4741 Bias and Variance HW 4, Question 4 Aman jain (aj644@cornell.edu)

a)

```
In [117]: using Plots using Statistics
```

Suppose we have a sinusoid function

```
In [2]: f(x) = 10*\sin(x)
```

Out[2]: f (generic function with 1 method)

Our dataset $\mathcal D$ will consist of seven datapoints drawn from the following probabilistic model.

For each datapoint we randomly draw x_i uniformly in [0,6] and observe a noisy $y_i = f(x_i) + \epsilon_i$, where ϵ_i is some noise drawn from a standard normal distribution $\mathcal{N}(0,1)$.

Generate a sample dataset from this distribution.

```
In [147]: using Random
d = 6*rand()
d
```

Out[147]: 5.840382702837614

```
In [19]:  h = [[1,3],[4,5]]  h[1][1]
```

Out[19]: 1

```
In [370]:
           n = 7
           D = zeros(n, 2)
           for i = 1:n
               D[i,1] = 6*rand()
               D[i,2] = f(D[i,1]) + randn() # yi = f(xi) + \epsilon i
           end
           D
Out[370]: 7×2 Array{Float64,2}:
            0.0991965
                       0.642399
            2.07209
                        7.69209
                       -9.2755
            5.02041
            5.704
                       -3.88525
            2.86417
                        3.12726
            3.367
                        0.262463
                       -5.40448
            3.64152
           Plot the dataset \mathcal{D} and the true function f(x).
In [150]:
           """plot function y=f(x)"""
           function plotfunc first(f;
               xmin=0,xmax=6,nsamples=1000)
               xsamples = range(xmin,stop=xmax,length=nsamples)
               plot(xsamples, [f(x) for x in xsamples], color="black") ## only dif
           end
Out[150]: plotfunc first
           """plot function y=f(x)"""
In [151]:
           function plotfunc(f;
               xmin=0,xmax=6,nsamples=1000)
               xsamples = range(xmin,stop=xmax,length=nsamples)
               plot!(xsamples, [f(x) for x in xsamples], color="black")
           end
```

Out[151]: plotfunc

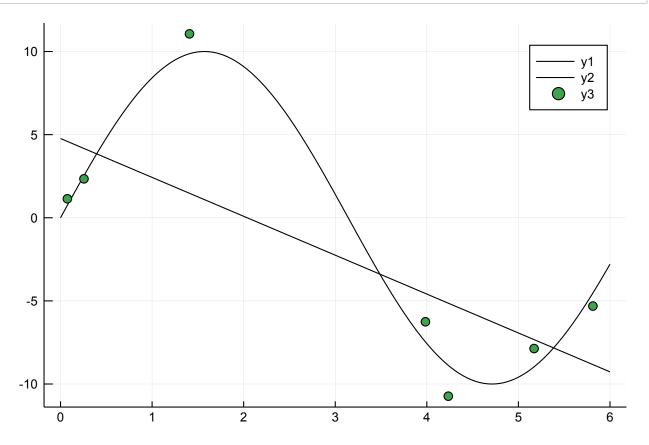
b)

Fit a linear model to \mathcal{D}

Plot the linear model l(x) together with \mathcal{D} and f(x). Feel free to use our method plotfunc(f).

```
In [156]: plotfunc_first(1)
    plotfunc(f)
    scatter!(D[:,1],D[:,2])
```

Out[156]:



c)

