Non-Gaussian Noise

Review: Probabilistic Classification



Last time, we talked about *probabilistic classification*, i.e. regression with *classification noise*.

$$\hat{\mu} = \operatorname*{argmin}_{m \in \mathcal{M}} \frac{1}{n} \sum_{i=1}^n \big\{ Y_i - m(X_i) \big\}^2 \quad \text{ where } \quad Y_i = \mu(X_i) + \varepsilon_i \text{ for } \varepsilon_i = \begin{cases} 1 - \mu(X_i) & \text{ w.p. } \mu(X_i) \\ -\mu(X_i) & \text{ w.p. } 1 - \mu(X_i) \end{cases}$$

By comparing widths, we showed that this is easier than regression with ...

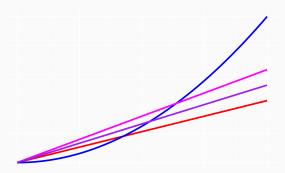
- 1. random sign noise, $s_i = \pm 1$ each w.p. 1/2.
- 2. gaussian noise σg_i of standard deviation $\sigma = 1.25$.

Easier in the sense that our crossing-point argument gives us a better error bound.

$$\frac{s^2}{2} \ge 1.25 \, \mathrm{w}(\mathcal{M}_s) \stackrel{\Longrightarrow}{\ge} \mathrm{w}_s(\mathcal{M}_s) \stackrel{\Longrightarrow}{\ge} \mathrm{w}_{\varepsilon}(\mathcal{M}_s)$$



Starting Point

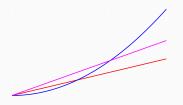


Today, we're going to generalize that result to regression with *any kind of noise*. We'll start with the same abstract bound. It applies no matter how noise is distributed.

$$\begin{split} &\|\hat{\mu} - \mu^\star\|_{L_2(\mathbf{P_n})} < s + 2\sqrt{\frac{2\Sigma_n}{\delta n}} \quad \text{w.p. } 1 - \delta \text{ for } \quad \frac{s^2}{2} \geq \mathbf{w_e}(\mathcal{M}_s) \\ &\text{where} \quad \mathbf{w}_\epsilon(\mathcal{V}) = \mathbf{E}\max_{\mathbf{r} \in \mathcal{N}} \langle \epsilon, v \rangle_{L_2(\mathbf{P_n})} \quad \text{and} \quad \Sigma_n = \mathbf{E}\max_{\mathbf{r} \in \mathcal{N}} \varepsilon_i^2. \end{split}$$

This bound depends on the model $\mathcal M$ and the distribution of the noise ε in a complex, entangled way: through the width $w_\varepsilon(\mathcal M_s)$.

Our Approach



To disentangle the impact of the model and noise distribution, we'll bound this width in terms of gaussian width.

$$w_{\epsilon}(\mathcal{M}_s) \leq \alpha w(\mathcal{M}_s)$$

for α depending on ε but not $\mathcal M$ or s.

At the heart of this comparison $\mathbf{w}_{\epsilon}(\cdot) \leq \alpha \mathbf{w}(\cdot)$ are two ideas.

1. Symmetrization. We'll substitute for ϵ_i a variant that's symmetric around zero.

$$\epsilon_i
ightarrow \epsilon_i - \epsilon_i'$$
 where ϵ_i' is an independent copy of ϵ_i

This substitution increases width: $w_{\epsilon}(\cdot) \leq w_{\epsilon-\epsilon'}(\cdot)$.

2. Contraction. We'll substitute a gaussian vector¹ for our symmetrized noise $\epsilon - \epsilon'$. We can bound the impact of this substitution in a model-invariant way.

$$\mathbf{w}_{\epsilon-\epsilon'}(\cdot) \leq 2M_n \, \mathbf{w}_s(\cdot) \leq \sqrt{2\pi} M_n \times \mathbf{w}(\cdot) \quad \text{for} \quad M_n = \mathbf{E} \max_{i \in 1...n} |\varepsilon_i|$$

This lets us re-use our gaussian width calculations to analyze regression with any noise distribution.

