

Homework Assignment - 05

Critical Thinking Group 5

Arindam Barman

Mohamed Elmoudni

Shazia Khan

Kishore Prasad

Contents

Overview	2
Objective	2
1 Data Exploration Analysis	2
1.1 Variable identification	2
1.2 Variable Relationships	3
1.3 Data summary analysis	4
1.4 Outliers Identification	7
2. Data Preparation	20
2.1 Missing Flags	20
2.2 Missing values treatment	21
2.3 Outliers treatment	21
2.4 Dummy Variables	22
2.5 Correlation for new variables	22
3. Build Models	23
3.1 Poisson models	25
3.2 Negative Binomial models	26
3.3 Linear Regression models	27
3.3.1 Linear Regression Model 5	27
3.3.1 Linear Regression Model 6	29

Overview

In this homework assignment, we will explore, analyze and model a data set containing information on approximately 12795 commercially available wines using 16 variables. The variables are mostly related to the chemical properties of the wine being sold. The response variable is the number of sample cases of wine that were purchased by wine distribution companies after sampling a wine. These cases would be used to provide tasting samples to restaurants and wine stores around the United States. The more sample cases purchased, the more likely is a wine to be sold at a high end restaurant. A large wine manufacturer is studying the data in order to predict the number of wine cases ordered based upon the wine characteristics. If the wine manufacturer can predict the number of cases, then that manufacturer will be able to adjust their wine offering to maximize sales.

Objective

Our objective is to build a count regression model to predict the number of cases of wine that will be sold given certain properties of the wine. Using the training data set, we will build at least two different Poisson regression models, at least two different negative binomial regression models, and at least two multiple linear regression models, using different variables (or the same variables with different transformations).

To attain our objective, we will be following the below best practice steps and guidelines:

- 1 -Data Exploration
- 2 -Data Preparation
- 3 -Build Models
- 4 -Select Models

1 Data Exploration Analysis

In section we will explore and gain some insights into the dataset by pursuing the below high level steps and inquiries:

- Variable identification
- Variable Relationships
- Data summary analysis
- Outliers and Missing Values Identification

1.1 Variable identification

First we look the variables' datatypes and their roles.

Variable	Datatype	Role
INDEX	int	none
TARGET	int	response
FixedAcidity	num	predictor
VolatileAcidity	num	predictor
CitricAcid	num	predictor
ResidualSugar	num	predictor

Variable	Datatype	Role
Chlorides	num	predictor
FreeSulfurDioxide	num	predictor
TotalSulfurDioxide	num	predictor
Density	num	predictor
pH	num	predictor
Sulphates	num	predictor
Alcohol	num	predictor
LabelAppeal	int	predictor
AcidIndex	int	predictor
STARS	int	predictor

From the Table 1 above, we see that that all variables are quantitative mainly of numeric and integer datatype. Also, we will ignore the INDEX variable as it is just a unique identifier for each row. However, we will use the TARTGET variable as response variable and the remaining variables as predictors.

1.2 Variable Relationships

Next let's display and examine the variable relationships as shown in table 2.

Table 2: Variable Description

VARIABLE	DEFINITION	THEORETICAL.EFFECT
INDEX	Identification Variable (do not use)	None
TARGET	Number of Cases Purchased	None
AcidIndex	Proprietary method of testing total acidity of wine by using a weighted average	
Alcohol	Alcohol Content	
Chlorides	Chloride content of wine	
CitricAcid	Citric Acid Content	
Density	Density of Wine	
FixedAcidity	Fixed Acidity of Wine	
FreeSulfurDioxide	Sulfur Dioxide content of wine	
LabelAppeal	Marketing Score indicating the appeal of label design for consumers. High numbers suggest customers like the label design. Negative numbers suggest customers don't like the design.	Many consumers purchase based on the visual appeal of the wine label design. Higher numbers suggest better sales.
ResidualSugar	Residual Sugar of wine	
STARS	Wine rating by a team of experts. 4 Stars = Excellent, 1 Star = Poor	A high number of stars suggests high sales
Sulphates	Sulfate conten of wine	
TotalSulfurDioxide	Total Sulfur Dioxide of Wine	
VolatileAcidity	Volatile Acid content of wine	
pH	pH of wine	

