# Simultaneous adaptation for several criteria using an extended Lepskii principle 

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## Setting: linear regression in Hilbert space

We consider the observation model

$$
Y_{i}=\left\langle f_{\circ}, X_{i}\right\rangle+\xi_{i}
$$

where

- $X_{i}$ takes its values in a Hilbert space $\mathcal{H}$, with $\left\|X_{i}\right\| \leq 1$ a.s.;
- $\xi_{i}$ is a random variable with $\mathbb{E}\left[\xi_{i} \mid X_{i}\right]=0, \mathbb{E}\left[\xi^{2} \mid X_{i}\right] \leq \sigma^{2},|\xi| \leq M$ a.s.;
- $\left(X_{i}, \xi_{i}\right)_{1 \leq i \leq n}$ are i.i.d.

The goal is to estimate $f_{0}$ (in a sense to be specified) from the data.
Note that if $\operatorname{dim}(\mathcal{H})=\infty$, this is essentially a non-parametric model.

## Why this model?

- Hilbert-space valued variables appear in standard models of Functional Data Analysis, where the observed data are modeled (idealized) as function-valued.
- Such models also appear in reproducing kernel Hilbert space (RKHS) methods in machine learning:
- assume observations $X_{i}$ take valued in some space $\mathcal{X}$
- let $\Phi: \mathcal{X} \rightarrow \mathcal{H}$ be a "feature mapping" in a Hilbert space $\mathcal{H}$, and $\widetilde{X}=\Phi(X)$, then one considers the model

$$
Y_{i}=\left\langle f_{0}, \widetilde{X}_{i}\right\rangle+\xi_{i}=\widetilde{f}_{0}\left(X_{i}\right)+\xi_{i},
$$

where $\widetilde{f} \in \widetilde{H}:=\{x \mapsto\langle f, \Phi(x)\rangle ; f \in \mathcal{H}\}$ is a nonparametric model of functions (nonlinear in $x$ !).

- Usually all computations don't require explicit knowledge of $\Phi$ but only access to the kernel $k\left(x, x^{\prime}\right)=\left\langle\Phi(x), \Phi\left(x^{\prime}\right)\right\rangle$.


## Why this model (II) - inverse learning

Of interest is also the inverse learning problem:

- $X_{i}$ takes value in $\mathcal{X}$;
- if $A$ is a linear operator from a Hilbert space $\mathcal{H}$ to a real function space on $\mathcal{X}$;
- inverse regression learning model:

$$
Y_{i}=\left(A f_{\circ}\right)\left(X_{i}\right)+\xi_{i} .
$$

- If $A$ is a Carleman operator (i.e. evaluation functionals $f \mapsto(A f)(x)$ are continuous for all $x$ ), then this can be isometrically reduced to a reproducing kernel learning setting (De Vito, Rosasco, Caponnetto 2006; Blanchard and Mücke, 2017).


## Two notions of risk

We will consider two notions of error (risk) for a candidate estimate $\widehat{f}$ of $f_{\circ}$ :

- Squared prediction error:

$$
\mathcal{E}(\widehat{f}):=\mathbb{E}\left[(\langle\widehat{f}, X\rangle-Y)^{2}\right] .
$$

- The associated (excess error) risk is

$$
\mathcal{E}(\widehat{f})-\mathcal{E}\left(f_{\circ}\right)=\mathbb{E}\left[\left(\left\langle\widehat{f}-f_{\circ}, X\right\rangle\right)^{2}\right]=\left\|\widehat{f}^{*}-f_{\circ}^{*}\right\|_{2, x^{\prime}}^{2}
$$

Reconstruction error risk:

$$
\left\|\widehat{f}-f_{\circ}\right\|_{\mathcal{H}}^{2} .
$$

The goal is to find a suitable estimator $\widehat{f}$ of $f_{0}$ from the data having "optimal" convergence properties with respect to these two risks.