¹or a random-sign vector

An Example

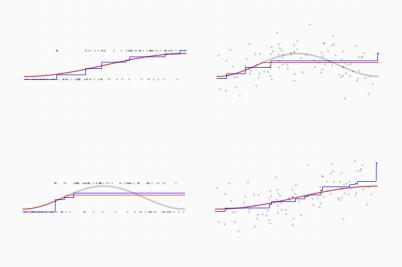


Figure 1: real noise \rightarrow symmetrized noise \downarrow scaled sign noise \leftarrow scaled gaussian noise

Here's the same signal with 4 types of noise.

Symmetrization

$$w_{\varepsilon}(\mathcal{V}) \leq w_{s(\varepsilon - \varepsilon')}(\mathcal{V}) \leq 2 w_{s\varepsilon}(\mathcal{V})$$

$$E \max_{v \in \mathcal{V}} \sum_{i=1}^{n} \varepsilon_{i} v_{i} = E \max_{v \in \mathcal{V}} \sum_{i=1}^{n} (\varepsilon_{i} - E \varepsilon'_{i}) v_{i}$$

$$\stackrel{(a)}{\leq} \operatorname{EE'} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} (\varepsilon_i - \varepsilon_i') v_i$$

$$= E_s E E' \max_{v \in \mathcal{V}} \sum_{i=1}^n s_i (\varepsilon_i - \varepsilon_i') v_i$$

$$\overset{(b)}{\leq} \operatorname{E}_s \operatorname{E} \max_{v \in \mathcal{V}} \sum_{i=1}^n s_i \varepsilon_i + \operatorname{E}_s \operatorname{E}' \max_{v \in \mathcal{V}} \sum_{i=1}^n s_i \varepsilon_i' v_i = 2 \operatorname{E}_s \operatorname{E} \max_{v \in \mathcal{V}} \sum_{i=1}^n \varepsilon_i s_i v_i.$$





Symmetrization

$$w_{\varepsilon}(\mathcal{V}) \leq w_{s(\varepsilon - \varepsilon')}(\mathcal{V}) \leq 2 w_{s\varepsilon}(\mathcal{V})$$

$$\begin{split} \mathbf{E} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} \varepsilon_{i} v_{i} &= \mathbf{E} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} (\varepsilon_{i} - \mathbf{E} \, \varepsilon_{i}') v_{i} \\ &\overset{(a)}{\leq} \mathbf{E} \, \mathbf{E}' \max_{v \in \mathcal{V}} \sum_{i=1}^{n} (\varepsilon_{i} - \varepsilon_{i}') v_{i} \\ &= \mathbf{E}_{s} \, \mathbf{E} \, \mathbf{E}' \max_{v \in \mathcal{V}} \sum_{i=1}^{n} s_{i} (\varepsilon_{i} - \varepsilon_{i}') v_{i} \\ &\overset{(b)}{\leq} \mathbf{E}_{s} \, \mathbf{E} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} s_{i} \varepsilon_{i} + \mathbf{E}_{s} \, \mathbf{E}' \max_{v \in \mathcal{V}} \sum_{i=1}^{n} s_{i} \varepsilon_{i}' v_{i} = 2 \, \mathbf{E}_{s} \, \mathbf{E} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} \varepsilon_{i} s_{i} v_{i}. \end{split}$$

- (a) Replacing ε_i with $s_i(\varepsilon_i \varepsilon_i')$ is 'free'.
 - · We stopped here in our classification example because $\varepsilon_i \varepsilon_i'$ was easy to bound.
 - Generally, we take an extra step to express things in terms of ε_i again.
- (b) Replacing ε_i with $s_i\varepsilon_i$ increases width by at most $2\times$.

Contraction

$$\begin{split} \mathbf{w}_{\eta}(\mathcal{V}) &= \mathbf{w}_{s\eta}(\mathcal{V}) \leq \mathbf{E} \|\eta\|_{\infty} \, \mathbf{w}_{\eta}(\mathcal{V}) \quad \text{if} \quad \eta \overset{dist}{=} -\eta. \\ \mathbf{E}_{s} \, \mathbf{E}_{\eta} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} \eta_{i} s_{i} v_{i} \leq \mathbf{E}_{\eta} \max_{u \in \mathbb{R}^{n} \|u_{i}| \leq \|\eta\|_{\infty}} \mathbf{E}_{s} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} u_{i} s_{i} v_{i} \\ &= \mathbf{E}_{\eta} \|\eta\|_{\infty} \max_{u \in [-1,1]^{n}} \mathbf{E}_{s} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} u_{i} s_{i} v_{i} \\ &= \mathbf{E}_{\eta} \|\eta\|_{\infty} \times \max_{u \in \{-1,1\}^{n}} \mathbf{E}_{s} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} u_{i} s_{i} v_{i} \\ &= \mathbf{E}_{\eta} \|\eta\|_{\infty} \times \mathbf{E}_{s} \max_{v \in \mathcal{V}} \sum_{i=1}^{n} s_{i} v_{i} \end{split}$$