At first glance, we can easily deduce that that the FreeSulfurDioxide (Sulfur Dioxide content of wine) can be

derived from the TotalSulfurDioxide (Total Sulfur Dioxide of Wine). However, looking closer at the role of the sulfur dioxide SO_2 , as it is used as a preservative because of its anti-oxidative and anti-microbial properties in wine and also as a cleaning agent for barrels and winery facilities, we realize that when a winemaker says his/her wine has 100 ppm (part per million) of SO_2 , he/she is most probably referring to the total amount of SO_2 in his wine, and that means:

total SO_2 = free SO_2 + bound SO_2 .

free SO_2 : molecular SO_2 + bisulfites + sulfites

bound SO_2 : sulfites attached to either sugars, acetaldehyde or phenolic compounds

In this case the free SO_2 portion (not associated with wine molecules) is effectively the buffer against microbes and oxidation... Hence without knowing the bound SO_2 , we won't be able to derive FreeSulfurDioxide from TotalSulfurDioxide.

Also, looking briefly at the VolatileAcidity (Volatile Acid content of wine) and FixedAcidity (Fixed Acidity of Wine), we can easily deduce AcidIndex as the Acid index = Total acid (g/L) - pH. where Total acidity = Volatile Acid + Fixed Acidity. However, in our case the index is weighted average and we don't know the weighted average of either Volatile Acid or Fixed Acidity. Hence we will assume these variable do not have strict arithmetic relationships.

1.3 Data summary analysis

In this section, we will create summary data to better understand the initial relationship variables have with our dependent variable using correlation, central tendency, and dispersion As shown in table 3.

```
## 'data.frame': 12795 obs. of 15 variables:
## $ TARGET : int 3 3 5 3 4 0 0 4 3 6 ...
## $ FixedAcidity : num 3.2 4.5 7.1 5.7 8 11.3 7.7 6.5 14.8 5.5 ...
## $ VolatileAcidity : num 1.16 0.16 2.64 0.385 0.33 0.32 0.29 -1.22 0.27 -0.22 ...
## $ CitricAcid : num -0.98 -0.81 -0.88 0.04 -1.26 0.59 -0.4 0.34 1.05 0.39 ...
## $ ResidualSugar : num 54.2 26.1 14.8 18.8 9.4 ...
## $ Chlorides : num -0.567 -0.425 0.037 -0.425 NA 0.556 0.06 0.04 -0.007 -0.277 ...
## $ FreeSulfurDioxide : num NA 15 214 22 -167 -37 287 523 -213 62 ...
## $ TotalSulfurDioxide: num 268 -327 142 115 108 15 156 551 NA 180 ...
## $ Density : num 0.993 1.028 0.995 0.996 0.995 ...
## $ pH : num 3.33 3.38 3.12 2.24 3.12 3.2 3.49 3.2 4.93 3.09 ...
## $ Sulphates : num -0.59 0.7 0.48 1.83 1.77 1.29 1.21 NA 0.26 0.75 ...
## $ Alcohol : num 9.9 NA 22 6.2 13.7 15.4 10.3 11.6 15 12.6 ...
## $ LabelAppeal : int 0 -1 -1 -1 0 0 0 1 0 0 ...
## $ AcidIndex : int 8 7 8 6 9 11 8 7 6 8 ...
## $ STARS : int 2 3 3 1 2 NA NA 3 NA 4 ...
```

