

1 Introduction and motivations

Regression and classification from an infinite dimensional predictor

Settings

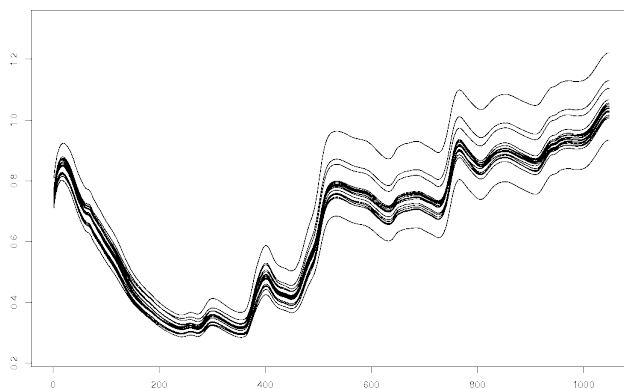
(X, Y) is a random pair of variables where

- $Y \in \{-1, 1\}$ (binary classification problem) or $Y \in \mathbb{R}$
- $X \in (X, \langle \cdot, \cdot \rangle_X)$, an infinite dimensional Hilbert space.

We are given a **learning set** $S_n = \{(X_i, Y_i)\}_{i=1}^n$ of n i.i.d. copies of (X, Y) .

Purpose: Find $\phi_n : X \rightarrow \{-1, 1\}$ or \mathbb{R} , that is universally consistent: *Classification case:* $\lim_{n \rightarrow +\infty} \mathbb{P}(\phi_n(X) \neq Y) = L^*$ where $L^* = \inf_{\phi: X \rightarrow \{-1, 1\}} \mathbb{P}(\phi(X) \neq Y)$ is the **Bayes risk**. *Regression case:* $\lim_{n \rightarrow +\infty} \mathbb{E}([\phi_n(X) - Y]^2) = L^*$ where $L^* = \inf_{\phi: X \rightarrow \mathbb{R}} \mathbb{E}([\phi(X) - Y]^2)$ will also be called the Bayes risk.

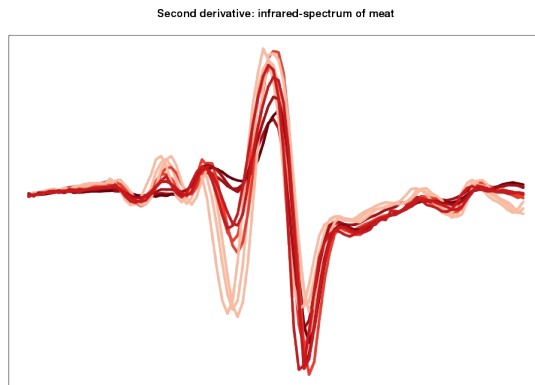
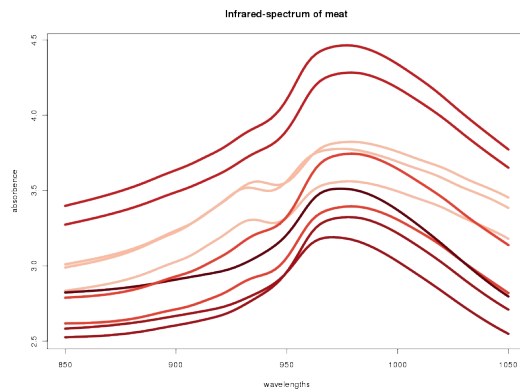
An example



Predicting the **rate of yellow berry in durum wheat** from its **NIR spectrum**.

Using derivatives

Practically, $X^{(m)}$ is often more relevant than X for the prediction.