## Finite-dimensional case

- The final dimensional case: $\mathcal{X}=\mathbb{R}^{p}, f_{\circ}$ now denoted $\beta$ 。
- In usual matrix form:

$$
Y=X \beta_{\circ}+\xi .
$$

- $X_{i}^{T}$ form the lines of the $(n, p)$ design matrix $X$
- $Y=\left(Y_{1}, \ldots, Y_{n}\right)^{T}$
- $\xi=\left(\xi_{1}, \ldots, \xi_{n}\right)^{T}$
- "Reconstruction" risk corresponds to $\left\|\beta_{\circ}-\widehat{\beta}\right\|^{2}$.
- Prediction risk corresponds to

$$
\mathbb{E}\left[\left\langle\beta_{\circ}-\widehat{\beta}, X\right\rangle^{2}\right]=\left\|\Sigma^{1 / 2}\left(\beta_{\circ}-\widehat{\beta}\right)\right\|^{2},
$$

where $\Sigma:=\mathbb{E}\left[X X^{\top}\right]$.

- In Hilbert space, same relation with $\Sigma:=\mathbb{E}\left[X \otimes X^{*}\right]$.


## The founding fathers of machine learning


A.M. Legendre

C.F. Gauß

The "ordinary" least squares (OLS) solution:

$$
\widehat{\beta}_{O L S}=\left(\boldsymbol{X}^{\top} \boldsymbol{X}\right)^{-1} \boldsymbol{X}^{\top} \boldsymbol{Y}
$$

## Convergence of OLS in finite dimension

- The "ordinary" least squares (OLS) solution:

$$
\widehat{\beta}_{O L S}=\left(\boldsymbol{X}^{\top} \boldsymbol{X}\right)^{-1} \boldsymbol{X}^{\top} \boldsymbol{Y}
$$

- We want to understand the behavior of $\widehat{\beta}_{O L S}$, when the data size $n$ grows large. Will we be close to the truth $\beta_{0}$ ?
- Recall

$$
\widehat{\beta}_{O L S}=\left(\boldsymbol{X}^{\top} \boldsymbol{X}\right)^{-1} \boldsymbol{X}^{\top} \boldsymbol{Y}=(\underbrace{\frac{1}{n} \boldsymbol{X}^{\top} \boldsymbol{X}}_{:=\widehat{\Sigma}})^{-1}(\underbrace{\frac{1}{n} \boldsymbol{X}^{\top} \boldsymbol{Y}}_{:=\widehat{\gamma}})=\widehat{\Sigma}^{-1} \widehat{\gamma},
$$

- Observe by a vectorial LLN, as $n \rightarrow \infty$ :

$$
\begin{aligned}
& \widehat{\Sigma}:=\frac{1}{n} \boldsymbol{X}^{\top} \boldsymbol{X}=\frac{1}{n} \sum_{i=1}^{n} \underbrace{X_{i} X_{i}^{\top}}_{=: Z_{i}^{\prime}} \longrightarrow \mathbb{E}\left[X_{1} X_{1}^{\top}\right]=: \Sigma ; \\
& \widehat{\gamma}:=\frac{1}{n} \boldsymbol{X}^{\top} \boldsymbol{Y}=\frac{1}{n} \sum_{i=1}^{n} \underbrace{X_{i} Y_{i}}_{=: Z_{i}} \longrightarrow \mathbb{E}\left[X_{1} Y_{1}\right]=\Sigma \beta_{\circ}=: \gamma ;
\end{aligned}
$$

- Hence $\widehat{\beta}=\widehat{\Sigma}^{-1} \widehat{\gamma} \rightarrow \Sigma^{-1} \gamma=\beta_{0} . \quad$ (Assuming $\Sigma$ invertible.)


## From OLS to Hilbert-space regression

- For ordinary linear regression with $\mathcal{X}=\mathbb{R}^{p}$ (fixed $\left.p, n \rightarrow \infty\right)$ :
- LLN implies $\widehat{\beta}_{O L S}\left(=\widehat{\Sigma}^{-1} \widehat{\gamma}\right) \rightarrow \beta_{0}\left(=\Sigma^{-1} \gamma\right)$;
- CLT+Delta Method imply asymptotic normality and convergence in $\mathcal{O}\left(n^{-\frac{1}{2}}\right)$.

How to generalize to $\mathcal{X}=\mathcal{H}$ ?

- Main issue: $\Sigma=\mathbb{E}\left[X \otimes X^{*}\right]$ does not have a continuous inverse. ( $\rightarrow$ ill-posed problem)
- Need to consider a suitable approximation $\zeta(\widehat{\Sigma})$ of $\Sigma^{-1}$ (regularization), where

$$
\widehat{\Sigma}:=\frac{1}{n} \sum_{i=1}^{m} X_{i} \otimes X_{i}^{*}
$$

is the empirical second moment operator.