Table 3: Data Summary

	mean	sd	median	trimmed
TARGET	3.0290739	1.9263682	3.00000	3.0538244
FixedAcidity	7.0757171	6.3176435	6.90000	7.0736739
VolatileAcidity	0.3241039	0.7840142	0.28000	0.3243890
CitricAcid	0.3084127	0.8620798	0.31000	0.3102520
ResidualSugar	5.4187331	33.7493790	3.90000	5.5800410
Chlorides	0.0548225	0.3184673	0.04600	0.0540159
FreeSulfurDioxide	30.8455713	148.7145577	30.00000	30.9334877
TotalSulfurDioxide	120.7142326	231.9132105	123.00000	120.8895367

	mean	sd	median	trimmed
Density	0.9942027	0.0265376	0.99449	0.9942130
pH	3.2076282	0.6796871	3.20000	3.2055706
Sulphates	0.5271118	0.9321293	0.50000	0.5271453
Alcohol	10.4892363	3.7278190	10.40000	10.5018255
LabelAppeal	-0.0090660	0.8910892	0.00000	-0.0099639
AcidIndex	7.7727237	1.3239264	8.00000	7.6431572
STARS	2.0417550	0.9025400	2.00000	1.9711258

Below is the missing values and correlation table of the predictor variables to the response variables.

Table 4: Missing Data and Data Correlation

	Missing	Correlation
TARGET	0	1.0000000
FixedAcidity	0	-0.0490109
VolatileAcidity	0	-0.0887932
CitricAcid	0	0.0086846
ResidualSugar	616	0.0164913
Chlorides	638	-0.0382631
FreeSulfurDioxide	647	0.0438241
TotalSulfurDioxide	682	0.0514784
Density	0	-0.0355175
pH	395	-0.0094448
Sulphates	1210	-0.0388496
Alcohol	653	0.0620616
LabelAppeal	0	0.3565005
AcidIndex	0	-0.2460494
STARS	3359	0.5587938

Missing Values and Correlation Interpretation

From tables 3 and 4 above, we observe the followings:

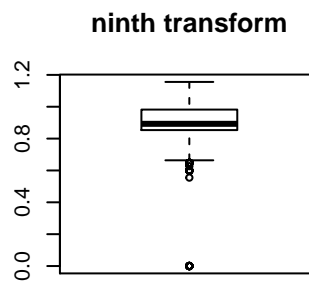
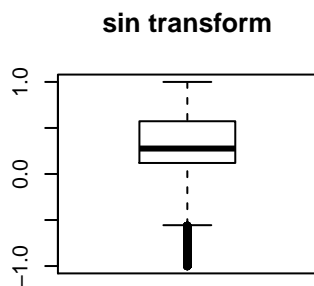
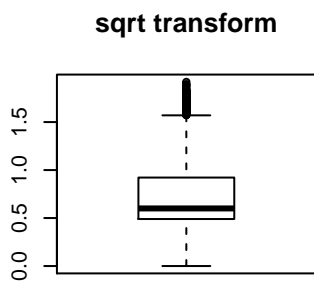
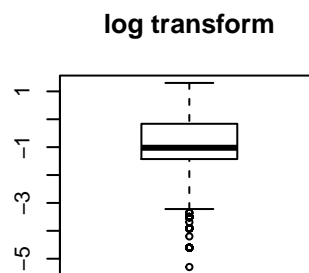
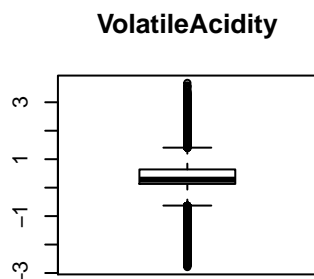
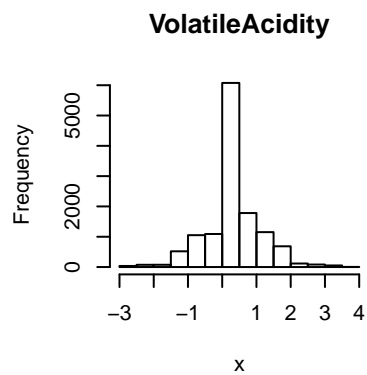
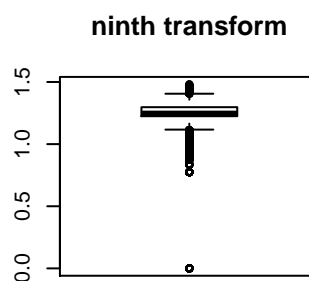
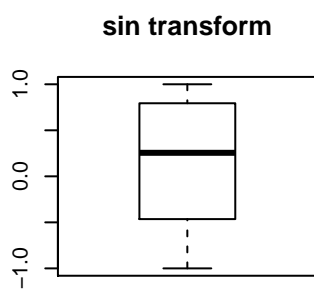
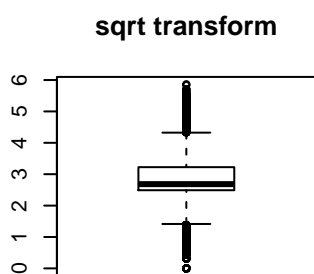
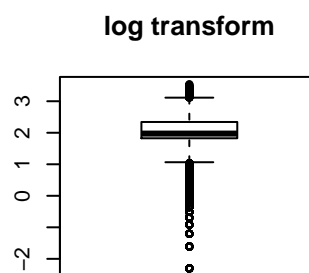
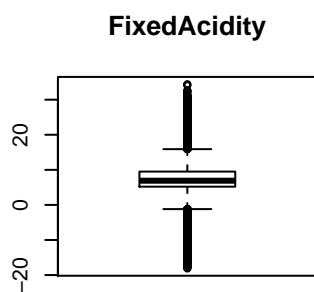
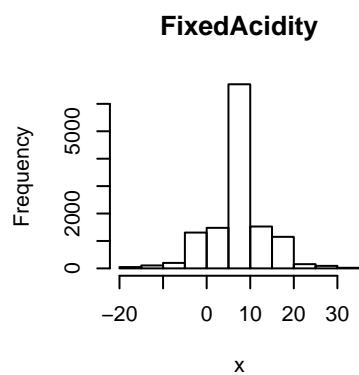
- Variable ResidualSugar has 616 and 0.0164913 correlation. Given the low correlation we will try some imputation techniques to handle the missing the values and replace missing values with their respective value.
- variable Chlorides 638 -0.0382631 correlation. . Given the low negative correlation we will try we would replace missing values with their respective value
- Variable FreeSulfurDioxide 647 0.0438241. Given the low correlation we will impute the missing values with their respective value
- Variable TotalSulfurDioxide has 682 missing values with 0.0514784 correlation. Given the low correlation we will impute the missing values with their respective value.
- Variable Alcohol has 653 missing values with 0.0620616 correlation. Given the low correlation we will impute the missing values with their respective value.