But $X \rightarrow X^{(m)}$ induces **information loss** and

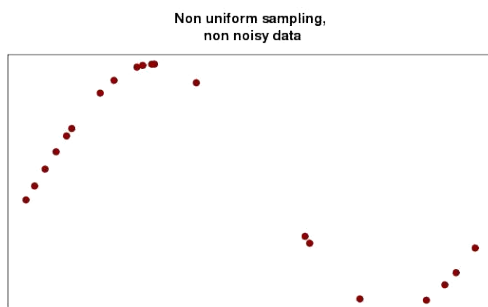
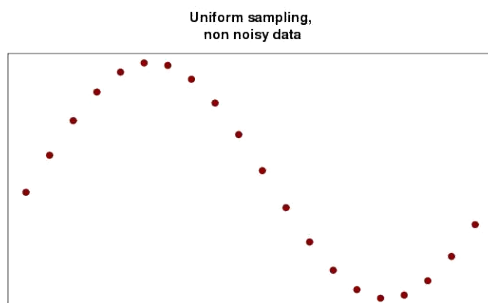
$$\inf_{\phi: D^m \mathcal{X} \rightarrow \{-1,1\}} \mathbb{P}(\phi(X^{(m)}) \neq Y) \geq \inf_{\phi: \mathcal{X} \rightarrow \{-1,1\}} \mathbb{P}(\phi(X) \neq Y) = L^*$$

and

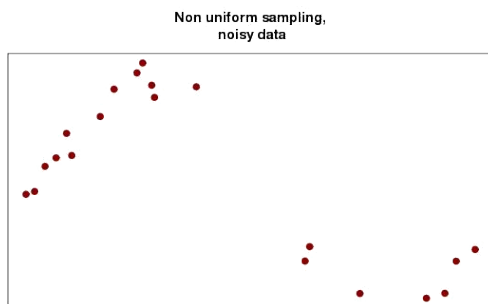
$$\inf_{\phi: D^m \mathcal{X} \rightarrow \mathbb{R}} \mathbb{E}([\phi(X^{(m)}) - Y]^2) \geq \inf_{\phi: \mathcal{X} \rightarrow \mathbb{R}} \mathbb{E}([\phi(X) - Y]^2) = L^*.$$

Sampled functions

Practically, (X_i) are not perfectly known; only a discrete sampling is given: $\mathbf{X}_i^{\tau_d} = (X_i(t))_{t \in \tau_d}$ where $\tau_d = \{t_1^{\tau_d}, \dots, t_{|\tau_d|}^{\tau_d}\}$.



The sampling can be non uniform...



... and the data can be corrupted by noise.
Then, $X_i^{(m)}$ is estimated from $\mathbf{X}_i^{\tau_d}$, by $\widehat{X}_{\tau_d}^{(m)}$, which also induces **information loss**:

$$\inf_{\phi: D^m \mathcal{X} \rightarrow \{-1,1\}} \mathbb{P}(\phi(\widehat{X}_{\tau_d}^{(m)}) \neq Y) \geq \inf_{\phi: D^m \mathcal{X} \rightarrow \{-1,1\}} \mathbb{P}(\phi(X^{(m)}) \neq Y) \geq L^*$$

and

$$\inf_{\phi: D^m \mathcal{X} \rightarrow \mathbb{R}} \mathbb{E} \left(\left[\phi(\widehat{X}_{\tau_d}^{(m)}) - Y \right]^2 \right) \geq \inf_{\phi: D^m \mathcal{X} \rightarrow \mathbb{R}} \mathbb{E} \left(\left[\phi(X^{(m)}) - Y \right]^2 \right) \geq L^*.$$

Purpose of the presentation

Find a classifier or a regression function ϕ_{n,τ_d} built from $\widehat{X}_{\tau_d}^{(m)}$ such that the risk of ϕ_{n,τ_d} **asymptotically reaches** the Bayes risk L^* :

$$\lim_{|\tau_d| \rightarrow +\infty} \lim_{n \rightarrow +\infty} \mathbb{P}(\phi_{n,\tau_d}(\widehat{X}_{\tau_d}^{(m)}) \neq Y) = L^*$$

or

$$\lim_{|\tau_d| \rightarrow +\infty} \lim_{n \rightarrow +\infty} \mathbb{E} \left([\phi_{n,\tau_d}(\widehat{X}_{\tau_d}^{(m)}) - Y]^2 \right) = L^*$$

Main idea: Use a relevant way to estimate $X^{(m)}$ from \mathbf{X}^{τ_d} (by smoothing splines) and combine the consistency of splines with the consistency of a $\mathbb{R}^{|\tau_d|}$ -classifier or regression function.

2 A general consistency result

Basics about smoothing splines I

Suppose that \mathcal{X} is the Sobolev space

$$\mathcal{H}^m = \{h \in L^2_{[0,1]} \mid \forall j = 1, \dots, m, D^j h \text{ exists (weak sense) and } D^m h \in L^2\}$$

equipped with the scalar product

$$\langle u, v \rangle_{\mathcal{H}^m} = \langle D^m u, D^m v \rangle_{L^2} + \sum_{j=1}^m B^j u B^j v$$

where B are m boundary conditions such that $\text{Ker} B \cap \mathbb{P}^{m-1} = \{0\}$.

