DSC 441 – FUNDAMENTAL OF DATA SCIENCE

HOMEWORK 4

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PROBLEM 1

For this problem, you will tune and apply kNN and compare it to other classifiers. We will use the wine quality data, which has a number of measurements about chemical components in wine, plus a quality rating. There are separate files for red and white wines, so the first step is some data preparation

a. For this problem, you will tune and apply kNN and compare it to other classifiers. We will use the wine quality data, which has a number of measurements about chemical components in wine, plus a quality rating. There are separate files for red and white wines, so the first step is some data preparation.

```
```{r}
Importing the datasets and separating them with a semicolon
winequality_white <- read.csv("winequality-white.csv", sep = '
winequality_red <- read.csv("winequality-red.csv", sep = ";")</pre>
summary(winequality_white)
 fixed.acidity
 volatile.acidity citric.acid
 Min. : 3.800
1st Qu.: 6.300
 Min. :0.00900
1st Qu.:0.03600
 Min. :0.0800
1st Qu.:0.2100
 Min. :0.0000
1st Qu.:0.2700
 Min. : 0.600
1st Qu.: 1.700
 Min.
 Median : 6.800
 Median :0.2600
 Median :0.3200
 Median : 5.200
 Median :0.04300
 :0.04577
 Mean
 : 6.855
 Mean
 :0.2782
 Mean
 :0.3342
 Mean
 : 6.391
 Mean
 3rd Qu.: 7.300
 3rd Qu.:0.3200
 3rd Qu.:0.3900
 3rd Qu.: 9.900
 3rd Qu.:0.05000
 :1.6600
 :0.34600
 :14.200
 Max.
 :1.1000
 Max.
 Max.
 :65.800
 Max.
 pH
Min.
 free.sulfur.dioxide total.sulfur.dioxide
 density
 sulphates
 Min.
 Min.
 2.00
 Min.
 :0.9871
 :2.720
 Min.
 1st Qu.:0.9917
Median :0.9937
 1st Qu.:0.4100
 1st Qu.: 23.00
 1st Qu.:108.0
 1st Qu.:3.090
 Median: 34.00
 Median :134.0
 Median :3.180
 Median :0.4700
 Mean
 : 35.31
 Mean
 :138.4
 Mean
 :0.9940
 Mean
 :3.188
 Mean
 3rd Qu.: 46.00
 3rd Qu.:167.0
 3rd Qu.:0.9961
 3rd Qu.:3.280
 3rd Qu.:0.5500
 :289.00
 Max.
 Max.
 :440.0
 Max.
 :1.0390
 Max.
 :3.820
 Max.
 alcohol
 quality
 : 8.00
 Min. : 8.00
1st Qu.: 9.50
 Min. :3.000
1st Qu.:5.000
 Median:10.40
 Median :6.000
 :5.878
 Mean
 :10.51
 Mean
 3rd Ou.:11.40
 3rd Ou.:6.000
```

```
summary (winequality_red)
 volatile.acidity citric.acid
Min. :0.1200 Min. :0.000
1st Qu.:0.3900 1st Qu.:0.090
 fixed.acidity
 residual.sugar
 chlorides
 Min. : 4.60
1st Qu.: 7.10
Median : 7.90
Mean : 8.32
 Min. : 0.900
1st Qu.: 1.900
 Min. :0.1200
1st Qu.:0.3900
 Min. :0.01200
1st Qu.:0.07000
 Median :0.5200
 Median : 2.200
 Median :0.07900
 Median :0.260
 Mean
 :0.5278
 Mean
 Mean
 Mean
 Mean
 3rd Qu.: 9.20
 3rd Qu.:0.6400
 3rd Qu.:0.420
 3rd Qu.: 2.600
 3rd Qu.:0.09000
 Max.
 :1.5800 Max.
 :15.90
 Max. :15.500
 Max.
 :1.000
 Max. :0.61100
 pH
Min. :?
 free.sulfur.dioxide total.sulfur.dioxide
 density
 sulphates
 Min.
 Min. : 6.00
1st Qu.: 22.00
 :0.9901
 :2.740
 Min. : 1.00
1st Qu.: 7.00
 Min. :0.9901
1st Qu.:0.9956
 Min. :0.3300
1st Qu.:0.5500
 1st Qu.:3.210
 Median :14.00
 Median : 38.00
 Median :0.9968
 Median :3.310
 Median :0.6200
 Median :0.3355
Mean :0.9967
3rd Qu::0.9978
 Mean :15.87
3rd Qu.:21.00
 Mean : 46.47
3rd Qu.: 62.00
 Mean
 Mean :3.311
 Mean
 :0.6581
 3rd Qu.:3.400
 3rd Qu.:0.7300
 :289.00
 Max. :1.0037
 Max.
 quality
Min. :3.000
1st Qu.:5.000
 alcohol
 Min.
 : 8.40
 1st Qu.: 9.50
 Median :10.20
 Median :6.000
 :10.42
 :5.636
 Mean
 Mean
 3rd Qu.:11.10
 3rd Qu.:6.000
 Max.
 :14.90
 Max.
 :8.000
```

Now, check the type of data in the columns of the wines dataset