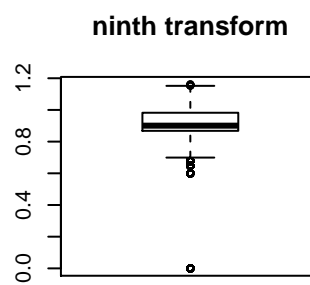
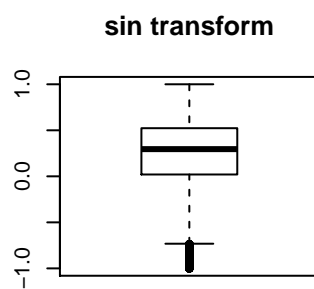
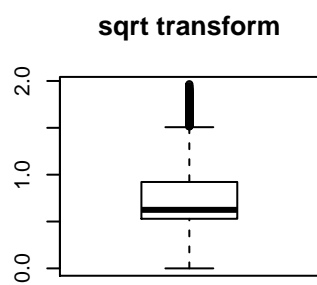
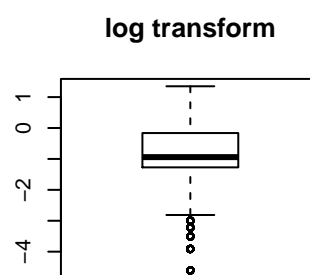
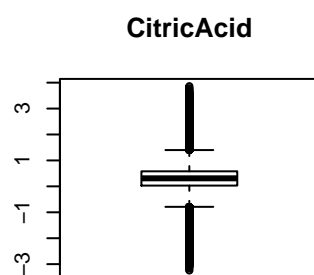
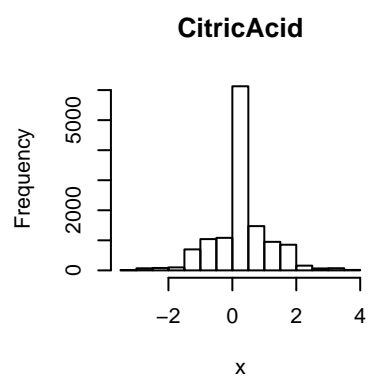
Please note that ResidualSugar, Chlorides, FreeSulfurDioxide, Alcohol, and TotalSulfurDioxide variables have similar number of missing values. They are chemically related. However, we don't think they are arithmetically related.