$(\mathcal{H}^m, \langle \cdot, \cdot \rangle_{\mathcal{H}^m})$ is a **RKHS**: $\exists k_0 : \mathbb{P}^{m-1} \times \mathbb{P}^{m-1} \rightarrow \mathbb{R}$ and $k_1 : \text{Ker} B \times \text{Ker} B \rightarrow \mathbb{R}$ such that

$$\forall u \in \mathbb{P}^{m-1}, t \in [0, 1], \langle u, k_0(t, \cdot) \rangle_{\mathcal{H}^m} = u(t)$$

and

$$\forall u \in \text{Ker} B, t \in [0, 1], \langle u, k_1(t, \cdot) \rangle_{\mathcal{H}^m} = u(t)$$

See **[Berlinet and Thomas-Agnan, 2004]** for further details.

Basics about smoothing splines II

A simple example of boundary conditions:

$$h(0) = h^{(1)}(0) = \dots = h^{(m-1)}(0) = 0.$$

Then,

$$k_0(s, t) = \sum_{k=0}^{m-1} \frac{t^k s^k}{(k!)^2}$$

and

$$k_1(s, t) = \int_0^1 \frac{(t-w)_+^{m-1} (s-w)_+^{m-1}}{(m-1)!} dw.$$

Estimating the predictors with smoothing splines I

Assumption (A1)

- $|\tau_d| \geq m - 1$
- sampling points are distinct in $[0, 1]$
- B^i are linearly independent from $h \rightarrow h(t)$ for all $t \in \tau_d$

[Kimeldorf and Wahba, 1971]: for \mathbf{x}^{τ_d} in $\mathbb{R}^{|\tau_d|}$, $\exists ! \hat{x}_{\lambda, \tau_d} \in \mathcal{H}^m$ solution of

$$\arg \min_{h \in \mathcal{H}^m} \frac{1}{|\tau_d|} \sum_{l=1}^{|\tau_d|} (h(t_l) - \mathbf{x}^{\tau_d})^2 + \lambda \int_{[0,1]} (h^{(m)}(t))^2 dt.$$

and $\hat{x}_{\lambda, \tau_d} = \mathcal{S}_{\lambda, \tau_d} \mathbf{x}^{\tau_d}$ where $\mathcal{S}_{\lambda, \tau_d} : \mathbb{R}^{|\tau_d|} \rightarrow \mathcal{H}^m$.

These assumptions are fulfilled by the previous simple example as long as $0 \notin \tau_d$.

Estimating the predictors with smoothing splines II

$\mathcal{S}_{\lambda, \tau_d}$ is given by:

$$\begin{aligned} \mathcal{S}_{\lambda, \tau_d} &= \omega^T (U(K_1 + \lambda \mathbb{I}_{|\tau_d|}) U^T)^{-1} U (K_1 + \lambda \mathbb{I}_{|\tau_d|})^{-1} \\ &\quad + \eta^T (K_1 + \lambda \mathbb{I}_{|\tau_d|})^{-1} (\mathbb{I}_{|\tau_d|} - U^T (U(K_1 + \lambda \mathbb{I}_{|\tau_d|})^{-1} U (K_1 + \lambda \mathbb{I}_{|\tau_d|})^{-1}) \\ &= \omega^T M_0 + \eta^T M_1 \end{aligned}$$

with

- $\{\omega_1, \dots, \omega_m\}$ is a basis of \mathbb{P}^{m-1} , $\omega = (\omega_1, \dots, \omega_m)^T$ and $U = (\omega_i(t))_{i=1, \dots, m}^T$ $t \in \tau_d$;
- $\eta = (k_1(t, \cdot))_{t \in \tau_d}^T$ and $K_1 = (k_1(t, t'))_{t, t' \in \tau_d}$.

The observations of the **predictor X (NIR spectra)** are then estimated from their sampling \mathbf{X}^{τ_d} by $\hat{\mathbf{X}}_{\lambda, \tau_d}$.