```
> typeof(winequality_red$chlorides)
[1] "double"
> typeof(winequality_red$citric.acid)
[1] "double"
> typeof(winequality_red$residual.sugar)
[1] "double"
> typeof(winequality_red$alcohol)
[1] "double"
```

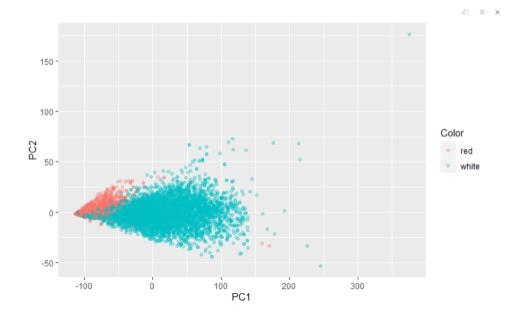
Note that there is no NAs in the dataset. Now, let's add a column for the types of wines.

```
Add the column for type of wine
winequality_red$type <- c('red')
winequality_white$type <- c('white')
Now combine both tables using the full_join
wines <- full_join(winequality_red, winequality_white)
head(wines)</pre>
```

	fixed.acidity «dbl>	volatile.acidity «dbl»	citric.acid <dbl></dbl>	residual.sugar <dbl></dbl>	chlorides «dol»	free.sulfur.dioxide	total.sulfur.dioxide <dbl></dbl>	density «dbl»	<dbl></dbl>
1	7.4	0.70	0.00	1.9	0.076	11	34	0.9978	3.51
2	7.8	0.88	0.00	2.6	0.098	25	67	0.9968	3.20
3	7.8	0.76	0.04	2.3	0.092	15	54	0.9970	3.26
4	11.2	0.28	0.56	1.9	0.075	17	60	0.9980	3.16
5	7.4	0.70	0.00	1.9	0.076	11	34	0.9978	3.51
6	7.4	0.66	0.00	1.8	0.075	13	40	0.9978	3.51

b. Use PCA to create a projection of the data to 2D and show a scatterplot with color showing the wine type.

```
fraction of the dataset
dummy <- dummyVars(type ~ ., data = wines)
dummies <- as.data.frame(predict(dummy, newdata = wines))
set.seed(123)
Calculate PCA
pca = prcomp(dummies)
Save as data frame
rotated_data = as.data.frame(pca$x)
Add the original label 'type' as a reference
rotated_data$Color <- wines$type
Plot and color the labels based on wine type red (or) white
ggplot(data = rotated_data, aes(x = PC1, y = PC2, col = Color)) + geom_point(alpha = 0.3)</pre>
```



# c. We are going to try kNN, SVM and decision trees on this data. Based on the 'shape' of the data in the visualization from (b), which do you think will do best and why?

I think the best method in this case may be to use KNN. The wines dataset is somewhat large so KNN may be faster and more efficient solution. Also, the non-linear relationship will allow for a more accurate prediction by using the nearest K neighbour to predict the type of wine. I've performed some sample work below that helped me also get to this conclusion, although SVM and KNN performed similarly in regards to accuracy. However, SVM's patterns may be limited in this case (we can discover this later).

#### Finding the K nearest neighbour:

```
k-Nearest Neighbors

6497 samples
12 predictor
2 classes: 'red', 'white'

Pre-processing: centered (12), scaled (12)
Resampling: Cross-Validated (10 fold)
Summary of sample sizes: 5848, 5847, 5847, 5848, 5847, 5848, ...
Resampling results across tuning parameters:

k Accuracy Kappa
5 0.9923051 0.9792070
7 0.9929205 0.9808827
9 0.9930741 0.9813154

Accuracy was used to select the optimal model using the largest value. The final value used for the model was k = 9.
```

#### **SVM:**

```
Fit the Model
svm1 <- train(type ~., data = wines, method = "svmLinear")
Evaluate the Fit
svm1

Support Vector Machines with Linear Kernel

6497 samples
12 predictor
2 classes: 'red', 'white'

No pre-processing
Resampling: Bootstrapped (25 reps)
Summary of sample sizes: 6497, 6497, 6497, 6497, 6497, ...
Resampling results:
 Accuracy Kappa
 0.994937 0.9862915

Tuning parameter 'C' was held constant at a value of 1</pre>
```

#### **Decision Tree:**