$$\begin{split} \mathbf{w}_{\eta}(\mathcal{V}) &= \mathbf{w}_{s\eta}(\mathcal{V}) \leq \mathbf{E} \|\eta\|_{\infty} \, \mathbf{w}_{\eta}(\mathcal{V}) \quad \text{if} \quad \eta \overset{dist}{=} -\eta. \\ \mathbf{E}_{s} \, \mathbf{E}_{\eta} \, \max_{v \in \mathcal{V}} \sum_{i=1}^{n} \eta_{i} s_{i} v_{i} \leq \mathbf{E}_{\eta} \, \max_{u \in \mathbb{R}^{n} \mid u_{i} \leq \|\eta\|_{\infty}} \mathbf{E}_{s} \, \max_{v \in \mathcal{V}} \sum_{i=1}^{n} u_{i} s_{i} v_{i} \\ &= \mathbf{E}_{\eta} \|\eta\|_{\infty} \, \max_{u \in \{-1,1\}^{n}} \mathbf{E}_{s} \, \max_{v \in \mathcal{V}} \sum_{i=1}^{n} u_{i} s_{i} v_{i} \\ &= \mathbf{E}_{\eta} \|\eta\|_{\infty} \times \max_{u \in \{-1,1\}^{n}} \mathbf{E}_{s} \, \max_{v \in \mathcal{V}} \sum_{i=1}^{n} u_{i} s_{i} v_{i} \\ &= \mathbf{E}_{\eta} \|\eta\|_{\infty} \times \mathbf{E}_{s} \, \max_{v \in \mathcal{V}} \sum_{i=1}^{n} s_{i} v_{i} \end{split}$$

- We can 'contract out' any symmetrically distributed noise vector η by ...
 - 1. multiplying in independent random signs s_i . Symmetry $\implies s_i \eta_i \stackrel{dist}{=} \eta_i$.
 - 2. maximizing over a cube containing η .
- · We just have to use a big enough cube.
 - · In our classification example, $\eta=\varepsilon-\varepsilon'$ was in the unit cube $[-1,1]^n$ deterministically.
 - Generally, we maximize over a random cube $[-\|\eta\|_{\infty}, \|\eta\|_{\infty}]^n$.
 - · And we can pull out the cube's radius $\|\eta\|_{\infty}$ as a multiplicative factor.

Symmetrization, Contraction, and Gaussian Noise

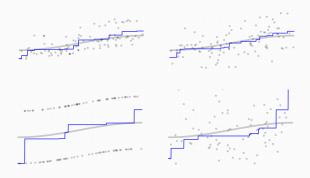


Figure 2: real noise \rightarrow symmetrized noise \downarrow scaled sign noise \leftarrow scaled gaussian noise

After symmetrizing and introducing random signs, i.e. making the substitution

$$\varepsilon_i \to s_i(\varepsilon_i - \varepsilon_i'),$$

we 'contract out' the symmetrized noise $\varepsilon - \varepsilon'$ to get a bound in terms of random-sign width.

$$w_{\varepsilon}(\mathcal{V}) \leq w_{s(\varepsilon - \varepsilon')}(\mathcal{V}) \leq \|\varepsilon - \varepsilon'\|_{\infty} w_{s}(\mathcal{V}) \leq \|\varepsilon - \varepsilon'\|_{\infty} 1.25 w(\mathcal{V})$$

We can substitute 1.25 times gaussian width because that's at least as large as random sign width.

$$\mathrm{E}\max_{v\in\mathcal{V}}\sum_{i=1}^n g_iv_i = \mathrm{E}_s\,\mathrm{E}_g\max_{v\in\mathcal{V}}\sum_{i=1}^n |g_i|\; s_iv_i \geq \mathrm{E}_s\max_{v\in\mathcal{V}}\sum_{i=1}^n \mathrm{E}_g|g_i|\; s_iv_i.$$