Two important consequences

1. No information loss

$$\inf_{\phi: \mathcal{H}^m \rightarrow \{-1,1\}} \mathbb{P}(\phi(\widehat{X}_{\lambda, \tau_d}) \neq Y) = \inf_{\phi: \mathbb{R}^{|\tau_d|} \rightarrow \{-1,1\}} \mathbb{P}(\phi(\mathbf{X}^{\tau_d}) \neq Y)$$

and

$$\inf_{\phi: \mathcal{H}^m \rightarrow \{-1,1\}} \mathbb{E}\left([\phi(\widehat{X}_{\lambda, \tau_d}) - Y]^2\right) = \inf_{\phi: \mathbb{R}^{|\tau_d|} \rightarrow \{-1,1\}} \mathbb{E}\left([\phi(\mathbf{X}^{\tau_d}) - Y]^2\right)$$

2. Easy way to use derivatives:

$$\begin{aligned} (\mathbf{Q}_{\lambda, \tau_d} \mathbf{u}^{\tau_d})^T (\mathbf{Q}_{\lambda, \tau_d} \mathbf{v}^{\tau_d}) (\mathbf{u}^{\tau_d})^T \mathbf{M}_{\lambda, \tau_d} \mathbf{v}^{\tau_d} (\mathbf{u}^{\tau_d})^T M_0^T W M_0 \mathbf{v}^{\tau_d} + (\mathbf{u}^{\tau_d})^T M_1^T K_1 M_1 \mathbf{v}^{\tau_d} \langle \mathbf{S}_{\lambda, \tau_d} \mathbf{u}^{\tau_d}, \mathbf{S}_{\lambda, \tau_d} \mathbf{v}^{\tau_d} \rangle_{\mathcal{H}^m} &= \langle \widehat{\mathbf{u}}_{\lambda, \tau_d}, \widehat{\mathbf{v}}_{\lambda, \tau_d} \rangle \\ &\simeq \langle \widehat{\mathbf{u}}_{\lambda, \tau_d}^{(m)}, \widehat{\mathbf{v}}_{\lambda, \tau_d}^{(m)} \rangle \end{aligned}$$

where K_1 , M_0 and M_1 have been previously defined and $W = (\langle \omega_i, \omega_j \rangle_{\mathcal{H}^m})_{i,j=1,\dots,m}$. where $\mathbf{M}_{\lambda, \tau_d}$ is symmetric, definite positive. where $\mathbf{Q}_{\lambda, \tau_d}$ is the Choleski triangle of $\mathbf{M}_{\lambda, \tau_d}$: $\mathbf{Q}_{\lambda, \tau_d}^T \mathbf{Q}_{\lambda, \tau_d} = \mathbf{M}_{\lambda, \tau_d}$. **Remark:** $\mathbf{Q}_{\lambda, \tau_d}$ is calculated only from the RKHS, λ and τ_d : it does not depend on the data set.

Classification and regression based on derivatives

Suppose that we know a **consistent classifier or regression function in $\mathbb{R}^{|\tau_d|}$** that is based on $\mathbb{R}^{|\tau_d|}$ scalar product or norm. The **corresponding derivative based classifier or regression function** is given by using the norm induced by $\mathbf{Q}_{\lambda, \tau_d}$:

Example: Nonparametric kernel regression

$$\Psi : u \in \mathbb{R}^{|\tau_d|} \rightarrow \frac{\sum_{i=1}^n T_i K\left(\frac{\|u - U_i\|_{\mathbb{R}^{|\tau_d|}}}{h_n}\right)}{\sum_{i=1}^n K\left(\frac{\|u - U_i\|_{\mathbb{R}^{|\tau_d|}}}{h_n}\right)}$$

where $(U_i, T_i)_{i=1, \dots, n}$ is a learning set in $\mathbb{R}^{|\tau_d|} \times \mathbb{R}$.

$$\begin{aligned} \phi_{n,d} = \Psi \circ \mathbf{Q}_{\lambda, \tau_d} : x \in \mathcal{H}^m &\rightarrow \frac{\sum_{i=1}^n Y_i K \left(\frac{\|\mathbf{Q}_{\lambda, \tau_d} \mathbf{x}^{\tau_d} - \mathbf{Q}_{\lambda, \tau_d} \mathbf{X}_i^{\tau_d}\|_{\mathbb{R}^{|\tau_d|}}}{h_n} \right)}{\sum_{i=1}^n K \left(\frac{\|\mathbf{Q}_{\lambda, \tau_d} \mathbf{x}^{\tau_d} - \mathbf{Q}_{\lambda, \tau_d} \mathbf{X}_i^{\tau_d}\|_{\mathbb{R}^{|\tau_d|}}}{h_n} \right)} \\ &\simeq \frac{\sum_{i=1}^n Y_i K \left(\frac{\|\mathbf{x}^{(m)} - \mathbf{X}_i^{(m)}\|_{L^2}}{h_n} \right)}{\sum_{i=1}^n K \left(\frac{\|\mathbf{x}^{(m)} - \mathbf{X}_i^{(m)}\|_{L^2}}{h_n} \right)} \end{aligned}$$