```
Evaluation Method
train control = trainControl(method = "cv", number = 10)
Fit the Model
tree1 <- train(type ~., data = wines, method = "rpart", trControl = train control)
tree1
 CART
 6497 samples
 12 predictor
 2 classes: 'red', 'white'
 No pre-processing
 Resampling: Cross-Validated (10 fold)
 Summary of sample sizes: 5847, 5848, 5847, 5848, 5847, ...
 Resampling results across tuning parameters:
 cp Accuracy Kappa
0.06253909 0.9465933 0.8498495
 0.06754221 0.9348963 0.8147365
 0.70043777 0.7855788 0.1500825
 Accuracy was used to select the optimal model using the largest value.
 The final value used for the model was cp = 0.06253909.
```

d. Use kNN (tune k), use decision trees (basic rpart method is fine), and SVM (tune C) to predict type from the rest of the variables. Compare the accuracy values – is this what you expected? Can you explain it? Note: you will need to fix the columns names for rpart because it is not able to handle the underscores. This code will do the trick (assuming you called your data wine\_quality): colnames(wine\_quality) <- make.names(colnames(wine\_quality)).</p>

#### KNN:

```
k-Nearest Neighbors
6497 samples
 12 predictor
 2 classes: 'red', 'white'
Pre-processing: centered (12), scaled (12)
Resampling: Cross-Validated (10 fold)
Summary of sample sizes: 5847, 5847, 5848, 5849, 5847, 5847, ...
Resampling results across tuning parameters:
 Accuracy
 Kappa
 0.9936890
 0.9933813
 0.9821286
 11
 0.9927659 0.9805034
 0.9926121
 0.9801038
 0.9923044 0.9792783
 0.9927659
 0.9804950
 19
 0.9926116
 0.9800879
 0.9929198
 0.9809104
 0.9924582
 0.9796711
 0.9923044
 0.9792539
 0.9923044 0.9792539
 31
 0.9916885 0.9776041
 33 0.9918428 0.9780181
Accuracy was used to select the optimal model using the largest value.
```

#### For k = 7:

```
fit <- kmeans(dummies, centers = 7, nstart = 25) # Display the kmeans object information fit
```

The final value used for the model was k = 7.

```
K-means clustering with 7 clusters of sizes 856, 910, 354, 717, 1382, 1131, 1147
 425 426 427 428 44 4 2 4 4 5 4 4 5 4 4 5 4 4 5 4 4 5 4 4 5 4 4 5 4 4 5
 429
2
448
2
467
4
486
2
505
2
524
 430
2
449
2
468
2
487
2
506
 431
2
450
2
469
2
488
2
507
 432
451
2
470
4
489
4
508
 434
2
453
2
472
4
491
4
510
4
529
 452
 471
4
490
2
509
 525
 2 538 2 2 5577 2 2 5587 4 595 5 2 6114 4 652 2 709 2 2 747 4 6 6 4 785 6 804 823
 5
544
4
563
 4 548 2 5866 2 605 4 4 624 4 4 681 2 700 2 719 2 776 2 2 776 2 2 833
 7
543
 2
562
5
581
2
600
2
619
2
638
 2
575
6
594
2
613
2
632
2
651
2
670
2
689
 5
582
2
601
2
620
2
639
2
658
4
677
2
696
 17
6
36
2
55
4
74
4
93
5
112
6
131
5
 610
2
629
2
648
2
667
 16 7 7 35 4 4 54 4 54 4 55 4 4 1300 2 2 1499 2 2 225 4 2 2 263 3 2 2 282 2
 4
40
6
59
4
78
2
97
2
 4
628
2
647
2
666
 7
657
2
676
2
695
5
714
4
733
2
752
2
771
6
790
5
809
828
 685
 686
 2
705
2
724
5
743
2
762
2
781
4
800
2
819
 715
4
734
4
753
4
772
 4
135
 746
2
765
4
784
2
803
6
 169
4
188
2
207
2
226
 2
154
6
173
2
192
4
211
2
230
2
 7
791
6
810
2
829
 245
245
264
 838
6857
46876
2895
6914
2933
4952
2971
2990
 4
841
2
860
2
879
4
898
4
917
4
936
4
955
2
974
2
 2
842
2
861
6
880
6
899
4
918
4
937
4
956
2
975
2
 4
852
2
871
4
890
6
909
4
928
4
947
2
966
4
985
 282 283

2 4

301 302

4 2

320 321

4 4

339 340

6 2

358 359

2 2

377 378

4 2

396 397

2 5

415 416

5 5
 2
287
4
306
2
325
4
344
2
363
2
382
4
 Within cluster sum of squares by cluster:

[1] 330867.7 108090.8 367269.3 147020.0 301721.7 206162.7 340030.8

(between_SS / total_SS = 92.2 %)
 Available components:
 [1] "cluster"
[7] "size"
 "totss"
 "centers"
 "withinss"
 "tot.withinss" "betweenss"
SVM:
 # I decided to use the Grid Search here to try different values of C
 grid \leftarrow expand.grid(C = 10^seq(-5, 2, 0.5))
 Fit the Model
 svm_grid <- train(type ~., data = wines, method = "svmLinear",
 trControl = train_control, tuneGrid = grid)
 # View grid search result
 svm grid
 Support Vector Machines with Linear Kernel
 6497 samples
 12 predictor
2 classes: 'red', 'white'
 No pre-processing
Resampling: Cross-Validated (10 fold)
Summary of sample sizes: 5847, 5847, 5847, 5848, 5847, 5847, ...
Resampling results across tuning parameters:
 Accuracy
0.7538866
0.7541943
 1.000000e-05
3.162278e-05
 0.000000000
0.001878654
 0.9302726
0.9839943
 0.792017114
0.956278082
 1.000000e-04
 3.162278e-04
 1.000000e-03
 0.9906118
 0.974623044
 3.162278e-03
1.000000e-02
3.162278e-02
 0.9915354
0.9926128
 0.977177217
 0.9938435
 0.983410595
 1.000000e-01
3.162278e-01
 0.9946132
0.9947671
```

0.985887888

0.987123844

0.986706680

Accuracy was used to select the optimal model using the largest value. The final value used for the model was  $C\,=\,1.$ 

0.9950750 0.986706680 0.9950750 0.986706680

0.9952289

0.9950750

**Decision Trees:** 

1.000000e+00 3.162278e+00

1.000000e+01

3.162278e+01 1.000000e+02

```
Evaluation Method
train_control = trainControl(method = "cv", number = 10)
Fit the Model
tree1 <- train(type ~., data = wines, method = "rpart", trControl = train_control)
Evaluate the Fit
tree1</pre>
```

```
CART

6497 samples
12 predictor
2 classes: 'red', 'white'

No pre-processing
Resampling: Cross-Validated (10 fold)
Summary of sample sizes: 5847, 5847, 5847, 5848, 5847, ...
Resampling results across tuning parameters:

cp Accuracy Kappa
0.06253909 0.9448994 0.8427741
0.06754221 0.9344300 0.8125794
0.70043777 0.8370148 0.3857512

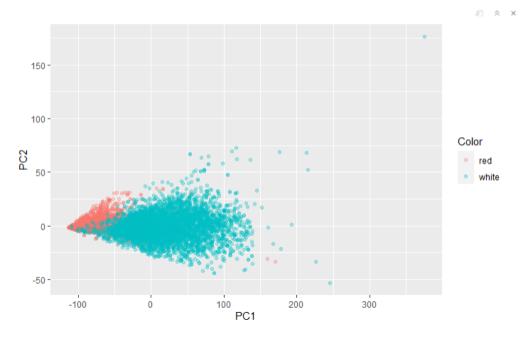
Accuracy was used to select the optimal model using the largest value.
The final value used for the model was cp = 0.06253909.
```

Accuracy seems to be comparable in SVM and KNN, although not decision trees. I think the dataset may be too complex/large for a decision tree to handle. However, SVM and KNN seem to perform better. Although, due to the non-linear nature of this dataset, probably a KNN solution may be the best.

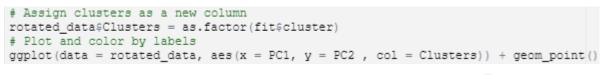
e. Use the same already computed PCA again to show a scatter plot of the data and to visualize the labels for kNN, decision tree and SVM. Note that you do not need to recreate the PCA projection, you have already done this in 1b. Here, you just make a new visualization for each classifier using its labels for color (same points but change the color). Map the color results to the classifier, that is use the "predict" function to predict the class of your data, add it to your data frame and use it as a color. This is done for KNN in the tutorial, it should be similar for the others. Consider and explain the differences in how these classifiers performed.

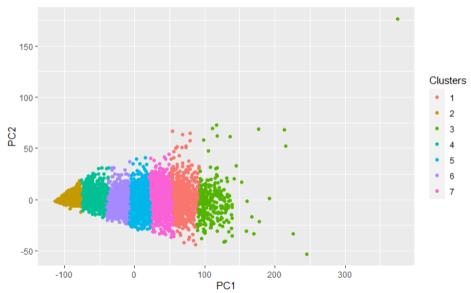
#### **Using PCA:**

```
Plot and color the labels based on Wine type red (or) white
ggplot(data = rotated_data, aes(x = PC1, y = PC2, col = Color)) + geom_point(alpha = 0.3)
```



#### **Using KNN:**





PCA seemed to depict the clusters into two uneven clusters, although we are aware that there is far more white wine samples than there is red wine in the dataset. The KNN clustering method provided an interesting distribution, although I prefer PCA as it seems to make more sense with the predicting of the red (or) white wine type. I think trying a different K value, perhaps smaller, may provide us with the better results. For the sake of testing it, I will try k=2 below, which seems to cluster the data more realistically:

```
fit2 <- kmeans(dummies, centers = 2, nstart = 25)
Assign clusters as a new column
rotated data$Clusters = as.factor(fit2$cluster)
Plot and color by the labels
ggplot(data = rotated_data, aes(x = PC1, y = PC2, col = Clusters)) + geom_point()
 150
 100
 Clusters
 PC2
 • 1
 • 2
 -50
 -100
 100
 200
 300
 PC1
```

#### PROBLEM - 2

In this question we will use the Sacramento data, which covers available housing in the region of that city. The variables include numerical information about the size of the housing and its price, as well as categorical information like zip code (there are a large but limited number in the area), and the type of unit (condo vs house (coded as residential))

```
'``{r}
data("Sacramento")
Remove the Zipcode, and lon for simplicity
cpsacramento <- Sacramento %>% select(-c("latitude", "longitude", "zip"))
'``
```

a. Load the data from the tidyverse library with the data ("Sacramento") command and you should have a variable Sacramento. Because we have categoricals, convert them to dummy variables.