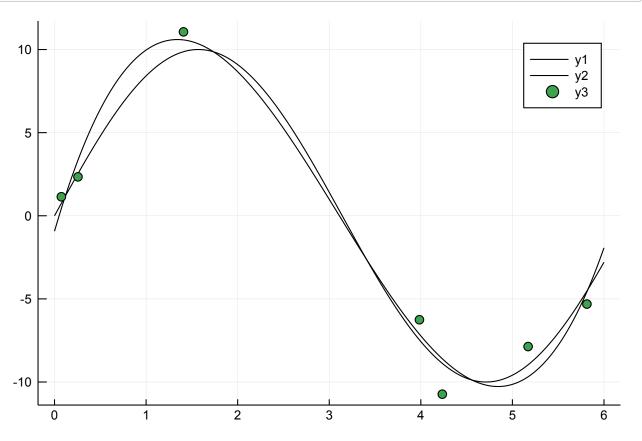
Fit a cubic model c(x) to ${\mathcal D}$

```
In [157]: # first, construct a Vandermonde matrix
          max order = 3
           x = D[:,1]
           V = zeros(n, max order+1)
           for k=0:max order
               V[:,k+1] = x.^k
           end
           # solve least squares problem
           w c = V \setminus D[:,2]
           W C
Out[157]: 4-element Array{Float64,1}:
           -0.9290704184314104
           18.94177395481076
           -9.017525168290522
             0.9721339006409941
In [158]: | function c(x; order = max_order, w = w_c)
               y = 0
               for k=0:order
                   y += w[k+1]*x^k
               end
               return y
           end
Out[158]: c (generic function with 1 method)
In [159]: c(0.5)
Out[159]: 6.408952004481463
```

Plot the cubic model with \mathcal{D} and f(x)

```
In [160]: plotfunc_first(c)
    plotfunc(f)
    scatter!(D[:,1],D[:,2])
```

Out[160]:



d)

Repat the parts (b) and (c) for 1000 different sets \mathcal{D} . Compute \bar{l} and \bar{c} , the average linear and average cubic models. (Please name \bar{l} as I_avg(x) and \bar{c} as c_avg(x) in following codes)

```
In [384]: max iter = 1000
          w l bar_all_iter = zeros((2, max_iter))
          for j = 1:max iter
              D l bar = zeros((n, 2))
              for i = 1:n
                  D \mid bar[i,1] = 6*rand()
                  D_l_{bar}[i,2] = f(D_l_{bar}[i,1]) + randn() # yi = f(xi) + \epsilon i
              end
              X_1_{bar} = [D_1_{bar}; 1] ones(7)]
              y_l_bar = D_l_bar[:,2]
              w l iter = X l bar\y l bar
              w l bar all_iter[:,j] = w_l_iter
          end
          w l bar all iter
Out[384]: 2×1000 Array{Float64,2}:
           -3.53113 -1.29912 -1.83742 -4.51579 ... -6.83591 -2.72421 -3.97
          078
            9.97079 -0.797265 6.57384 14.8425
                                                        19.7701
                                                                   7.50876 13.15
          63
In [385]: w l avg = mean(w l bar all iter, dims=2)
Out[385]: 2×1 Array{Float64,2}:
           -3.7246094918268278
           11.442821369125939
In [386]:
          function 1 avg(x)
              y_pred = [x 1]w_l_avg
              return( y pred[1])
          end
Out[386]: 1 avg (generic function with 1 method)
In [387]: 1 avg(0.5)
Out[387]: 9.580516623212525
```

```
In [388]:
          max iter = 1000
          max order = 3
          w c bar all iter = zeros((4, max iter))
          n = 7
           for j = 1:max iter
              D_c_{bar} = zeros((n, 2))
               for i = 1:n
                   D_c_{bar}[i,1] = 6*rand()
                   D c bar[i,2] = f(D c bar[i,1])+randn() # yi = f(xi) + \epsilon i
              end
              x = D c bar[:,1]
              #print(x)
              V = zeros(n, max order+1)
              for k=0:max order
                   V[:,k+1] = x.^k
              end
              # solve least squares problem
              w c iter = V \setminus D c bar[:,2]
              #print(w_c_iter)
              #w c iter
              w c bar all iter[:,j] = w c iter
          end
          w c bar all iter
Out[388]: 4×1000 Array{Float64,2}:
           -2.02624
                      -0.6236
                                   -7.48199 ... -1.2245
                                                            0.397898
                                                                        -2.13843
           17.8872
                     15.1926
                                   28.2868
                                                16.2158
                                                           15.085
                                                                        21.5857
                                                           -7.60372
           -8.62625 -7.29606
                                  -12.9076
                                                -7.51311
                                                                       -10.1894
            0.946467 0.786004
                                                0.789725
                                                            0.825389
                                    1.47059
                                                                         1.1185
In [389]: w c avg = mean(w c bar all iter, dims=2)
Out[389]: 4×1 Array{Float64,2}:
           -1.8907775436853607
           20.198840015115497
           -9.348139968761021
            0.9911502188652426
In [390]: c_{avg}(x) = c(x; order = max_order, w = w c avg)
Out[390]: c avg (generic function with 1 method)
```