Remark for consistency

Classification case (approximatively the same is true for regression):

$$\mathbb{P}(\phi_{n, \tau_d}(\widehat{X}_{\lambda, \tau_d}) \neq Y) - L^* = \mathbb{P}(\phi_{n, \tau_d}(\widehat{X}_{\lambda, \tau_d}) \neq Y) - L_d^* + L_d^* - L^*$$

where $L_d^* = \inf_{\phi: \mathbb{R}^{|\tau_d|} \rightarrow \{-1, 1\}} \mathbb{P}(\phi(\mathbf{X}^{\tau_d}) \neq Y)$.

1. For all fixed d ,

$$\lim_{n \rightarrow +\infty} \mathbb{P}(\phi_{n, \tau_d}(\widehat{X}_{\lambda, \tau_d}) \neq Y) = L_d^*$$

as long as the $\mathbb{R}^{|\tau_d|}$ -classifier is consistent because there is a one-to-one mapping between \mathbf{X}^{τ_d} and $\widehat{X}_{\lambda, \tau_d}$.

2. $L_d^* - L^* \leq \mathbb{E} \left(\left| \mathbb{E}(Y | \widehat{X}_{\lambda, \tau_d}) - \mathbb{E}(Y | X) \right| \right)$ with consistency of spline estimate $\widehat{X}_{\lambda, \tau_d}$ and assumption on the regularity of $\mathbb{E}(Y | X = \cdot)$, consistency would be proved. **But** continuity of $\mathbb{E}(Y | X = \cdot)$ is a strong assumption in infinite dimensional case, and is not easy to check.

Spline consistency

Let λ depends on d and denote $(\lambda_d)_d$ the series of regularization parameters. Also introduce $\bar{\Delta}_{\tau_d} := \max\{t_1, t_2 - t_1, \dots, 1 - t_{|\tau_d|}\}$, $\underline{\Delta}_{\tau_d} := \min_{1 \leq i < |\tau_d|} \{t_{i+1} - t_i\}$

Assumption (A2) $\bar{\Delta}_{\tau_d} / \underline{\Delta}_{\tau_d} \leq R$ for all d ;

- $\lim_{d \rightarrow +\infty} |\tau_d| = +\infty$;
- $\lim_{d \rightarrow +\infty} \lambda_d = 0$.

[Ragozin, 1983]: Under (A1) and (A2), $\exists A_{R,m}$ and $B_{R,m}$ such that for any $x \in \mathcal{H}^m$ and any $\lambda_d > 0$,

$$\|\hat{x}_{\lambda_d, \tau_d} - x\|_{L^2}^2 \leq \left(A_{R,m} \lambda_d + B_{R,m} \frac{1}{|\tau_d|^{2m}} \right) \|D^m x\|_{L^2}^2 \xrightarrow{d \rightarrow +\infty} 0$$

Bayes risk consistency**Assumption (A3a)**

$\mathbb{E}(\|D^m X\|_{L^2}^2)$ is finite and $Y \in \{-1, 1\}$.

or

Assumption (A3b)

$\tau_d \subset \tau_{d+1}$ for all d and $\mathbb{E}(Y^2)$ is finite.

Under (A1)-(A3), $\lim_{d \rightarrow +\infty} L_d^* = L^*$.

Proof under assumption (A3a)**Assumption (A3a)**

$\mathbb{E}(\|D^m X\|_{L^2}^2)$ is finite and $Y \in \{-1, 1\}$.