Implications for Regression



$$\begin{split} \mathbf{w}_{\varepsilon}(\mathcal{V}) &\leq M \, \mathbf{w}_{s}(\mathcal{V}) \leq 1.25 M \, \mathbf{w}(\mathcal{V}) \\ \text{for } & \mathsf{M} = \mathbf{E} \| \varepsilon - \varepsilon' \|_{\infty} \leq 2 \, \mathbf{E} \| \varepsilon \|_{\infty}. \end{split}$$

In terms of our crossing-point bounds, regression with arbitrary independent noise,

i.e.
$$Y_i = \mu(X_i) + \varepsilon_i$$
 where $\varepsilon_1 \dots \varepsilon_n$ are independent,

is no harder than with scaled random sign noise or with gaussian noise

i.e.
$$Y_i=\mu(X_i)+Ms_i$$
 for $s_i=egin{cases} +1 & \text{w.p. } 1/2 \\ -1 & \text{w.p. } 1/2 \end{cases}$ or $=\mu(X_i)+1.25Mg_i$ for $g_i\sim N(0,1)$







The scale factor is $2 \times$ the expected magnitude of our noise vector's largest element.

Lost Precision



Figure 3: standard gaussian noise →scaled random sign noise →scaled gaussian noise

- · This isn't the absolute best bound we can get.
- · For example, if we start with standard gaussian noise, we lose ...
- · ...a factor of roughly $7\sqrt{\log(2n)}$ going to random sign width and back.

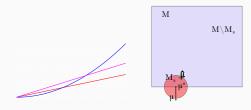
$$\mathrm{w}_{\varepsilon}(\mathcal{V}) \leq 2\,\mathrm{w}_{s\varepsilon}(\mathcal{V}) \leq 2 \times 2 \sqrt{2\log(2n)}\,\mathrm{w}_{s}(\mathcal{V}) \leq 4 \sqrt{2\log(n)} \times \sqrt{\frac{\pi}{2}}\,\mathrm{w}_{\varepsilon}(\mathcal{V}) \approx 7 \sqrt{\log(2n)}\,\mathrm{w}_{\varepsilon}(\mathcal{V}).$$

- (a) 'Symmetrization' cost us a factor of 2.
- (b) Contraction costs us a factor of $\mathrm{E}\max_{i < n} |\varepsilon_i| \le 2\sqrt{2\log(2n)}$. (See HW Appendix B)
- (c) Converting random signs back to gaussians costs us a factor of $\sqrt{\frac{\pi}{2}} \approx 1.25$.

We're in the right ballpark. For sample sizes n between 50 and 50 million, that factor is between 15 and 30. But if we want a more precise error bound, we need to be a little more careful.

Sampling

What We've Done



We have a bound that's valid for any signal μ and any vector of independent noise ε .

$$\|\hat{\mu} - \mu^\star\|_{L_2(\mathbf{P_n})} < 2\sqrt{\Sigma_n} \left(s + \sqrt{\frac{2}{\delta n}}\right) \quad \text{w.p. } 1 - \delta \text{ for } \quad \frac{s^2}{2} \geq \mathbf{w}_s(\mathcal{M}_s)$$

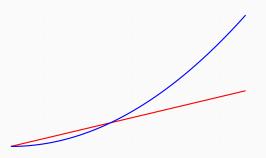
- It depends on the model's size through the critical radius of random-sign width.
 - $s \quad \text{satisfying} \quad s^2/2 \geq \underline{\mathbf{w}}_s(\mathcal{M}_s) \quad \text{for} \quad \mathcal{M}_s = \{m \in \mathcal{M} \, : \, \|m \mu^\star\|_{L_2(\mathbf{P_n})} \leq s\}$
 - · This is a one-number summary of the random-sign width of neighborhoods ...
 - \cdot ...of the model's best approximation to the signal. It's the summary that matters.
- It depends on the noise's size through the expected maximum square.

$$\Sigma_n = \mathrm{E} \max_{i \in 1 \dots n} |\varepsilon_i|^2$$

What does this tell us?

Bounds like this say how close $\hat{\mu}$ and μ^\star are, on average, on our sample $X_1 \dots X_n$.

$$\frac{1}{n}\sum_{i=1}^{n} \{\hat{\mu}(X_i) - \mu^{\star}(X_i)\} < \dots$$



It doesn't tell us how close they are in the gaps between those points.