## Regularization methods

- Main idea: replace $\widehat{\Sigma}^{-1}$ by an approximate inverse, such as
- Ridge regression/Tikhonov:

$$
\widehat{f}_{\text {Ridge }(\lambda)}=\left(\widehat{\Sigma}+\lambda I_{p}\right)^{-1} \widehat{\gamma}
$$

- PCA projection/spectral cut-off: restrict $\widehat{\Sigma}$ on its $k$ first eigenvectors

$$
\widehat{f}_{P C A(k)}=(\widehat{\Sigma})_{\mid k}^{-1} \widehat{\gamma}
$$

- Gradient descent/Landweber Iteration/ $L^{2}$ boosting:

$$
\begin{aligned}
\widehat{f}_{L W(k)} & =\widehat{f}_{L W(k-1)}+\left(\widehat{\gamma}-\widehat{\Sigma} \widehat{f}_{L W(k-1)}\right) \\
& =\sum_{i=0}^{k}(I-\widehat{\Sigma})^{k} \widehat{\gamma}
\end{aligned}
$$

(assuming $\|\widehat{\Sigma}\|_{o p} \leq 1$ ).

## General form spectral linearization

Bauer, Rosasco, Pereverzev 2007

- General form regularization method:

$$
\widehat{f}_{\lambda}=\zeta_{\lambda}(\widehat{\Sigma}) \widehat{\gamma}
$$

for some well-chosen function $\zeta_{\lambda}: \mathbb{R}_{+} \rightarrow \mathbb{R}_{+}$acting on the spectrum and "approximating" the function $x \mapsto x^{-1}$.

- $\lambda>0$ : regularization parameter; $\lambda \rightarrow 0 \Leftrightarrow$ less regularization

Notation of (autoadjoint) functional calculus, i.e.

$$
\widehat{\Sigma}=Q^{T} \operatorname{diag}\left(\mu_{1}, \mu_{2}, \ldots\right) Q \Rightarrow \zeta(\widehat{\Sigma}):=Q^{\top} \operatorname{diag}\left(\zeta\left(\mu_{1}\right), \zeta\left(\mu_{2}\right), \ldots\right) Q
$$

- Examples (revisited):
- Tikhonov: $\zeta_{\lambda}(t)=(t+\lambda)^{-1}$
- Spectral cut-off: $\zeta_{\lambda}(t)=t^{-1} 1\{t \geq \lambda\}$
- Landweber iteration: $\zeta_{k}(t)=\sum_{i=0}^{k}(1-t)^{i}$.


## Assumptions on regularization function

Standard assumptions on the regularization family $\zeta_{\lambda}:[0,1] \rightarrow \mathbb{R}$ are:
(i) There exists a constant $D<\infty$ such that

$$
\sup _{0<\lambda \leq 10<t \leq 1} \sup _{0}\left|t \zeta_{\lambda}(t)\right| \leq D,
$$

(ii) There exists a constant $E<\infty$ such that

$$
\sup _{0<\lambda \leq 10<t \leq 1} \sup _{1} \lambda\left|\zeta_{\lambda}(t)\right| \leq E,
$$

(iii) Qualification: for residual $r_{\lambda}(t):=1-t \zeta_{\lambda}(t)$,

$$
\forall \lambda \leq 1: \quad \sup _{0<t \leq 1}\left|r_{\lambda}(t)\right| t^{v} \leq \gamma_{\nu} \lambda^{v},
$$

holds for $v=0$ and $v=q>0$.

## Structural Assumptions (I)

- Denote $\left(\mu_{i}\right)_{i \geq 1}$ the sequence of positive eigenvalues of $\Sigma$ in nonincreasing order.
- Assumptions on spectrum decay: for $s \in(0,1) ; \alpha>0$ :

$$
\mathbb{I P}^{<}(s, \alpha): \quad \mu_{i} \leq \alpha i^{-\frac{1}{s}}
$$

- This implies quantitative estimates of the "effective dimension"

$$
\mathcal{N}(\lambda):=\operatorname{Tr}\left((\Sigma+\lambda)^{-1} \Sigma\right) \lesssim \lambda^{-s} .
$$

## Structural Assumptions (II)

Denote $\left(\mu_{i}\right)_{i \geq 1}$ the sequence of positive eigenvalues of $\Sigma$ in nonincreasing order.