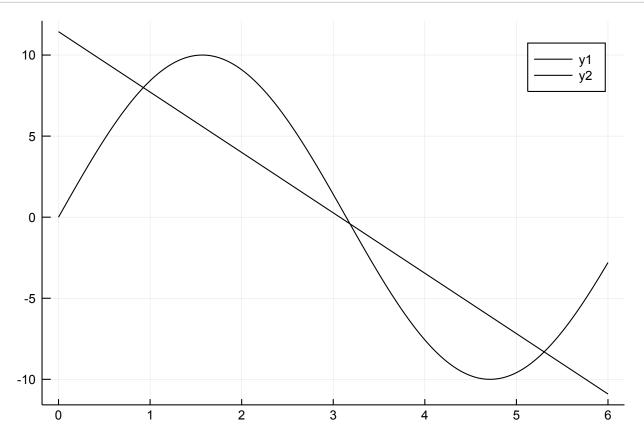
```
In [391]: c_avg(0.5)
```

Out[391]: 5.995501249040287

Plot \bar{l} together with f(x).

```
In [392]: plotfunc_first(l_avg)
    plotfunc(f)
```

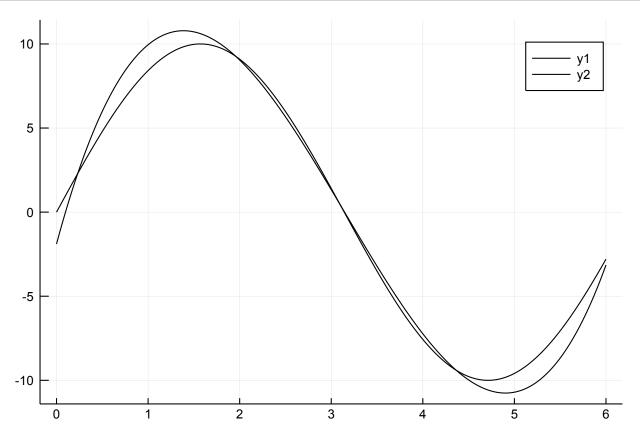
Out[392]:



Plot $\bar{c}(x)$ together with f(x).

```
In [393]: plotfunc_first(c_avg)
plotfunc(f)
```

Out[393]:



e)

Compute the bias of \overline{l} . You can use our integrate function.

```
In [436]: function integrate(f, a, b)
    n = 1000
    delta = (b - a)/n;  ## nothing to change below here
    xs = a*ones(n) + [0:1:n-1;] * delta;  ## n, right is 1:n * d
    fx = map(f, xs);
    return sum(fx) * delta
end
```

Out[436]: integrate (generic function with 1 method)

```
In [440]: g_1(x) = (f(x) - 1_avg(x))^2
```

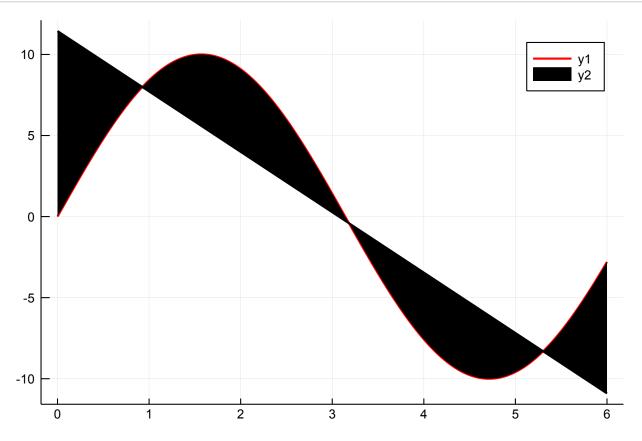
Out[440]: g_l (generic function with 1 method)

```
In [441]: bias_l_bar = integrate(g_l,0,6)
```