The proof is based on a result of **[Farágó and Györfi, 1975]**:

For a pair of random variables (X, Y) taking their values in $X \times \{-1, 1\}$ where X is an arbitrary metric space and for a series of functions $T_d : X \rightarrow X$ such that

$$\mathbb{E}(\delta(T_d(X), X)) \xrightarrow{d \rightarrow +\infty} 0$$

then $\lim_{d \rightarrow +\infty} \inf_{\phi: X \rightarrow \{-1, 1\}} \mathbb{P}(\phi(T_d(X)) \neq Y) = L^$.*

- T_d is the spline estimate based on the sampling;
- the inequality of **[Ragozin, 1983]** about this estimate is exactly the assumption of Farago and Györfi's Theorem.

Then the result follows.

Proof under assumption (A3b)**Assumption (A3b)**

$\tau_d \subset \tau_{d+1}$ for all d and $\mathbb{E}(Y^2)$ is finite.

Under (A3b), $(\mathbb{E}(Y|\widehat{X}_{\lambda_d, \tau_d}))_d$ is a uniformly bounded martingale and thus converges for the L^1 -norm. Using the consistency of $(\widehat{X}_{\lambda_d, \tau_d})_d$ to X ends the proof.

Concluding result (consistency)

Theorem

Under assumptions (A1)-(A3),

$$\lim_{|\tau_d| \rightarrow +\infty} \lim_{n \rightarrow +\infty} \mathbb{P}(\phi_{n,\tau_d}(\widehat{X}_{\lambda_d, \tau_d}) \neq Y) = L^*$$

and

$$\lim_{|\tau_d| \rightarrow +\infty} \lim_{n \rightarrow +\infty} \mathbb{E}\left(\left[\phi_{n,\tau_d}(\widehat{X}_{\lambda_d, \tau_d}) - Y\right]^2\right) = L^*$$

Proof: For a $\epsilon > 0$, fix d_0 such that, for all $d \geq d_0$, $L_d^* - L^* \leq \epsilon/2$.

Then, by consistency of the $\mathbb{R}^{|\tau_d|}$ -classifier or regression function, conclude.

A practical application to SVM I

Recall that, for a learning set $(U_i, T_i)_{i=1, \dots, n}$ in $\mathbb{R}^p \times \{-1, 1\}$, gaussian SVM is the classifier

$$u \in \mathbb{R}^p \rightarrow \text{Sign} \left(\sum_{i=1}^n \alpha_i T_i e^{-\gamma \|u - U_i\|_{\mathbb{R}^p}^2} \right)$$

where $(\alpha_i)_i$ satisfy the following quadratic optimization problem:

$$\arg \min_w \sum_{i=1}^n |1 - T_i w(U_i)|_+ + C \|w\|_{\mathcal{S}}^2$$

where $w(u) = \sum_{i=1}^n \alpha_i e^{-\gamma \|u - U_i\|_{\mathbb{R}^p}^2}$ and \mathcal{S} is the RKHS associated with the gaussian kernel and C is a **regularization parameter**.

Under suitable assumptions, [Steinwart, 2002] proves the consistency of SVM classifiers.

A practical application to SVM II

Additional assumptions related to SVM: Assumptions (A4) such that $\lim_{n \rightarrow +\infty} n C_n^d = +\infty$ and $C_n^d = \mathcal{O}_n(n^{\beta_d - 1})$ for a $0 < \beta_d < 1/d$.

- For all d , there is a bounded subset of $\mathbb{R}^{|\tau_d|}$, \mathcal{B}_d , such that \mathbf{X}^{τ_d} belongs to \mathcal{B}_d .

Result: Under assumptions (A1)-(A4), the SVM $\phi_{n,d} : x \in \mathcal{H}^m \rightarrow$

$$\text{Sign} \left(\sum_{i=1}^n \alpha_i Y_i e^{-\gamma \|\mathbf{Q}_{\lambda_d, \tau_d} \mathbf{x}^{\tau_d} - \mathbf{Q}_{\lambda_d, \tau_d} \mathbf{X}_i^{\tau_d}\|_{\mathbb{R}^d}^2} \right) \simeq \text{Sign} \left(\sum_{i=1}^n \alpha_i Y_i e^{-\gamma \|x^{(m)} - X_i^{(m)}\|_{L^2}^2} \right)$$

is consistent: $\lim_{|\tau_d| \rightarrow +\infty} \lim_{n \rightarrow +\infty} \mathbb{P}(\phi_{n,\tau_d}(\widehat{X}_{\lambda_d, \tau_d}) \neq Y) = L^*$.