- Source condition for the signal: for $r>0$, define

$$
\mathrm{SC}(r, R): \quad f_{\circ}=\Sigma^{r} h_{\circ} \text { for some } h_{\circ} \text { with }\left\|h_{\circ}\right\| \leq R,
$$

or equivalently, as a Sobolev-type regularity

$$
\mathbf{S C}(r, R): \quad f_{\circ} \in\left\{f \in \mathcal{H}: \sum_{i \geq 1} \mu_{i}^{-2 r} f_{i}^{2} \leq R^{2}\right\}
$$

where $f_{i}$ are the coefficients of $h$ in the eigenbasis of $\Sigma$.

- Under $(\mathrm{SC})(r, R)$ it is assumed that the qualification $q$ of the regularization method satisfies $q \geq r+\frac{1}{2}$.


## A general upper bound risk estimate

## Theorem

Assume the source condition (SC) $(r, R)$ holds.
If $\lambda$ is such that $\lambda \gtrsim\left(\mathcal{N}(\lambda) \vee \log (\eta)^{2}\right) / n$, then with probability at least $1-\eta$, it holds:

$$
\begin{aligned}
&\left\|(\Sigma+\lambda)^{1 / 2}\left(f_{\circ}-\widehat{f}_{\lambda}\right)\right\|_{\mathcal{H}} \\
& \lesssim \log (\eta)^{2}\left(R \lambda^{r+\frac{1}{2}}+\sigma \sqrt{\frac{\mathcal{N}(\lambda)}{n}}+\frac{1}{n \sqrt{\lambda}}+\mathcal{O}\left(n^{-\frac{1}{2}}\right)\right)
\end{aligned}
$$

This gives rise to estimates in both norms of interest since

$$
\left\|f_{\circ}-\widehat{f}_{\lambda}\right\|_{\mathcal{H}} \leq \lambda^{-\frac{1}{2}}\left\|(\Sigma+\lambda)^{1 / 2}\left(f_{\circ}-\widehat{f}_{\lambda}\right)\right\|_{\mathcal{H}^{\prime}}
$$

and

$$
\left\|f_{\circ}^{*}-\widehat{f}_{\lambda}^{*}\right\|_{L^{2}\left(P_{X}\right)}=\left\|\Sigma^{\frac{1}{2}}\left(f_{\circ}-\widehat{f}_{\lambda}\right)\right\|_{\mathcal{H}} \leq\left\|(\Sigma+\lambda)^{1 / 2}\left(f_{\circ}-\widehat{f}_{\lambda}\right)\right\|_{\mathcal{H}} .
$$

## Upper bound on rates

Optimizing the obtained bound over $\lambda$ (i.e. balancing the main terms) one obtains

## Theorem

Assume r, R, s, $\alpha$ are fixed positive constants and assume $\mathbb{P}_{X Y}$ satisfies $\left(\mathbf{I P}^{<}\right)(s, \alpha)$, (SC) $(r, R)$ and $\|X\| \leq 1,\|Y\| \leq M, \operatorname{Var}[Y \mid X]_{\infty} \leq \sigma^{2}$ a.s. Define

$$
\widehat{\beta}_{n}=\zeta_{\lambda_{n}}(\widehat{\Sigma}) \widehat{\gamma}
$$

using a regularization family $\left(\zeta_{\lambda}\right)$ satisfying the standard assumptions with qualification $q \geq r+\frac{1}{2}$, and the parameter choice rule

$$
\lambda_{n}=\left(R^{2} \sigma^{2} / n\right)^{-\frac{1}{2 r+1+s}}
$$

Then it holds for any $p \geq 1$ :

$$
\begin{aligned}
& \limsup _{n \rightarrow \infty} \mathbb{E}^{\otimes n}\left(\left\|f_{\circ}-\widehat{f}_{\lambda_{n}}\right\|^{p}\right)^{1 / p} / R\left(\frac{\sigma^{2}}{R^{2} n}\right)^{\frac{r}{2 r+1+s}} \leq C_{\mathbf{\Delta}} ; \\
\limsup & \mathbb{E}^{\otimes n}\left(\left\|f_{\circ}^{*}-\widehat{f}_{\lambda_{n}}\right\|_{2, X}^{p}\right)^{1 / p} / R\left(\frac{\sigma^{2}}{R^{2} n}\right)^{\frac{r+1 / 2}{2 r+1+s}} \leq C_{\mathbf{\Delta}} .
\end{aligned}
$$