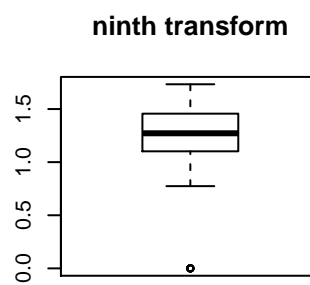
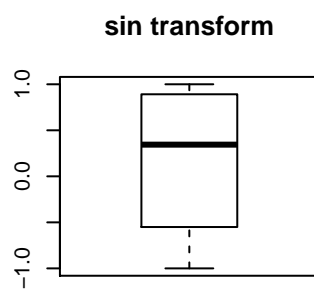
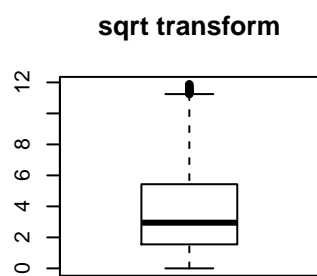
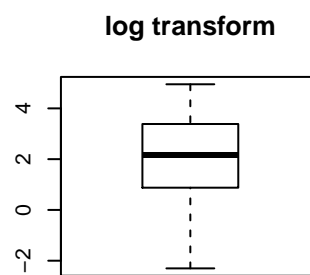
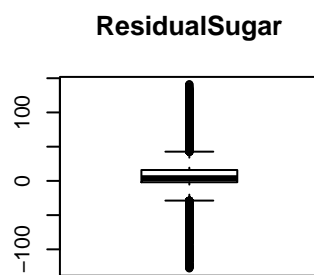
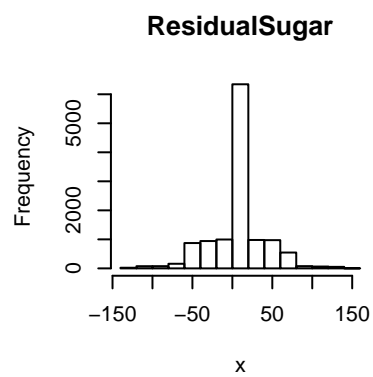
- In addition, variable pH has 395 missing values with negative correlation of -0.0094448. Again we may just ignore these missing values especially that it has very low negative correlation to the target variable.
- Variable Sulphates has much higher missing values of 1210 with low negative correlation of -0.0388496. We will be imputing this values with their respective value
- Now, variable STARS has the highest missing values of 3359 and highest correlation of 0.5587938. This is very important variable and it drives sales and consequently heavily impacts our response variable. We have to be careful in fixing the missing values as this variable STARS is rating score variable with 1 being the lowest and 4 the highest