Additional remark about the link between n and $|\tau_d|$

Under suitable (and usual) regularity assumptions on $\mathbb{E}(Y|X = \cdot)$ and if $n \sim \nu^{|\tau_d| \log |\tau_d|}$, the **rate of convergence** of this method is of order $d^{-\frac{2\nu}{2\nu+1}}$ where ν is either equal to m or to a Lipschitz constant related to $\mathbb{E}(Y|X = \cdot)$.

3 Examples

Chosen regression method: Regression with kernel ridge regression

Recall that **kernel ridge regression** in \mathbb{R}^p is given by solving

$$\arg \min_w \sum_{i=1}^n (T_i - w(U_i))^2 + C \|w\|_{\mathcal{S}}^2$$

where \mathcal{S} is a RKHS induced by a given kernel (such as the Gaussian kernel) and $(U_i, T_i)_i$ is a training sample in $\mathbb{R}^p \times \mathbb{R}$.

In the following examples, U_i is either:

- the original (sampled) functions \mathbf{X}_i (viewed as $\mathbb{R}^{|\tau_d|}$ vectors);
- $\mathbf{Q}_{\lambda, \tau_d} \mathbf{X}_i^{\tau_d}$ for derivatives of order 1 or 2.

Example 1: Predicting yellow berry in durum wheat from NIR spectra

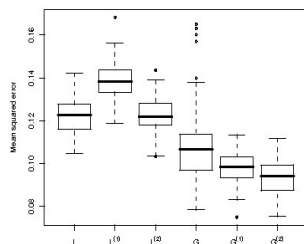
953 wheat samples were analyzed:

- **NIR spectrometry**: 1049 wavelengths regularly ranged from 400 to 2498 nm;
- **Yellow berry**: manual count (%) of affected grains.

Methodology for comparison:

- **Split the data** into train/test sets (50 times);
- **Train** 50 regression functions for the 50 train sets (hyper-parameters were tuned by CV);
- **Evaluate** these regression functions by calculating the **MSE** for the 50 corresponding test sets.

Kernel (SVM)	MSE on test (and sd $\times 10^{-3}$)
Linear (L)	0.122 (8.77)
Linear on derivatives ($L^{(1)}$)	0.138 (9.53)
Linear on second derivatives ($L^{(2)}$)	0.122 (1.71)
Gaussian (G)	0.110 (20.2)
Gaussian on derivatives ($G^{(1)}$)	0.098 (7.92)
Gaussian on second derivatives ($G^{(2)}$)	0.094 (8.35)



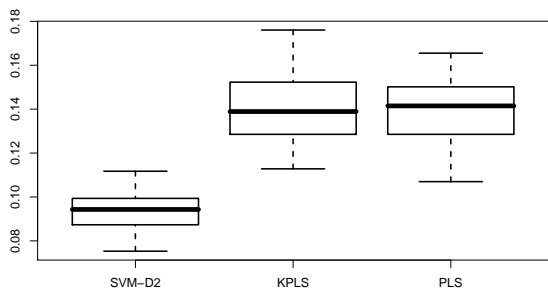
The differences are significant between $G^{(2)} / G^{(1)}$ and between $G^{(1)} / G$.

Comparison with PLS...

	MSE (mean)	MSE (sd)
PLS	0.154	0.012
Kernel PLS	0.154	0.013
KRR splines (reg. D^2)	0.094	0.008