## Towards adaptivity: existing approaches

- Cross-validation (or hold-out) will yield a tuning of the parameter which is adaptive in the prediction risk, it is based on a unbiased estimate of the risk (URE) principle.
- Standard Lepski's principle parameter selection can be applied for any fixed norm (provided a good estimate of the "variance" term $\sigma \sqrt{\mathcal{N}}(\lambda) / n$ is available)
- Despite the existence of a regularization parameter $\lambda$ being optimal for both norms, there is no guarantee that any (close to) optimal parameter for prediction risk (eg. selected by cross-validation) will be close to optimal in reconstruction risk, or vice-versa.
- We want to construct a simultaneously (for both norms) adaptive data-driven parameter selection.


## Generalized Lepskii's principle

We consider the following "deterministic" assumption to highlight the construction.

## Assumption

Let $\Lambda \subset \mathbb{R}_{+}$be a finite set of candidate regularization parameters,

$$
\Lambda:=\left\{\lambda_{j}, \quad \lambda_{0}>\lambda_{1}>\ldots>\lambda_{m}=\lambda_{\min }>0\right\}
$$

The (known) family of elements of $\mathcal{H},\left(f_{\lambda}\right)_{\lambda \in \Lambda}$, satisfies for any $\lambda \in \Lambda$ :

$$
\left\|(\Sigma+\lambda)^{1 / 2}\left(f_{\circ}-f_{\lambda}\right)\right\|_{\mathcal{H}} \leq C \sqrt{\lambda}(\mathcal{A}(\lambda)+\mathcal{S}(\lambda))
$$

where

- the function $\lambda \in \Lambda \mapsto \mathcal{A}(\lambda) \in \mathbb{R}_{+}$is non-decreasing with $\mathcal{A}(0)=0$ and possibly unknown;
- the function $\lambda \in \Lambda \mapsto \sqrt{\lambda} \mathcal{S}(\lambda) \in \mathbb{R}_{+}$is non-increasing and known.


## Generalized Lepskii's principle (II)

- Set

$$
\begin{array}{r}
\mathcal{M}(\Lambda):=\left\{\lambda \in \Lambda:\left\|\left(\Sigma+\lambda^{\prime}\right)^{1 / 2}\left(f_{\lambda}-f_{\lambda^{\prime}}\right)\right\|_{\mathcal{H}} \leq 4 C \sqrt{\lambda^{\prime}} \mathcal{S}\left(\lambda^{\prime}\right)\right. \\
\left.\forall \lambda^{\prime} \in \Lambda, \text { s.t. } \lambda^{\prime} \leq \lambda\right\} .
\end{array}
$$

- The balancing parameter is given as

$$
\hat{\lambda}:=\max \mathcal{M}(\Lambda) ;
$$

(this quantity is always well-defined since $\lambda_{\text {min }} \in \mathcal{M}(\Lambda)$.)

## Generalized Lepskii's principle: bound

## Theorem

Under the assumptions made previously, if

$$
\lambda_{*}:=\max \{\lambda \in \Lambda: \mathcal{A}(\lambda) \leq \mathcal{S}(\lambda)\},
$$

and $\widehat{\lambda}$ is the parameter choice defined previously, then:

- It holds

$$
\left\|\left(\Sigma+\lambda_{*}\right)^{\frac{1}{2}}\left(f_{0}-f_{\hat{\lambda}}\right)\right\|_{\mathcal{H}} \lesssim \sqrt{\lambda_{*}} \mathcal{S}\left(\lambda_{*}\right) ;
$$

- Assuming it holds $\mathcal{S}\left(\lambda_{k}\right) \leq C_{\mathcal{S}} \mathcal{S}\left(\lambda_{k-1}\right)$ for $k=1, \ldots, m$, then:

$$
\begin{aligned}
\left\|f_{\circ}-f_{\widehat{\lambda}}\right\|_{\mathcal{H}} & \lesssim \min _{\lambda \in \Lambda}(\mathcal{A}(\lambda)+\mathcal{S}(\lambda)) \\
\left\|\Sigma^{\frac{1}{2}}\left(f_{\circ}-f_{\hat{\lambda}}\right)\right\|_{\mathcal{H}} & \lesssim \min _{\lambda \in \Lambda} \sqrt{\lambda}(\mathcal{A}(\lambda)+\mathcal{S}(\lambda)) .
\end{aligned}
$$

