

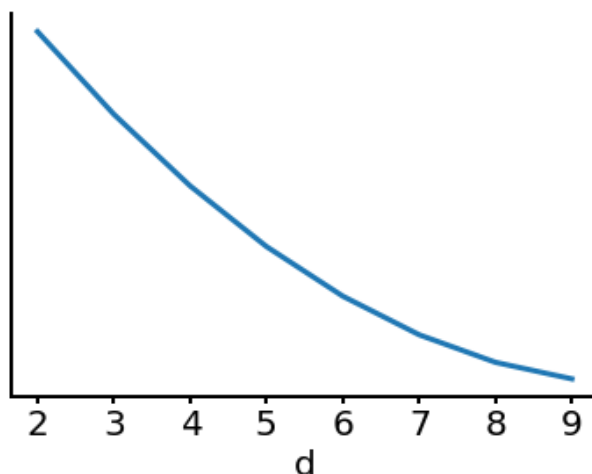
Discussion #10

Name:

Monday March 8th

Bias-Variance Trade-Off

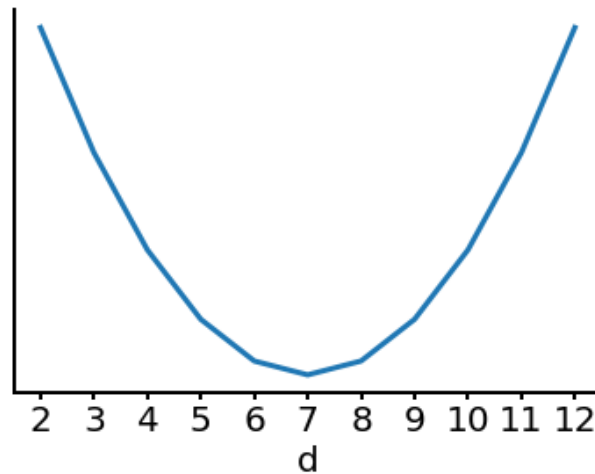
1. Your team would like to train a machine learning model in order to predict the next YouTube video that a user will click on based on the videos the user has watched in the past. We extract m attributes (such as length of video, view count etc) from each video and our model will be based on the previous d videos watched by that user. Hence the number of features for each data point for the model is $m \cdot d$. Currently, you're not sure how many videos to consider.
 - (a) Your colleague generates the following plot, where the value d is on the x-axis. However, they forgot to label the y-axis.



Which of the following could the y-axis represent? Select all that apply.

- ☐ A. Training Error
- ☐ B. Validation Error
- ☐ C. Bias
- ☐ D. Variance

- (b) Your colleague generates the following plot, where the value d is on the x-axis. However, they forgot to label the y-axis again.



Which of the following could the y axis represent? Select all that apply.

- ☐ A. Training Error
- ☐ B. Validation Error
- ☐ C. Bias
- ☐ D. Variance

2. We randomly sample some data $(x_i, y_i)_{i=1}^n$ and use it to fit a model $f_{\hat{\theta}}(x)$ according to some procedure (e.g. OLS, Ridge, LASSO). We then sample a new point that is independent from our existing points, but sampled from the same underlying truth as our data. Furthermore, assume that we have a function $g(x)$ and some noise generation process that produces ϵ such that $\mathbb{E}[\epsilon] = 0$ and $\text{var}(\epsilon) = \sigma^2$. Every time we query mother nature for Y at a given x , she gives us $Y = g(x) + \epsilon$. (The true function for our data is $Y = g(x) + \epsilon$.) A new ϵ is generated each time, independent of the last. In class, we showed that

$$\underbrace{\mathbb{E}[(Y - f_{\hat{\theta}}(x))^2]}_{\text{Total Error}} = \underbrace{\sigma^2}_{\text{Observation Variance}} + \underbrace{(g(x) - \mathbb{E}[f_{\hat{\theta}}(x)])^2}_{\text{Model Bias}^2} + \underbrace{\mathbb{E}[(f_{\hat{\theta}}(x) - \mathbb{E}[f_{\hat{\theta}}(x)])^2]}_{\text{Model Variance}}$$

- (a) Label each of the terms above.

Word Bank: observation variance, model variance, observation bias², model bias², model risk, empirical mean square error.

- (b) What is random in the equation above? Where does the randomness come from?
- (c) True or false and explain. $\mathbb{E}[\epsilon f_{\hat{\theta}}(x)] = 0$
- (d) Suppose you lived in a world where you could collect as many data sets you would like. Given a fixed algorithm to fit a model f_{θ} to your data e.g. linear regression, describe a procedure to get good estimates of $\mathbb{E}[f_{\hat{\theta}}(x)]$
- (e) If you could collect as many data sets as you would like, how does that affect the quality of your model $f_{\theta}(x)$?