- · Let's think about what happens when $X_1 \dots X_n$ is are drawn independently from some distribution P. Think sampling with replacement from a population.
- · We'll bound the population root mean squared error $\|\hat{\mu} \mu^{\star}\|_{L_2(\mathbf{P})}$.

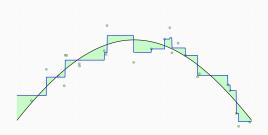
What Population Mean Squared Error Is

It's the mean squared error we make at random point X' distributed like $X_1 \ldots X_n$.

$$\|\hat{\mu} - \mu^*\|_{L_2(\mathbf{P})}^2 = \mathcal{E}_{X'} \left[\{\hat{\mu}(X') - \mu^*(X')\}^2 \right]$$

That's the integral of the squared distance between the two curves, multiplied by the density of X_i .

$$\|\hat{\mu} - \mu^{\star}\|_{L_2(\mathbf{P})}^2 = \int {\{\hat{\mu}(x) - \mu^{\star}(x)\}^2 p(x) dx}$$
 if X_i has the density $p(x)$.

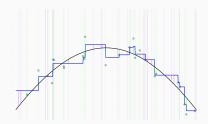


Why we care about Population Mean Squared Error: Generalization

If we're interested in average accuracy for a bunch of new points $X_1' \dots X_{n'}'$ distributed like $X_1 \dots X_n$, that's more or less exactly what it is.

$$\|\hat{\mu} - \mu\|_{L_2(\mathbf{P})}^2 = \mathbf{E}_{X'} \left[\left\{ \hat{\mu}(X') - \mu(X') \right\}^2 \right] \stackrel{LLN}{\approx} \frac{1}{n'} \sum_{i=1}^{n'} \left\{ \hat{\mu}(X_i') - \mu(X_i') \right\}^2.$$

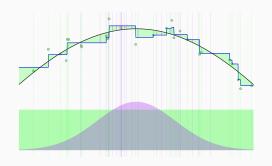
This can be a bit different from accuracy on our original sample $X_1 \dots X_n$.



- BV regression spends its 'variation budget' jumping to fit on the original sample.
- Between those points, it doesn't know whether it should jump or not.
 - · So we can get larger error at our new points.
 - · It's usually not much larger, but sometimes it is. We'll see why.

Why we care about Population Mean Squared Error: Generalization

If we're interested in average accuracy for new points from a different distribution Q, we can bound this by comparing this distribution's density to that of our observations.

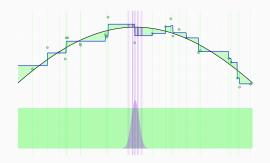


$$\frac{1}{n'} \sum_{i=1}^{n'} \{ \hat{\mu}(X_i') - \mu(X_i') \}^2 \approx \|\hat{\mu} - \mu\|_{L_2(\mathbb{Q})}^2 = \int \{ \hat{\mu}(x) - \mu(x) \}^2 \frac{q(x)}{p(x)} p(x) dx$$

$$\leq \max_x \frac{q(x)}{p(x)} \|\hat{\mu} - \mu\|_{L_2(\mathbb{P})}^2.$$

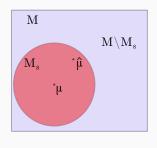
Why we care about Population Mean Squared Error: Generalization

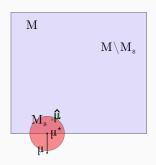
If we're interested in accuracy at a specific point x', we can think of this new distribution Q as a little bump around x'.



$$\{\hat{\mu}(x') - \mu(x')\}^2 \approx \|\hat{\mu} - \mu\|_{L_2(Q_\epsilon)} \quad \text{ for } \quad Q = N(x', \epsilon^2).$$

Same Argument, Different Neighborhood





- We want to show that $\hat{\mu}$ is in a population-distance neighborhood of μ .
- $\boldsymbol{\cdot}\,$ Or, if we've chosen the model wrong, at least its best population-distance approximation.

$$\hat{\mu} \in \mathcal{M}_s \quad \text{for} \quad \mathcal{M}_s = \{m \in \mathcal{M} \,:\, \|m - \mu^\star\|_{L_2(\mathbf{P})} \leq s\} \quad \text{for} \quad \mu^\star = \underset{m \in \mathcal{M}}{\operatorname{argmin}} \|m - \mu\|_{L_2(\mathbf{P})}$$

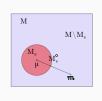
- · We'll do this using essentially the same argument we used to bound sample MSE.
 - 1. We know that the $\hat{\mu}$'s squared error loss is at least as good as μ^{\star} 's.
 - 2. We find a radius s for which every curve with this property is in the neighborhood \mathcal{M}_s .
- · It amounts to showing the loss difference $\ell(m)-\ell(\mu^\star)$ is positive outside this neighborhood.