```
Type is the largest variable to predict
dummy <- dummyVars(type ~., data = cpsacramento)
dummies <- as.data.frame(predict(dummy, newdata = cpsacramento))
head(dummies)</pre>
```



b. With kNN, because of the high dimensionality, which might be a good choice for the distance function?

It's hard to tell without actually trying the different metrics, although using Minkowski is a fairly common in high dimensional data.

c. Use kNN to classify this data with type as the label. Tune the choice of k plus the type of distance function. Report your results – what values for these parameters were tried, which were chosen, and how did they perform with accuracy?

```
Move the Type back to the dataset
Move the type back to the dataset
sacramento_dummies <- dummies
sacramento_dummies@type <- Sacramento@type
library(kknn)
library(kknn)
Setup a tuneGrid with the tuning parameters
tuneGrid <- expand.grid(kmax = 3:7, # Test a range of k values 3 to 7
kernel = c("rectangular","cos"), # Regular and cosine-based distance funtions
distance = 1:3) # Powers of Minkowski 1 to 3
 * Italia and It the model with 10-1010 cross valid

* Standardization, and our specialized tune grid

kknn_fit <- train(type ~.,

data = sacramento_dummies,
 method = 'kknn',
trControl = ctrl,
preProcess = c('center', 'scale'),
tuneGrid = tuneGrid)
 # Printing trained model provides report
 k-Nearest Neighbors
 932 samples
 1 predictor
3 classes: 'Condo', 'Multi_Family', 'Residential'
 Pre-processing: centered (41), scaled (41)
Resampling: Cross-Validated (10 fold)
Summary of sample sizes: 838, 839, 839, 838, 838, 840, ...
Resampling results across tuning parameters:
 distance Accuracy
1 0.9367516
2 0.9367631
 kmax kernel
 Kappa
0.3884981
0.3763622
0.3763622
0.4495401
0.4298051
0.4205730
0.3884981
0.3763622
0.3763622
0.4619788
 rectangular 1
rectangular 2
 0.9367631
0.93567631
0.9355756
0.9335256
0.9324500
0.9367516
0.9367631
0.9367631
0.9388907
0.9399660
 rectangular 3
 COS
 cos
 0.9399889
 0.4391495
0.3360441
 rectangular 1
 0.9346011
 0.9346011
0.9303458
0.9303458
0.9388907
0.9356992
0.9378612
0.9346011
0.9303458
 rectangular
rectangular
cos
 0.2866676
 cos
rectangular
 0.2866676
 rectangular
 rectangular
 0.9303458
 0.2866676
 cos
 0.9367631
 0.4273555
 0.9346240 0.3705842
0.9378612 0.3943742
0.9324620 0.3059469
0.9335716 0.2803458
0.9346354 0.2835647
0.9367631 0.4273555
0.9346240 0.3705842
0.9389482 0.4006952
 rectangular 1
rectangular 2
rectangular 3
 COS
 Accuracy was used to select the optimal model using the largest value. The final values used for the model were kmax = 4, distance = 3 and kernel = cos.
```

Here we tried different Minkowski distances from 1 to 3, and tested regular and cosine distance functions. K-max of 6, with a distance of 2 (Euclidean), and a cosine based distance function were the results. The cosine's distance function kappa value generally performed better, although accuracy was comparable on both.

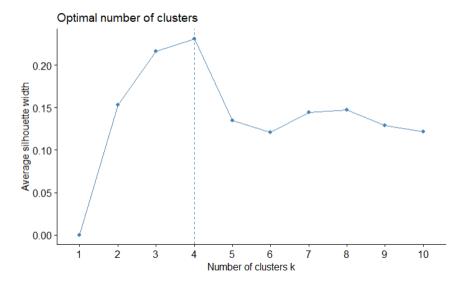
#### PROBLEM - 3

In this problem we will continue with the wine quality data from Problem 1, but this time we will use clustering. Do not forget to remove the type variable before clustering because that would be cheating by using the label to perform clustering.

```
Copy wines dataset and remove type cpwines <- wines cpwines <- cpwines %>% select(-c("type"))
```

a. Use k-means to cluster the data. Show your usage of silhouette and the elbow method to pick the best number of clusters. Make sure it is using multiple restarts.

```
df <- cpwines
Set seed
set.seed(123)
Center scale allows us to standardize the data
preproc <- preProcess(df, method = c("center", "scale"))
We have to call predict to fit our data based on preprocessing
predictors <- predict(preproc, df)
Find the knee
fviz nbclust(predictors, kmeans, method = "wss")
 Optimal number of clusters
 80000
 70000
 Total Within Sum of Square
 60000
 50000
 40000
 2
 5
 6
 8
 10
 Number of clusters k
 fviz nbclust(predictors, kmeans, method = "silhouette")
```



The silhouette method suggests a k = 4, and the elbow method also suggests a k = 4 may be reasonable.