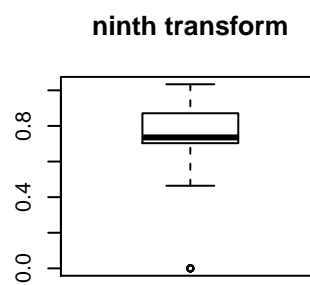
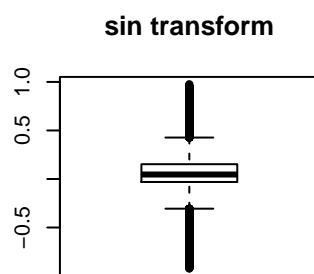
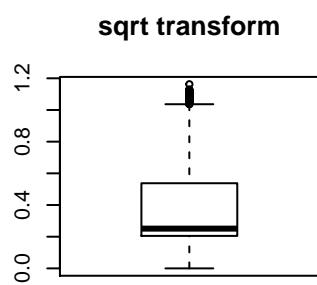
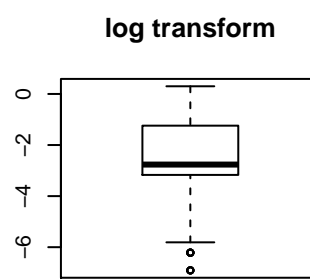
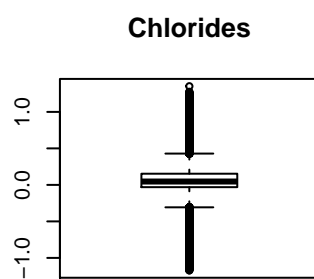
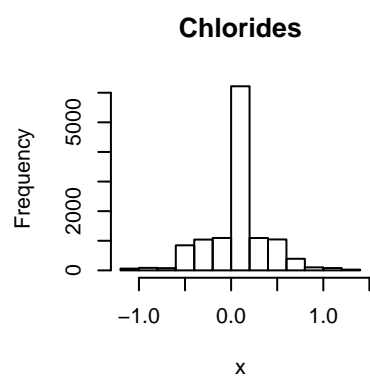
1.4 Outliers Identification

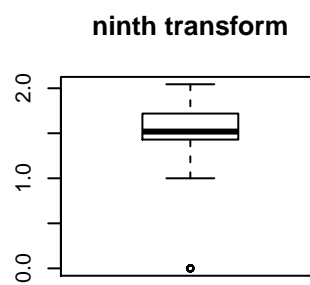
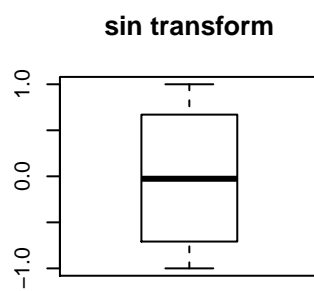
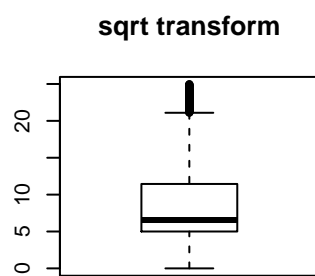
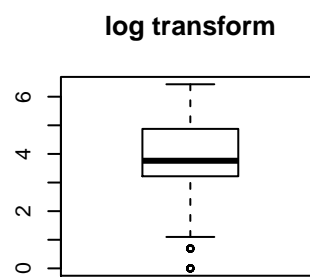
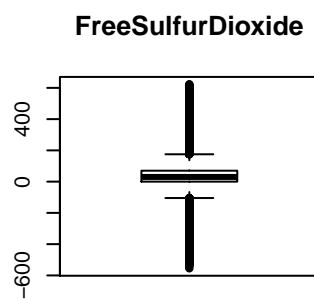
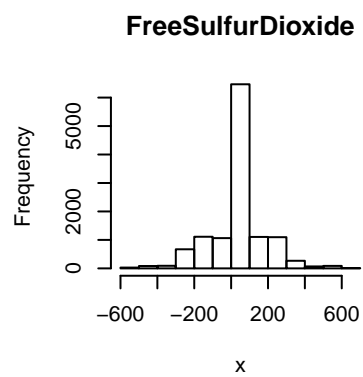
In this section we look at boxplots to determine the outliers in variables and decide on whether to act on the outliers. Lets do some univariate analysis. We will look at the Histogram and Boxplot for each variable to detect outliers if any and treat it accordingly.

