y}$, the LOOCV has a simple form

$$CV(\hat{f}) = \frac{1}{n} \sum_{i=1}^{n} \left(y_i - \hat{f}^{-(i)}(x_i) \right)^2 = \frac{1}{n} \sum_{i=1}^{n} \left(\frac{y_i - \hat{f}(x_i)}{1 - S_{ii}} \right)^2$$

This form holds for linear smoothers based on basis expansion, e.g., projection smoother or shrinkage smoother; and kernel regression and k-NN regression.

GCV (generalized-cross-validation)

A computationally simpler approximation to LOOCV.

$$GCV(\hat{f}) = \frac{1}{n} \sum_{i=1}^{n} \left(\frac{y_i - \hat{f}(x_i)}{1 - \nu/n} \right)^2 = (1 - \nu/n)^{-2} \overline{err}$$

where $\nu = \operatorname{tr}(S)$ is the effective degrees of freedom and \overline{err} is the training error.

This can be of computational advantage in some cases where tr(S) is easier to compute than individual elements S_{ii} .

Tensor-product basis expansion

Tensor-product basis expansion

Suppose $X = (X_1, X_2) \in \mathbb{R}^2$, a basis of functions $h_{1k}(X_1), k = 1, \ldots, M_1$ for representing functions of coordinate X_1 , and likewise a set of M_2 functions $h_{2k}(X_2)$ for coordinate X_2 .

The $M_1 \times M_2$ dimensional **tensor product basis** defined by

$$g_{jk}(X) = h_{1j}(X_1) h_{2k}(X_2), j = 1, \dots, M_1, k = 1, \dots, M_2$$

can be used for representing a two-dimensional function:

$$g(X) = \sum_{j=1}^{M_1} \sum_{k=1}^{M_2} \theta_{jk} g_{jk}(X)$$

A generalization on \mathbb{R}^p is straightforward,

$$g(X) = \sum_{j_1=1}^{M_1} \sum_{j_2=1}^{M_2} \cdots \sum_{j_p=1}^{M_p} \theta_{j_1 \dots j_p} h_{1j_1}(X_1) \cdots h_{pj_p}(X_p)$$

example: tensor-product B-splines

$$f(X_{1},...,X_{p}) = \sum_{j_{1}=1}^{J_{1}} \cdots \sum_{j_{p}=1}^{J_{p}} \theta_{j_{1},...,j_{p}} B_{j_{1},q_{1}}(X_{1}) \cdots B_{j_{p},q_{p}}(X_{p})$$

$$= \sum_{j_{1}=1}^{J_{1}} \cdots \sum_{j_{p}=1}^{J_{p}} \theta_{j_{1},...,j_{p}} B_{j_{1},...,j_{p},q_{1},...,q_{p}}(X_{1},...,X_{p})$$

$$= \mathbf{b}_{J_{1},...,J_{p},q_{1},...,q_{p}}^{\top}(X_{1},...,X_{p}) \boldsymbol{\theta}$$

the elements in $\mathbf{b}_{J_1,\dots,J_p,q_1,\dots,q_p}$ and $\boldsymbol{\theta}$ are ordered lexicographically.

- ▶ **J** denotes the vector (J_1, \ldots, J_p)
- ightharpoonup **q** denotes (q_1, \ldots, q_p) .
- $f(x) = \boldsymbol{b}_{\mathbf{J},\mathbf{q}}^{\top}(x)\boldsymbol{\theta}$

Convergence rates: B-splines

Suppose f is s-times differentiable ($s \ge 1$). Using B-splines of order $q \ge s$ and suitable J, one can obtain the optimal MSE $n^{-2s/(2s+p)}$.

$$\sup_{x} |\hat{f}(x) - f(x)| = O_p\left(\sqrt{\frac{\log n}{n}J^p} + J^{-s}\right)$$

- ightharpoonup J is like the 1/h term
- ▶ $J^p \log n/n$ is the variance term, J^{-s} is the bias term.

with the optimal choice $J \simeq (n/\log n)^{1/(2s+p)}$,

$$\sup_{x} |\hat{f}(x) - f(x)| = O_p\left((\log n/n)^{(s)/(2s+p)} \right)$$

- ▶ $J < (n/\log n)^{1/(2s+p)}$: over-smoothing (relative to the optimal rate)-bias dominant.
- ▶ $J > (n/\log n)^{1/(2s+p)}$: under-smoothing (relative to the optimal rate)-variance dominant.