```
Fit the Data
 fit <- kmeans(predictors, centers = 4, nstart = 25)
 # Display the k means object information
K-means clustering with 4 clusters of sizes 661, 1935, 1066, 2835
 fixed.acidity volatile.acidity citric.acid residual.sugar chlorides free.sulfur.dioxide

1.95495218 0.4048784 1.02423958 -0.5697395 1.2673641 -0.89472554

-0.19278454 -0.3500900 0.24718059 1.1555186 -0.1081800 0.82522737

0.05693398 1.5882171 -1.20795561 -0.6076496 0.6266271 -0.76925628
 1.1525186 -0.1081800
-0.6076496 0.6266271
-0.4253161 -0.4572783
 -0.34563560
 -0.4526420 0.04668919
 -0.06538772
 -0.33263500 pH

total.sulfur.dioxide density pH

-1.24231768 0.9050474 -0.11205327

0.95020983 0.7315214 -0.38723328

-1.03454506 0.4579506 0.88616595

0.03010265 -0.8825064 -0.04277873
 -0.425324 -0.4371265 -0.4371265 -0.4251265 -0.425126 -0.23647183 -0.2693766 -0.80429691 -0.29268305 -0.3628512 -0.29754458 -0.57836111 -0.2700050 0.64611743 0.40873607
Within cluster sum of squares by cluster:
[1] 8421.783 13176.504 7011.810 17841.974
(between_SS / total_SS = 40.4 %)
Available components:
[1] "cluster"
[7] "size"
 "centers"
"iter"
 "withinss"
 "tot.withinss" "betweenss'
```

b. Use hierarchical agglomerative clustering (HAC) to cluster the data. Try at least 2 distance functions and at least 2 linkage functions (cluster distance functions), for a total of 4 parameter combinations. For each parameter combination, perform the clustering.

#### **Euclidean and complete linkage:**

```
dist_mat <- dist(predictors, method = 'euclidean')
Determine assembly/agglomeration method and run hclust
hfit1 <- hclust(dist_mat, method = 'complete')
hfit1

Call:
 hclust(d = dist_mat, method = "complete")

Cluster method : complete
 Distance : euclidean
 Number of objects: 6497</pre>
```

#### **Euclidean and average linkage:**

```
dist_mat <- dist(predictors, method = 'euclidean')
Determine assembly/agglomeration method and run hclust
hfit2 <- hclust(dist_mat, method = 'average')
hfit2

Call:
 hclust(d = dist_mat, method = "average")

Cluster method : average
 Distance : euclidean
 Number of objects: 6497</pre>
```

#### Manhattan and complete linkage:

```
dist_mat <- dist(predictors, method = 'manhattan')
Determine assembly/agglomeration method and run hclust (average uses mean)
hfit3 <- hclust(dist_mat, method = 'complete')
hfit3

Call:
 hclust(d = dist_mat, method = "complete")

Cluster method : complete
 Distance : manhattan
 Number of objects: 6497</pre>
```

#### Manhattan and average linkage:

```
dist_mat <- dist(predictors, method = 'manhattan')
Determine assembly/agglomeration method and run hclust (average uses mean)
hfit4 <- hclust(dist_mat, method = 'average')
hfit4

Call:
 hclust(d = dist_mat, method = "average")

Cluster method : average
 Distance : manhattan
 Number of objects: 6497

Build the new model
 h1 <- cutree(hfit1, k=4)
 h2 <- cutree(hfit2, k=4)
 h3 <- cutree(hfit3, k=4)
 h4 <- cutree(hfit4, k=4)</pre>
```

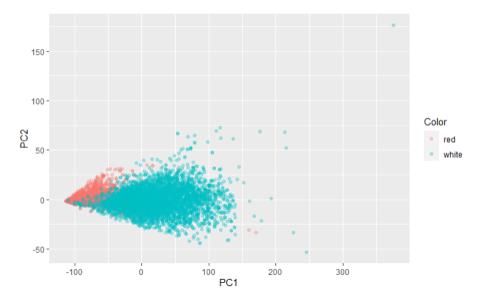
### c. Compare the k-means and HAC clusterings by creating a crosstabulation between their labels.