Out[441]: 104.67372562845834

```
In [442]: x_range = range(0,stop=6,length=1000)
x = [x for x in x_range]
y = f(x)
y_2 = [l_avg(x) for x in x_range]
plot(x, y, color="red", linewidth=2.0)
plot!(x, y_2, fillrange=[y y_2], color="black", linewidth=2.0)
```

Out[442]:



Compute the bias of the cubic model $\bar{c}(x)$.

```
In [443]: g_c(x) = (f(x) - c_avg(x))^2

Out[443]: g_c (generic function with 1 method)

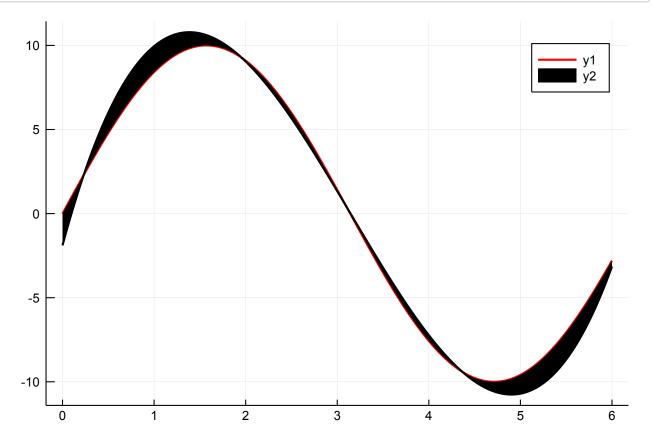
In [444]: bias_c_bar = integrate(g_c,0,6)
```

Out[444]: 4.63894175311889

We can interpret the bias as how far off our averaged model is from the true function. One way to visually see this is by plotting $\bar{l}(x)$ with f(x) and color in their difference. Try out the plotting function below.

```
In [445]: x_range = range(0,stop=6,length=1000)
    x = [x for x in x_range]
    y = f(x)
    y_2 = [c_avg(x) for x in x_range]
    plot(x, y, color="red", linewidth=2.0)
    plot!(x, y_2, fillrange=[y y_2], color="black", linewidth=2.0)
    #plot!([x y_2], fillrange=[y y_2], fillalpha=0.3, c=:orange)
```

Out[445]:



Which model has smaller bias?

Polynomial model has a smaller bias (of 4.6) as compared to Linear model (which has a bias of 104).



Next compute the variance of the linear model.

```
In [503]: x_range = range(0,stop=6,length=1000)
    x = [x for x in x_range]
    Var_l = 0

for i = 1:1000
    #print(x[i])
    x_l = [x[i] 1]
    #print(i, x_l)
    #print(i, w_l_bar_all_iter[:,i])
    y_jth_model_pred = x_l*w_l_bar_all_iter[:,i]
    #print(y_jth_model_pred)
    #print(w_l_avg)
    y_l_avg_pred = X_l*w_l_avg
    Var_l = Var_l + (y_jth_model_pred[1] - y_l_avg_pred[1])^2
    end

variance_linear = Var_l*6/1000 #integration factor (b-a/n)
```

Out[503]: 85.07503362489572

Compute the variance of the cubic model.

Out[504]: 343.0359823040535

1724

```
In [505]: total_error_linear = variance_linear + bias_l_bar ## I am doing square
    total_error_cubic = variance_cubic + bias_c_bar
    print("Total linear error: ",total_error_linear,"Total cubic error: ",to

Total linear error: 189.74875925335408Total cubic error: 347.674924057
```

Which model had higher variance? How do you interpret this? Which model has smaller overall error?

Cubic model has higher variance (343) as compared to linear model (85). Total error is high for cubic model (347) as compared to linear models (189).

This indicetes typical bias-variance tradeoff. Models with high values of bias (linear model in this case), undermine their training dataset (underfit) leading to not learn the trends in training dataset. consquently these models are simpler also. These models do not vary a great deal with change in training dataset.