## Applying Lepski's principle

Looking at the main error bound obtained earlier, with high probability the assumption

$$
\left\|(\Sigma+\lambda)^{1 / 2}\left(f_{\circ}-f_{\lambda}\right)\right\|_{\mathcal{H}} \leq C \sqrt{\lambda}(\mathcal{A}(\lambda)+\mathcal{S}(\lambda))
$$

is satisfied with

$$
\begin{aligned}
& \mathcal{A}(\lambda):=\left(R \lambda^{r+\frac{1}{2}}+\mathcal{O}\left(n^{-\frac{1}{2}}\right)\right) \\
& \mathcal{S}(\lambda):=\sigma \sqrt{\frac{\mathcal{N}(\lambda)}{n}}+\frac{1}{n \sqrt{\lambda}}
\end{aligned}
$$

## Remaining issues:

- $\Sigma$ is not known;
- $\mathcal{N}(\lambda)=\operatorname{Tr}\left((\Sigma+\lambda)^{-1} \Sigma\right)$ is not known;
- the noise variance $\sigma^{2}$ might not be known (issue ignored for now).


## Replacing $\Sigma, \mathcal{N}(\lambda)$ by empirical quantities

## Proposition

If $\lambda$ is such that $\lambda \gtrsim\left(\mathcal{N}(\lambda) \vee \log (\eta)^{2}\right) / n$, then with probability at least $1-\eta$, it holds:

$$
\left\|(\Sigma+\lambda)^{\frac{1}{2}}(\widehat{\Sigma}+\lambda)^{-\frac{1}{2}}\right\| \lesssim 1+\log \left(\eta^{-1}\right) .
$$

## Proposition

If $\lambda \gtrsim n^{-1}$, it holds with probability at least $1-\eta$, for $\widehat{\mathcal{N}}(\lambda):=\operatorname{Tr}\left(\widehat{\Sigma}(\widehat{\Sigma}+\lambda)^{-1}\right)$ :

$$
\max \left(\frac{\mathcal{N}(\lambda) \vee 1}{\widehat{\mathcal{N}}(\lambda) \vee 1}, \frac{\widehat{\mathcal{N}}(\lambda) \vee 1}{\mathcal{N}(\lambda) \vee 1}\right) \lesssim\left(1+\log \eta^{-1}\right)^{2}
$$

## Fully empirical procedure ( $\sigma, M$ known)

- Put $L:=2 \log (8 \log n /(\eta \log q))$ and let

$$
\widehat{\Lambda}:=\left\{\lambda_{i}=q^{-i}, i \in \mathbb{N}, \text { s.t. } \lambda_{i} \geq 100\left(\widehat{\mathcal{N}}(\lambda) \vee L^{2} / n\right)\right\} .
$$

- Define the parameter choice

$$
\begin{aligned}
& \widehat{\lambda}=\max \left\{\lambda \in \widehat{\Lambda}: \forall \lambda^{\prime} \in \widehat{\Lambda}, \text { s.t. } \lambda^{\prime} \leq \lambda:\right. \\
& \\
& \left.\qquad\left\|\left(\widehat{\Sigma}+\lambda^{\prime}\right)^{\frac{1}{2}}\left(\widehat{f}_{\lambda}-\widehat{f}_{\lambda^{\prime}}\right)\right\| \leq c L \sqrt{\lambda^{\prime}} \widehat{S}\left(\lambda^{\prime}\right)\right\}
\end{aligned}
$$

where

$$
\widehat{\mathcal{S}}(\lambda):=\frac{\sigma \sqrt{2(\widehat{\mathcal{N}}(\lambda) \vee 1)}+M / 5}{\sqrt{\lambda n}} .
$$

## Result for the empirical selection procedure

## Theorem

Assume the source condition (SC) ( $r, R$ ) holds.
Then for the generalized-Lepski parameter choice $\widehat{\lambda}$, with probability at least $1-\eta$ :

$$
\left\|(\Sigma+\lambda)^{\frac{1}{2}}\left(\widehat{f}_{\widehat{\lambda}}-f_{0}\right)\right\| \lesssim L^{3} \min _{\lambda \in\left[\lambda_{\text {min }}, 1\right]}\left(R \lambda^{r+\frac{1}{2}}+\sigma \sqrt{\frac{\mathcal{N}(\lambda)}{n}}+\frac{1}{n \sqrt{\lambda}}+\mathcal{O}\left(n^{-\frac{1}{2}}\right)\right) .
$$

where

$$
\lambda_{\min }=\min \left\{\lambda \in[0,1]: \lambda \gtrsim\left(\mathcal{N}(\lambda) \vee L^{2} / n\right)\right\} .
$$

Conclusion: as a direct byproduct we get the same rates (up to $\log \log n$ factor) as the optimal choice of $\lambda$ in the original bound, for both norms of interest.