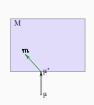
$$m \in \mathcal{M}_s$$
 if $m \in \mathcal{M}_s$ and $\ell(m) - \ell(\mu^*) > 0$ for all $m \in \mathcal{M} \setminus \mathcal{M}_s$

Reduction to a Maximal Inequality

$$\ell(m) - \ell(\mu^*) = \frac{1}{n} \sum_{i=1}^n Z_i(m) := \{ m(X_i) - \mu^*(X_i) \}^2 - 2 \{ Y_i - \mu^*(X_i) \} \{ m(X_i) - \mu^*(X_i) \}$$
$$= \mathbb{E} Z_i(m) + \frac{1}{n} \sum_{i=1}^n Z_i(m) - \mathbb{E} Z_i(m).$$

Convexity Helps as Usual.





1. The loss difference is positive outside the neighborhood if it's positive on its boundary.

$$m \in \mathcal{M}_s$$
 if $m \in \mathcal{M}_s$ and $\ell(m) - \ell(\mu^\star) > 0$ for all $m \in \mathcal{M} \setminus \mathcal{M}_s \mathcal{M}_s^\circ$

2. The projection theorem tells us an unwanted term in $\mathbf{E} Z_i(m)$ is non-negative.

$$\begin{split} -\operatorname{E}\left[\left\{Y_{i}-\mu^{\star}(X_{i})\right\}\left\{m(X_{i})-\mu^{\star}(X_{i})\right\}\right] &= -\operatorname{E}\left[\left\{\operatorname{E}[Y_{i}\mid X_{i}]-\mu^{\star}(X_{i})\right\}\left\{m(X_{i})-\mu^{\star}(X_{i})\right\}\right] \\ &= \left\langle\mu^{\star}-\mu, m-\mu^{\star}\right\rangle_{L_{2}(P)} \geq 0 \quad \text{for all} \quad m \in \mathcal{M} \end{split}$$

It follows that
$$m \in \mathcal{M}_s$$
 if $m \in \mathcal{M}_s$ and $s^2 > \max_{m \in \mathcal{M}_0^s} \frac{1}{n} \sum_{i=1}^n Z_i(m) - \operatorname{E} Z_i(m)$

Bounding the New Maximum

We show this maximum is approximately constant, i.e. close to its expectation.

$$\bar{Z} := \max_{m \in \mathcal{M}_{\mathcal{S}}^{\circ}} \frac{1}{n} \sum_{i=1}^{n} Z_i(m) - \operatorname{E} Z_i(m) \quad \text{ satisfies } \quad \bar{Z} \leq \operatorname{E} \bar{Z} + \sqrt{\frac{\operatorname{Var}(\bar{Z})}{\delta n}} \quad \text{ w.p. } 1 - \delta$$

We use symmetrization to bound its expectation in terms of random-sign width.

- (a) Write the centers $\mathrm{E}\,Z_i(v)$ in terms of an independent copy of our sample.
- (b) Compare the result to a maximum of an average of symmetric random variables.
- (c) Introduce random signs and compare to two copies of a simpler maximum.

$$\begin{split} n \times & \to \bar{Z} \overset{(a)}{=} \to_{Z} \max_{m \in \mathcal{M}_{S}^{\circ}} \to_{Z'} \sum_{i=1}^{n} \left\{ Z_{i}(m) - Z'_{i}(m) \right\} \\ & \stackrel{(b)}{\leq} \to_{Z} \to_{Z'} \to_{s} \max_{m \in \mathcal{M}_{S}^{\circ}} \sum_{i=1}^{n} s_{i} \left\{ Z_{i}(m) - Z'_{i}(m) \right\} \\ & \stackrel{(c)}{\leq} \to_{Z} \to_{Z'} \to_{s} \max_{m, m' \in \mathcal{M}_{S}^{\circ}} \sum_{i=1}^{n} s_{i} Z_{i}(m) + (-s_{i}) Z_{i}(m') \\ & = 2 \to_{Z} \to_{s} \max_{m \in \mathcal{M}_{S}^{\circ}} \sum_{i=1}^{n} s_{i} Z_{i}(m) \end{split}$$