High-variance are more complex and represent their training set well (at times overfit). However, these models are not able to handle random variation in test dataset and hence do not generalize well.

Ideally, one wants to choose a model with training error being low and generalize on out of set data as well - Alas! not easy to get there.

g)

How do you think your results would depend on the number of points in the data set \mathcal{D} ? Feel free to perform an experiment to check. How many points would you need before the opposite model has smaller overall error?

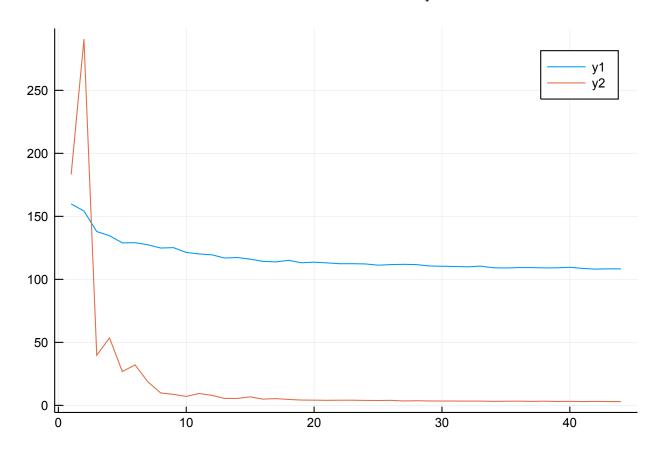
```
In [544]: 
    n_min = 7 # min data points
    n_max = 50 #max data points
    max_iter = 1000
    x_range = range(0, stop=6, length=max_iter)
    x = [x for x in x_range]
    n_iter = n_max-n_min+1
    max_order = 3
```

```
Variance matrix = zeros((2,n iter))
Bias matrix = zeros((2, n iter))
total error matrix = zeros((2, n iter))
for n = n min:n max
    ## Linear model calculations start here
    w l bar all iter = zeros((2, max iter))
    w c bar all iter = zeros((4, max iter))
    for j = 1:max iter
        D bar = zeros((n, 2))
        for i = 1:n
            D bar[i,1] = 6*rand()
            D_bar[i,2] = f(D_bar[i,1]) + randn() # yi = f(xi) + \epsilon i
        end
        X l bar = [D_bar[:,1] ones(n)]
        y 1 bar = D bar[:,2]
        w l iter = X l bar\y l bar
        w l bar all iter[:,j] = w l iter
        b = D bar[:,1]
        V = zeros(n, max order+1)
        for k=0:max order
            V[:,k+1] = b.^k
        end
        w c iter = V\D bar[:,2]
        w c bar all iter[:,j] = w c iter
    end
    w l avg = mean(w l bar all iter, dims=2)
    Var 1 = 0
   Bias l = 0
    w c avg = mean(w c bar all iter, dims=2)
    Var c = 0
    Bias c = 0
    for k = 1:max iter
        X 1 = [x[k] 1]
        y jth model pred = X 1*w l bar all iter[:,k]
        y l avg pred = X 1*w l avg
        Var_l = Var_l + (y_jth_model_pred[1] - y_l_avg_pred[1])^2
        Bias l = Bias l + (f(x[k][1])-y l avg pred[1])^2
        X c = x[k]
        y_jth_model_pred = c(X_c, w=w_c_bar_all_iter[:,k])
        y_c_avg_pred = c(X_c, w=w_c_avg)
```

```
Var c = Var_c + (y_jth_model_pred[1] - y_c_avg_pred[1])^2
        Bias_c = Bias_c + (f(x[k][1])-y_c_avg_pred[1])^2
    end
    Variance matrix[1, n-n min+1] = Var l*(6/max iter)
                                                         ##integration fa
                                                         ##integration fa
    Bias matrix[1,n-n min+1] = Bias l*(6/max iter)
    Variance matrix[2, n-n min+1] = Var c*(6/max iter)
    Bias matrix[2, n-n min+1] = Bias c*(6/max iter)
end
total error matrix = Variance matrix + Bias matrix
#print(Variance matrix)
#print(Bias matrix)
print(total error matrix[:,1:5])
plot(total error matrix[1,:])
plot!(total error matrix[2,:])
```