Ridge and LASSO Regression

3. Earlier, we posed the linear regression problem as follows: Find the θ value that minimizes the average squared loss. In other words, our goal is to find $\hat{\theta}$ that satisfies the equation below:

$$\hat{\theta} = \underset{\theta}{\operatorname{argmin}} L(\theta) = \underset{\theta}{\operatorname{argmin}} \frac{1}{n} \|\mathbb{Y} - \mathbb{X}\theta\|_2^2$$

Here, \mathbb{X} is a $n \times (p+1)$ matrix, θ is a $(p+1) \times 1$ vector and \mathbb{Y} is a $n \times 1$ vector. As we saw in lecture, the optimal $\hat{\theta}$ is given by the closed form expression $\hat{\theta} = (\mathbb{X}^T \mathbb{X})^{-1} \mathbb{X}^T \mathbb{Y}$.

To prevent overfitting, we saw that we can instead minimize the sum of the average squared loss plus a regularization function $\lambda \mathcal{S}(\theta)$.

- If use the function $\mathcal{S}(\theta) = \|\theta\|_2^2$, we have “ridge regression”.
- If we use the function $\mathcal{S}(\theta) = \|\theta\|_1$, we have “LASSO regression”.

For example, if we choose $\mathcal{S}(\theta) = \|\theta\|_2^2$, our goal is to find $\hat{\theta}$ that satisfies the equation below:

$$\begin{aligned} \hat{\theta} &= \underset{\theta}{\operatorname{argmin}} L(\theta) = \underset{\theta}{\operatorname{argmin}} \frac{1}{n} \|\mathbb{Y} - \mathbb{X}\theta\|_2^2 + \lambda \|\theta\|_2^2 \\ &= \underset{\theta}{\operatorname{argmin}} \frac{1}{n} \sum_{i=1}^n (y_i - \mathbb{X}_{i,\cdot}^T \theta)^2 + \lambda \sum_{j=0}^p \theta_j^2 \end{aligned}$$

Recall that λ is a hyperparameter that determines the impact of the regularization term. Though we did not discuss this in lecture, we can also find a closed form solution to ridge regression: $\hat{\theta} = (\mathbb{X}^T \mathbb{X} + n\lambda \mathbf{I})^{-1} \mathbb{X}^T \mathbb{Y}$. It turns out that $\mathbb{X}^T \mathbb{X} + n\lambda \mathbf{I}$ is guaranteed to be invertible (unlike $\mathbb{X}^T \mathbb{X}$ which might not be invertible).

- (a) As model complexity increases, what happens to the bias and variance of the model?
- (b) In terms of bias and variance, how does a regularized model compare to ordinary least squares regression?
- (c) In ridge regression, what happens if we set $\lambda = 0$? What happens as λ approaches ∞ ?
- (d) How does model complexity compare between ridge regression and ordinary least squares regression? How does this change for large and small values of λ ?
- (e) If we have a large number of features (10,000+) and we suspect that only a handful of features are useful, which type of regression (Lasso vs Ridge) would be more helpful in interpreting useful features?
- (f) What are the benefits of using ridge regression?

Cross Validation

4. After running 5-fold cross validation, we get the following mean squared errors for each fold and value of λ :

Fold Num	$\lambda = 0.1$	$\lambda = 0.2$	$\lambda = 0.3$	$\lambda = 0.4$	Row Avg
1	80.2	70.2	91.2	91.8	83.4
2	76.8	66.8	88.8	98.8	82.8
3	81.5	71.5	86.5	88.5	82.0
4	79.4	68.4	92.3	92.4	83.1
5	77.3	67.3	93.4	94.3	83.0
Col Avg	79.0	68.8	90.4	93.2	

How do we use the information above to choose our model? Do we pick a specific fold? a specific lambda? or a specific fold-lambda pair? Explain.

5. You build a model with two regularization hyperparameters λ and γ . You have 4 good candidate values for λ and 3 possible values for γ , and you are wondering which λ, γ pair will be the best choice. If you were to perform five-fold cross-validation, how many validation errors would you need to calculate?
6. In the typical setup of k-fold cross validation, we use a different parameter value on each fold, compute the mean squared error of each fold and choose the parameter whose fold has the lowest loss.
- ☐ A. True
- ☐ B. False