```
#Redefining the fit (I think I missed it previously)
 fit <- kmeans(predictors, centers = 4, nstart = 25)
 #Create a dataframe for the results
 result1 <- data.frame(WineType = wines$type, HAC1 = h1, Kmeans = fit$cluster)
 result2 <- data.frame (WineType = wines$type, HAC2 = h2, Kmeans = fit$cluster) result3 <- data.frame(WineType = wines$type, HAC3 = h3, Kmeans = fit$cluster)
 result4 <- data.frame(WineType = wines$type, HAC4 = h4, Kmeans = fit$cluster)
 #Crosstab for HAC
 result1 %>% group_by(HAC1) %>% select(HAC1, WineType) %>% table()
 WineType
 HAC1 red white
 1 1597 4896
 2
 0
 3
 0
 1
 4
 0
 1
result2 %>% group_by(HAC2) %>% select(HAC2, WineType) %>% table()
 WineType
 HAC2 red white
 1 1575 4895
 24
 3
 0
 1
 0
 1
result3 %>% group_by(HAC3) %>% select(HAC3, WineType) %>% table()
 WineType
 HAC3 red white
 1 1595 4889
 0
 3
 0
 8
 4
 0
 1
result4 %>% group_by(HAC4) %>% select(HAC4, WineType) %>% table()
 WineType
 HAC4 red white
 1 1576 4896
 2
 23
 0
 1
 3
 0
 0
 #Crosstab for K Means
 result <- data.frame(Type = wines$type, Kmeans = fit$cluster)
 result %>% group by (Kmeans) %>% select(Kmeans, Type) %>% table()
 Type
 Kmeans red white
 3 1932
 2 609
 52
 3
 927
 139
 60 2775
```

d. For comparison – use PCA to visualize the data in a scatterplot. Create 3 separate plots: use the color of the points to show (1) the type label, (2) the kmeans cluster labels and (3) the HAC cluster labels.

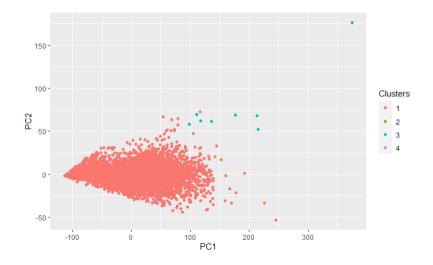
#### PCA:

```
#Recreating the PCA scatter plot
#Create Dummies
dummy <- dummyVars(type ~ ., data = wines)
dummies <- as.data.frame(predict(dummy, newdata = wines))
set.seed(123)
#Calculate PCA
pca = prcomp(dummies)
#Save as data frame
rotated_data = as.data.frame(pca$x)
#Add original label 'type' as a reference
rotated_data$Color <- wines$type
#Plot and color the labels based on wine type red (or) white
ggplot(data = rotated_data, aes(x = PC1, y = PC2, col = Color)) + geom_point(alpha = 0.3)</pre>
```



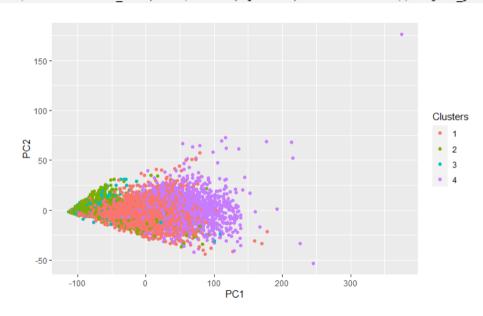
#### H3:

```
rotated_data$Clusters = as.factor(h3)
Plot and color by labels
ggplot(data = rotated_data, aes(x = PC1, y = PC2, col = Clusters)) + geom_point()
```



#### K-means:

```
rotated_data$Clusters = as.factor(fit$cluster)
Plot and color by labels
ggplot(data = rotated data, aes(x = PC1, y = PC2, col = Clusters)) + geom point()
```



## e. Consider the results of C and D and explain the differences between the clustering results in terms of how the algorithms work.

K-means uses a pre-specified K value, while HAC doesn't. We can see that the clustering method seems to be less random in the K-means method. Since HAC is arranged more like a 'tree', it seemed to struggling with clustering the data. K-means divided the data into non-overlapping clusters, which provided a more reasonable look at the dataset. Regardless, I still find PCA to be more sensible visually in this

case, even though I don't think it performed as well as k-means did in this case as it didn't handle the complexity of the dataset as well.

#### PROBLEM – 4

Back to the Starwars data from a previous assignment! Remember that the variable that lists the actual names and the variables that are actually lists will be a problem, so remove them (name, films, vehicles, starships). Make sure to double check the types of the variables, i.e., that they are numerical or factors as you expect.

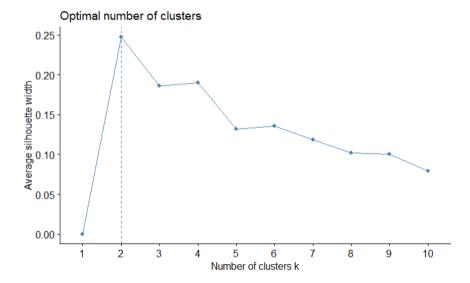
```
data("starwars")
#Copy Starwars
cpstarwars <- starwars
#Remove some columns
cpstarwars <- cpstarwars %>% select(-c("name", "vehicles", "starships", "films"))
#Remove NAs
cpstarwars <- na.omit(cpstarwars)
summary(cpstarwars)</pre>
```