[159.85614557980801 154.26763206676978 137.91061797595788 134.64880317 859985 128.9448314153106; 183.2304051817007 290.5232693648684 39.78711 971313455 53.50197959661283 26.843959071250108]

Out[544]:



As number of datapoints in D increases, bias comes down significantly. As n > 9, opposite model (in this case Cubic model) has lower total error than part (f). Errors for both models flattens out after 30 odd data points.

h)

Instead of sampling new data to compute the bias and variance of our model, we could use a bootstrap estimator to get more use out of the few data points we have. Try this for a few different data set sizes and report on your results. How big a data set is needed for the bootstrap to give a reliable estimate of the bias and variance?

We can borrow some learnings from part g, where we saw affect of size of dataset (n) on total error. Total errors flattens after a certain point (somewhere around n= 27).

We can further study impact of n on bootstrapping by taking three values of n. n = 7, 15, 25.

Also I will be considering number of datapoints taken, k = 100.

Below we will be looking at the Confidence interval for each value of n. Half width for 95% Confidence interval should keep reducing as n increases.

I will be restricting this exercise of bootstrapping to linear models and look at variation of first cofficent of w in those models. Conclusions from this should hold for other models and other values of w as well.

```
In [563]: n h = 35
           K = 1000
           max iter = K
           w l bar all iter = zeros((2, max iter))
           for j = 1:max iter
               D l bar = zeros((n h, 2))
               for i = 1:n h
                   D l bar[i,1] = 6*rand()
                    D_1_{bar[i,2]} = f(D_1_{bar[i,1]}) + randn() # yi = f(xi) + \epsilon i
               end
               X \mid bar = [D \mid bar[:,1] \text{ ones}(n \mid h)]
               y 1 bar = D 1 bar[:,2]
               w l iter = X l bar\y l bar
               w_l_bar_all_iter[:,j] = w l iter
           end
           z = w l bar all iter[1,:] #first cofficent of w
           UB = mean(z) + 1.96*sqrt(var(z)/K) # 95% confidence interval
           LB = mean(z) - 1.96*sqrt(var(z)/K)
           print("n = ", n h, "; k = ", K, "; Confidence interval = ", LB,",", UB,
```

n = 35; k = 1000; Confidence interval = -3.5192832189851346, -3.4575042 26763966; Half width = 0.030889496110584336

n = 7; k = 100; Confidence interval = -3.8342240783025403,-3.301831389430105; Half width = 0.26619634443621765

n = 15; k = 100; Confidence interval = -3.807226401341788,-3.4520754000707012; Half width = 0.17757550063554328

n = 25; k = 100; Confidence interval = -3.6357753693738997,-3.4047754358186473; Half width = 0.1154999667776262

n = 35; k = 100; Confidence interval = -3.5271356160037763,-3.3509984924327982; Half width = 0.08806856178548905

n = 100; k = 100; Confidence interval = -3.524302390585087,-3.4115248830183442; Half width = 0.05638875378337138

n = 35; k = 10; Confidence interval = -3.7530071105002247,-3.0721762268779265; Half width = 0.34041544181114913

n = 35; k = 50; Confidence interval = -3.588653833079747,-3.2770852642801134; Half width = 0.15578428439981673

n = 35; k = 100; Confidence interval = -3.6330291669220243,-3.4462648808811265; Half width = 0.0933821430204489

n = 35; k = 1000; Confidence interval = -3.5192832189851346,-3.457504226763966; Half width = 0.030889496110584336

As we can see from above data points:

- 1. As n and k increases half width reduces, which means with more confidence we can locate true value of w[1].
- 2. this makes intuitive sense, as higher the K (number of samples withdrawn), more close the estimates will be to actual.
- 3. For our use case n = 25-35 and K~100 should give 95% confidence interval of half width within 1.5% of the mean can be achieved.

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