## Can we estimate the noise variance $\sigma^{2}$ ?

- Observe that in general, there is no identifiability in the model

$$
y_{i}=f\left(x_{i}\right)+\sigma \xi_{i},
$$

if the function $f$ can be "arbitrary".

- There is a hope when we assumed that $f$ has some regularity (here: linearity)
- Idea:
- Take $\lambda$ small so that the "bias" $\mathcal{A}(\lambda)$ is expected to be much lower than the "variance" $\mathcal{S}(\lambda)$ (e.g., close to $\widehat{\lambda}_{\text {min }}$.
- Split the sample into two subsamples giving rise to $\widehat{f}_{\lambda}^{(1)}, \widehat{f}_{\lambda}^{(2)}$.
- The hope is that by considering $\left\|\hat{f}_{\lambda}^{(1)}-\widehat{f}_{\lambda}^{(2)}\right\|^{2}$ in a suitable norm, we cancel the bias and observe twice the "variance".
- Need somewhat precise concentration (upper and lower) for this quantity.


## Estimation of the variance $\sigma^{2}$

- Assume we have two independent sample of the same size $n$, giving rise to estimators $\widehat{f}_{\lambda}^{(1)}, \widehat{f}_{\lambda}^{(2)}$ (using the same regularization parameter $\lambda>0$ ).
- Consider the statistic

$$
\begin{aligned}
\Delta^{2} & :=\left\|\frac{1}{2}\left(\widehat{\Sigma}^{(1)}+\widehat{\Sigma}^{(2)}+\lambda\right)^{\frac{1}{2}}\left(\widehat{f}_{\lambda}^{(1)}-\widehat{f}_{\lambda}^{(2)}\right)\right\|_{\mathcal{H}}^{2} \\
& \left.=\frac{1}{2 n} \sum_{i=1}^{2 n}\left(\widehat{f}_{\lambda}^{(1)}-\widehat{f}_{\lambda}^{(2)}\right)^{2}\left(x_{i}\right)+\lambda \| \widehat{f}_{\lambda}^{(1)}-\widehat{f}_{\lambda}^{(2)}\right) \|_{\mathcal{H}^{\prime}}^{2}
\end{aligned}
$$

and (rescaling so that the estimate is approximately unbiased)

$$
\widehat{\sigma}^{2}:=\frac{\Delta^{2}}{\sum_{i, j=1}^{2}\left\|A_{i j}\right\|_{H S}^{2}}
$$

where $A_{i j}=\left(\widehat{\Sigma}^{(i)}+\lambda\right)^{\frac{1}{2}} \zeta_{\lambda}\left(\widehat{\Sigma}^{(j)}\right)\left(\widehat{\Sigma}^{(j)}\right)^{\frac{1}{2}}$.

## Estimation of the variance $\sigma^{2}$

## Theorem

If $\lambda \geq \hat{\lambda}_{\text {min }}$, where

$$
\hat{\lambda}_{\text {min }}=\min \left\{\lambda>0: \lambda \geq 100\left(\widehat{\mathcal{N}}(\lambda) \vee \log \left(\eta^{-1}\right) / 2\right)\right\}
$$

then with probability at least $1-\eta$, it holds

$$
\widehat{\sigma}^{2} \in\left[\sigma^{2} \pm\left(\lambda \sigma^{2}+F(\lambda) \log \left(\eta^{-1}\right)\right)\right]
$$

with $F(\lambda) \rightarrow 0$ as $\lambda \rightarrow 0$.
(Proof: based on previously established tools in this setting+Hanson-Wright inequality)
Conclusion: the estimator $\widehat{\sigma}^{2}$ is consistent, and can be used as a proxy for $\sigma^{2}$ in the procedure, with the same conclusions (up to changes in numerical constants, and for $n$ big enough).

## THANK YOU FOR YOUR ATTENTION

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