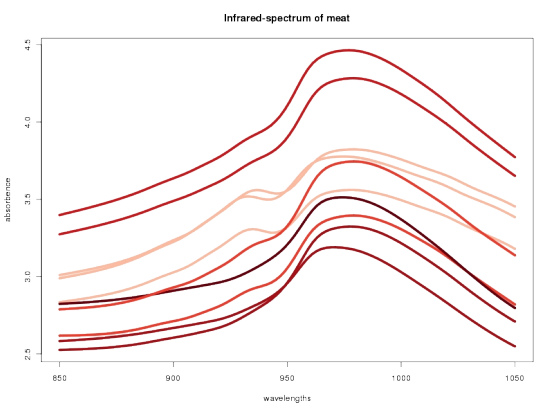
Error decrease: almost

40 %

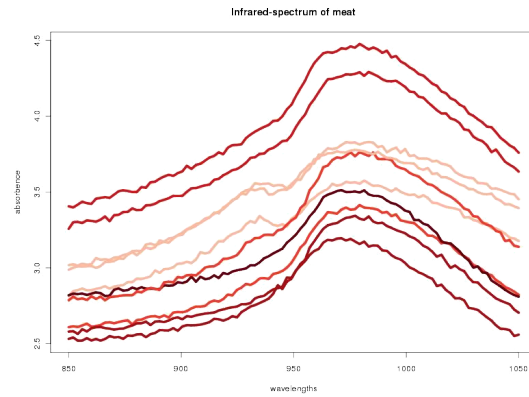


Example 2: Simulated noisy spectra

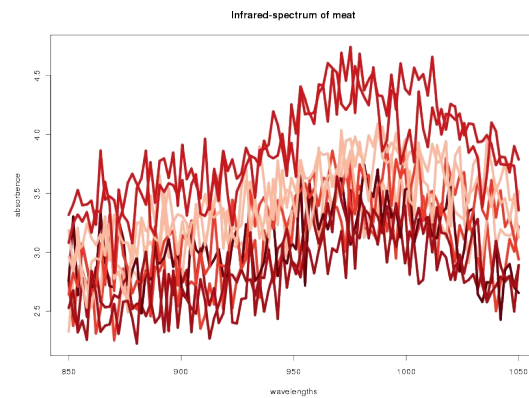
Original data:



Variable to predict: Fat content of pieces of meat. Noisy data: $X_i^b(t) = X_i(t) + \epsilon_{it}$, $\epsilon_{it} \sim \mathcal{N}(0, 0.01)$, i.i.d.:



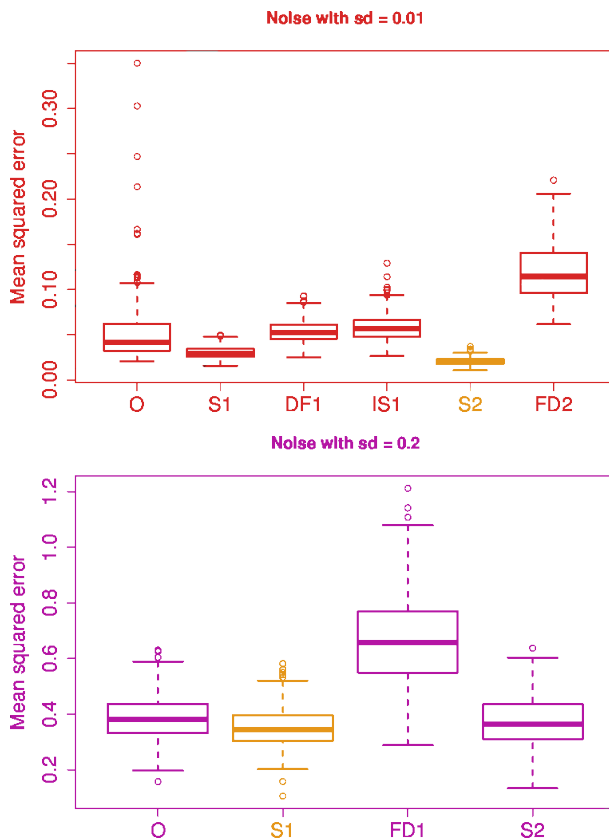
Worse noisy data: $X_i^b(t) = X_i(t) + \epsilon_{it}$, $\epsilon_{it} \sim \mathcal{N}(0, 0.2)$, i.i.d.:



Methodology for comparison

- **Split the data** into train/test sets (250 times);
- **Train** 250 regression functions for the 250 train sets (hyper-parameters were tuned by CV) with the predictors being
 - the original (sampled) functions \mathbf{X}_i (viewed as $\mathbb{R}^{|\tau_d|}$ vectors);
 - $\mathbf{Q}_{\lambda, \tau_d} \mathbf{X}_i^{\tau_d}$ for derivatives of order 1 or 2: **smoothing splines derivatives**;
 - $\mathbf{Q}_{0, \tau_d} \mathbf{X}_i^{\tau_d}$ for derivatives of order 1 or 2: **interpolating splines derivatives**;
 - derivatives of order 1 or 2 evaluated by $\frac{X_i(t_{j+1}) - X_i(t_j)}{t_{j+1} - t_j}$: **finite differences derivatives**;
- **Evaluate** these regression functions by calculating the **MSE** for the 50 corresponding test sets.

Performances



References

References

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Any question?