We can use the Efron-Stein inequality to bound the variance. Come back and try it later!

$$\operatorname{Var}(\bar{Z}) \overset{\text{why?}}{\leq} \frac{1}{n^2} \sum_{i=1}^n \operatorname{E}\left\{Z_i(\hat{m}) - Z_i'(\hat{m})\right\}_+^2 \quad \text{for} \quad \hat{m} = \operatorname*{argmax}_{m \in \mathcal{M}_\delta^0} \sum_{i=1}^n Z_i(m) - \operatorname{E}Z_i(m)$$

Contracting Out Lipschitz Functions

What we get is $2\times$ the expected random-sign width of some set of vectors, but it's not just the set of the vectors in our neighborhood $\mathcal{M}_s-\mu^\star$.

$$n \times \mathbf{E} Z \le 2 \mathbf{E} \mathbf{E}_{s} \max_{m \in \mathcal{M}_{s}^{o}} \sum_{i=1}^{n} s_{i} Z_{i}(m) = \left\{ m(X_{i}) - \mu^{*}(X_{i}) \right\}^{2} - 2 \left\{ Y_{i} - \mu^{*}(X_{i}) \right\} \left\{ m(X_{i}) - \mu^{*}(X_{i}) \right\}$$

$$\leq 4 \operatorname{E} \left\{ \max_{m \in \mathcal{M}_{s}^{\circ}} \|m - \mu\|_{L_{\infty}(\operatorname{Pn})} + \|\varepsilon\|_{L_{\infty}(\operatorname{Pn})} \right\} \operatorname{E}_{s} \max_{m \in \mathcal{M}_{s}^{\circ}} \sum_{i=1}^{n} s_{i} \{m(X_{i}) - \mu^{\star}(X_{i})\}$$

We've compared that to the width of the neighborhood itself using ... Lemma (Lipschitz Comparison)

$$\mathbf{E}_s \max_{v \in \mathcal{V}} \sum_{i=1}^n s_i \psi_i(v_i) \leq L \, \mathbf{E}_s \max_{v \in \mathcal{V}} \sum_{i=1}^n s_i v_i \ \text{if} \ |\psi_i(u_i) - \psi_i(v_i)| \leq L |u_i - v_i| \ \text{for all} \ u, v \in \mathcal{V}.$$

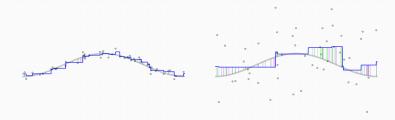
For
$$\psi_i(v) = v_i^2 - 2\{Y_i - \mu^{\star}(X_i)\}v_i$$
 and $V = \{m(X_1) - \mu^{\star}(X_1) \dots m(X_n) - \mu^{\star}(X_n) : m \in \mathcal{M}_s^{\circ}\},$ that's $\mathbf{E}_s \max_{m \in \mathcal{M}_s^{\circ}} \sum_{i=1}^n s_i \psi_i \{m(X_i) - \mu^{\star}(X_i)\} \leq L \max_{m \in \mathcal{M}_s^{\circ}} \sum_{i=1}^n s_i \{m(X_i) - \mu^{\star}(X_i)\}$

where
$$L = \max_{i} \max_{m \in \mathcal{M}_{s}^{0}} |\psi_{i}'\{m(X_{i}) - \mu^{\star}(X_{i})\}|$$

$$= \max_{i} \max_{m \in \mathcal{M}_{s}^{0}} |2\{m(X_{i}) - \mu^{\star}(X_{i})\} - 2\{Y_{i} - \mu^{\star}(X_{i})\}|$$

$$\leq 2 \max_{m \in \mathcal{M}_{0}} \|m - \mu\|_{L_{\infty}(\mathbf{P_n})} + 2\|\varepsilon\|_{L_{\infty}(\mathbf{P_n})}.$$

Interpretation



$$\begin{split} \|\hat{\mu} - \mu^\star\|_{L_2(\mathbf{P})} &\leq s \times 2\Big\{\sqrt{\Sigma_n} + B\Big\} + \sqrt{\frac{\mathrm{Var}(\bar{Z})}{\delta}} \quad \text{w.p. } 1 - \delta \\ &\text{if} \quad \frac{s^2}{2} \geq \mathrm{E}\, \mathrm{w}_s(\mathcal{M}_s) \quad \text{and} \quad \|m - \mu\|_\infty \leq B \end{split}$$

This is the bound we'd get on sample MSE with additional scaled random-sign noise,

i.e. if we'd observed
$$Y_i = \mu(X_i) + \varepsilon_i + B s_i$$

Left: With little noise, our estimator $\hat{\mu}$ fits substantially better at the sample points X_i . Right: With more, it doesn't. The observations are far enough from μ that we can't estimate it all that precisely even where we have some data.

