Homework 3: Model Order Selection

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- (a) (i) No.
 - (ii) There is no underfitting because the power of the model is higher than the true function.
 - (iii) $\vec{w} = [1, 2, 0]$ (bold and arrow used interchangeably)
- (b) (i) No.
 - (ii) There is underfitting because the power of the model is lower than the true function.
- (c) (i) Yes. The true function has x_1, x_2 , but, according to the footnote, our example only has 1 feature.
 - (ii) There is underfitting because the true function has the term x_1x_2 , which the model lacks.

(a)
$$x_0 = 1, x_1 = \text{cancer volume}, x_2 = \text{patient's age}, x_3 = \text{type of cancer}$$

 x_0 let's the x, w subscripts line up more, which looks better.

(Model 1)
$$h(\mathbf{x}) = w_0 + w_1 x_1$$

(Model 2)
$$h(\mathbf{x}) = w_0 + w_1 x_1 + w_2 x_2$$

(Model 3)
$$h(\mathbf{x}) = w_0 + (w_1 + x_3 w_3) x_1 + w_2 x_2$$

(b) There are 2 and 3 parameters, respectively. Model 2 is the most complex out of the 2, however Model 3 is the <u>most</u> complex.

(c) (Model 1)
$$X = \begin{bmatrix} 1 & 0.7 \\ 1 & 1.3 \\ 1 & 1.6 \end{bmatrix}$$

(Model 2)
$$X = \begin{bmatrix} 1 & 0.7 & 55 \\ 1 & 1.3 & 65 \\ 1 & 1.6 & 70 \end{bmatrix}$$

(Model 3)
$$X = \begin{bmatrix} 1 & 0.7 & 55 & 1 \\ 1 & 1.3 & 65 & 0 \\ 1 & 1.6 & 70 & 0 \end{bmatrix}$$

(d) Model 2 should be selected. It has similar numbers to Model 3, but Model 2 is a much simpler model. We want to avoid having an overly complicated model that doesn't even give us better numbers.

A is probably underfitting; this is because its error is much higher than the other models and because its error drops to around its final value at smaller sample sizes, implying a simpler model.

B is probably neither. If a model is overfitted, then its error would be extremely low for small N, and then much higher for larger N.

C is probably overfitted. It has very large fluctuating error, and we see that it is be most complex model because it has the highest N for which it still has 0 training error.

Flexible is good for its higher chance to find the true model, while less flexible is better for runtime. If you don't know (or have a good guess from the graphing the data or otherwise know something about the data's background), it'll probably a good idea to sacrifice some time to get a better model. On the other hand, if you somehow already know the true model, then you could use a less flexible approach since you wouldn't need to "consider" other models and instead just train a single model.

 $\bullet\,$ Use the Hoeffding Inequality. N=100

$$2e^{-2\cdot0.1^2\cdot100} \approx 0.270$$

• N = 200

$$2e^{-2\cdot 0.1^2\cdot 200}\approx 0.036$$

$$X = \begin{bmatrix} 1 & 6.6 & 1 & 4.0 \\ 1 & 6.4 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \\ 1 & 6.4 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \end{bmatrix}$$

$$w_{ridge} = \left(X^T X + N \lambda \mathbf{I}'\right)^{-1} X^T y$$

$$= \left(\begin{bmatrix} 1 & 6.6 & 1 & 4.0 \\ 1 & 6.4 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \end{bmatrix}^{T} \begin{bmatrix} 1 & 6.6 & 1 & 4.0 \\ 1 & 6.4 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \end{bmatrix}^{T} + 5 \cdot 0.1 \cdot \begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix} \right)^{-1} \begin{bmatrix} 1 & 6.6 & 1 & 4.0 \\ 1 & 6.4 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \end{bmatrix}^{T}$$

$$= \left(\begin{bmatrix}1 & 1 & 1 & 1 & 1 \\ 6.6 & 6.4 & 7.2 & 6.4 & 7.2 \\ 1 & 2 & 2 & 2 & 2 \\ 4.0 & 5.0 & 5.0 & 5.0 & 5.0\end{bmatrix}\begin{bmatrix}1 & 6.6 & 1 & 4.0 \\ 1 & 6.4 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0 \\ 1 & 7.2 & 2 & 5.0\end{bmatrix} + 5 \cdot 0.1 \cdot \begin{bmatrix}0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1\end{bmatrix}\right)^{-1} \begin{bmatrix}1 & 1 & 1 & 1 & 1 \\ 6.6 & 6.4 & 7.2 & 6.4 & 7.2 \\ 1 & 2 & 2 & 2 & 2 \\ 4.0 & 5.0 & 5.0 & 5.0 & 5.0\end{bmatrix}$$

We need to account for the L_2 term in the gradient.

$$\mathbf{w} = \mathbf{w} - \frac{\alpha}{N} \left(X^T \left(X \mathbf{w} - \mathbf{y} \right) + \lambda \mathbf{I}' \mathbf{w} \right)$$