```
height mass hair_color skin_color eye_color birth_year
Min.: 88 Min.: 20.00 Length:29 Length:29 Length:29 Min.: 8.00
1st Qu.:170 1st Qu.: 75.00 Class: character Class: character Class: character 1st Qu.: 31.00
Median: 180 Median: 79.00 Mode: character Mode: character Mode: character Median: 46.00
Mean: 178 Mean: 77.77
3rd Qu.:188 3rd Qu.: 83.00 3rd Qu.: 57.00
Max.: 228 Max.: 136.00 Max.: 238 Max.: 136.00
sex gender homeworld species
Length:29 Length:29 Length:29 Length:29
Class: character Class: character Class: character
Mode: character Mode: character Mode: character
Mode: character Mode: character Mode: character
```

a. Use hierarchical agglomerative clustering to cluster the Starwars data. This time we can leave the categorical variables in place, because we will use the gower metric from daisy in the cluster library to get the distances. Use average linkage. Determine the best number of clusters.

I had some trouble trying to run this without converting the categoricals to dummies, so I decided to proceed using the dummy variables.

```
library(cluster)
#Pass dataframe directly with mertic = gower
dist_mat <- daisy(dummies, metric = "gower")
#Center scale allows us to standardize the data
preproc <- predict(preproc, cpstarwars)
#Silhouette score comparison to find K
fviz_nbclust(predictors, FUN = hcut, method = "silhouette")</pre>
```



#### **Clusters:**

```
#Determine the assembly/agglomeration method and run hclust
hfit <- hclust(dist_mat, method = 'average')
#Build the new model
h2 <- cutree(hfit, k=2)
summary(h2)

Min. 1st Qu. Median Mean 3rd Qu. Max.</pre>
```

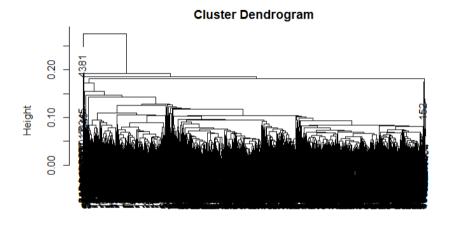
1 1

b. Produce the dendogram for (a). How might an anomaly show up in a dendogram? Do you see a Starwars character who does not seem to fit in easily? What is the advantage of considering anomalies this way as opposed to looking for unusual values relative to the mean and standard deviations, as we considered earlier in the course? Disadvantages?

1

1

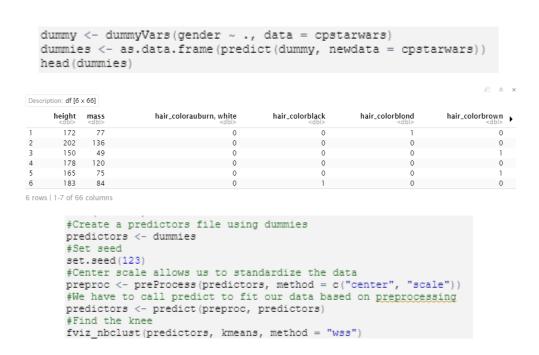
```
hfit <- hclust(dist_mat, method = 'average')
plot(hfit)</pre>
```

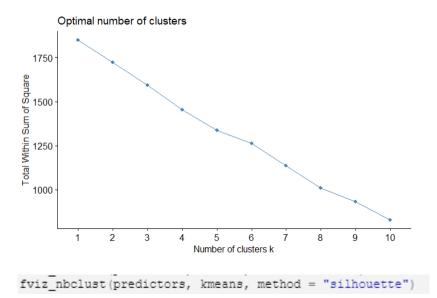


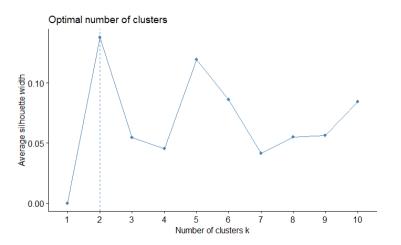
dist\_mat hclust (\*, "average")

Own branch with no relationship to other members of the dataset. Based on this dataset, 21, 9, and 18 may have anomalous features. The advantages of this dendogram is the ability to view anomalies directly without reviewing the actual dataset in a tabulation. However, we're only able to view the anomalies here based on height, which may be lacking in regards to other features like eye color (or) mass.

c. Use dummy variables to make this data fully numeric and then use k-means to cluster. Choose the best number of clusters.







#### K=2 will be the distance we are using here,

```
#Fit the data

fit <- kmeans (predictors, centers = 2, nstart = 25)

#Display the kmeans object information

fit

K-means clustering with 2 clusters of sizes 9, 20

Cluster means:
 height mass hair colorauburn, white hair colorblack hair colorblond hair colorbrown hair colorbrown, greys 1, 6126686 0,6114774 - 0.1356953 -0.532706 -0.2674319 -0.297273 -0.1356953 -0.13569
```

### d. Compare the HAC and k-means clusterings with a crosstabulation.

```
"``{r}

#Create a dataframe for results
result <- data.frame(Gender = cpstarwars$gender, HAC2 = h2, Kmeans = fit$cluster)
#Create a cross tab for HAC
result %>% group_by(HAC2) %>% select(HAC2, Gender) %>% table()
"``
```