Signal Recovery



Signal Recovery is regression without any noise at all. In that case ($\Sigma_n = 0$),

$$\begin{split} \|\hat{\mu} - \mu\|_{L_2(\mathbf{P})} &\leq s \times 2\Big\{\sqrt{\Sigma_n} + B\Big\} + \sqrt{\frac{\mathrm{Var}(\bar{Z})}{\delta}} \quad \text{ w.p. } 1 - \delta \end{split}$$
 if $\frac{s^2}{2} \geq \mathrm{E}\, \mathrm{w}_s(\mathcal{M}_s)$ and $\|m - \mu\|_{\infty} \leq B$

This is the bound we'd get on sample MSE with only scaled random-sign noise.

i.e. if we'd observed
$$Y_i = \mu(X_i) + \varepsilon_i + Bs_i$$

- This is an extreme case of the low-noise regime. And it's still hard.
- · When you want to estimate μ between the sample points $X_1 \dots X_n$, ...
- ...what you want to see obscured by bounded 'sampling noise' $\in [-B, B]$.

References

Chapter 6 of Talagrand's Upper and Lower Bounds for Stochastic Processes.

· Random Signs vs. Gaussians: Proposition 6.22

· Contraction: Lemma 6.4.5

· Lipschitz Contraction: Theorem 6.5.1

Appendices

Appendices

Boundedness



Our Population MSE bound introduces a new consideration: boundedness of $\|m-\mu\|_\infty$ in neighborhoods of μ^\star .

$$\begin{split} \|\hat{\mu} - \mu\|_{L_2(\mathbf{P})} &\leq s \times 2\Big\{\sqrt{\Sigma_n} + B\Big\} + \sqrt{\frac{\mathrm{Var}(\bar{Z})}{\delta}} \quad \text{ w.p. } 1 - \delta \\ &\text{if } \quad \frac{s^2}{2} \geq \mathrm{E}\, \mathrm{w}_s(\mathcal{M}_s) \quad \text{and} \quad \|m - \mu\|_{\infty} \leq B \end{split}$$

Getting a bound B can take a bit of work. There are options.

Option 1. Baking it into the Model.

$$\begin{split} \mathcal{M} &= \{m \ : \ \|m\|_{\infty} \leq B \quad \text{and} \quad \rho_{TV}(m) \leq B \} \\ &\implies \|m - \mu\|_{\infty} \leq \|m\|_{\infty} + \|\mu\|_{\infty} \leq B + \|\mu\|_{\infty} \\ \mathcal{M} &= \{m \ : \ m(0) = 0 \quad \text{and} \quad \rho_{TV}(m) \leq B \} \\ &\implies \end{split}$$

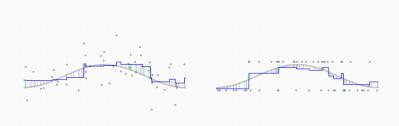


Option 2. Arguing Based on Bounded Data. In many models, you can show that $\hat{\mu}$ is will be within the range of the data.

i.e.
$$\min_{i \leq n} Y_i \leq \hat{\mu}(x) \leq \max_{i \leq n} Y_i$$

This is true, in particular, for Monotone and Bounded Variation Regression. We can add this constraint to our model when doing our analysis.

$$\begin{split} \|\hat{\mu} - \mu^\star\|_{L_2(\mathbf{P})} &< s \quad \text{if} \quad \ell(m) - \ell(\mu^\star) > 0 \quad \text{for all} \quad m \in \mathcal{M} \dots \\ \dots & \text{with} \quad \|m\|_\infty \leq B \quad \text{and} \quad \|m - \mu^\star\|_{L_2(\mathbf{P})} \geq s \end{split}$$



The are other options.