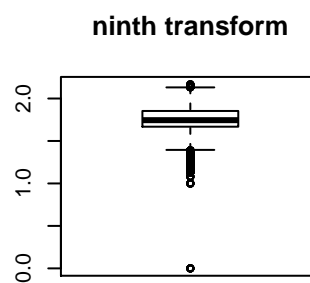
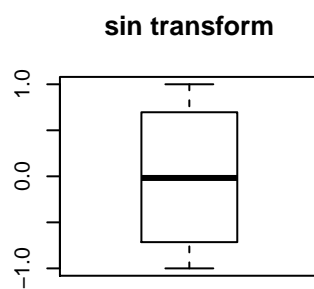
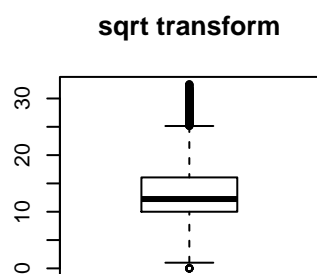
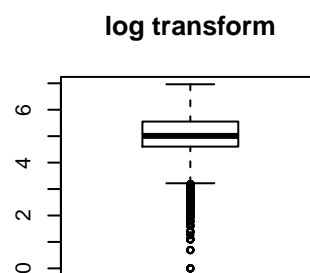
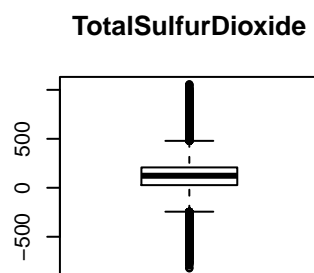
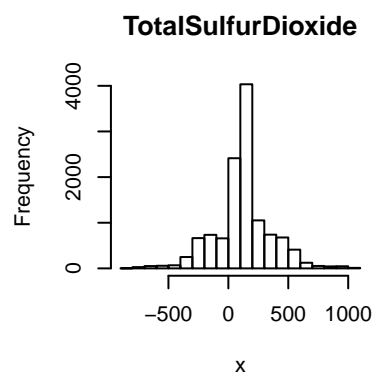


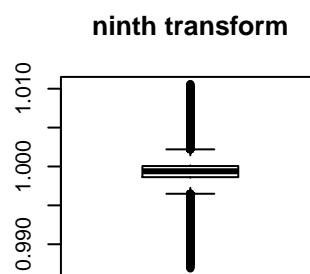
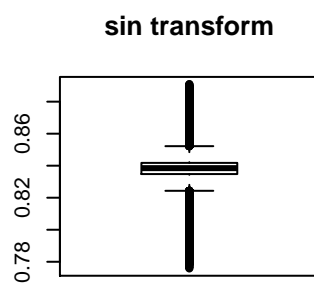
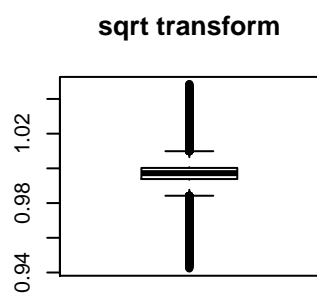
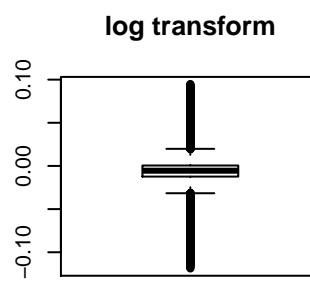
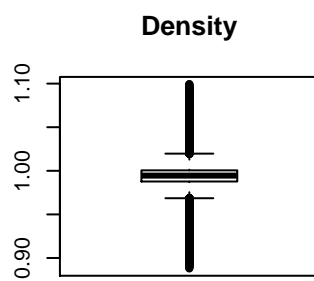
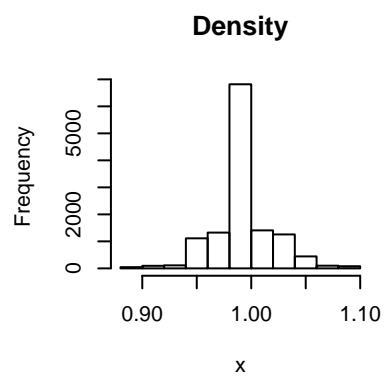


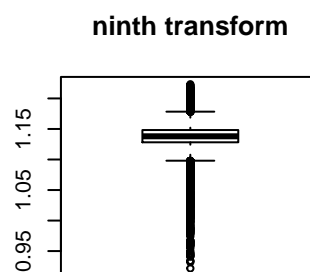
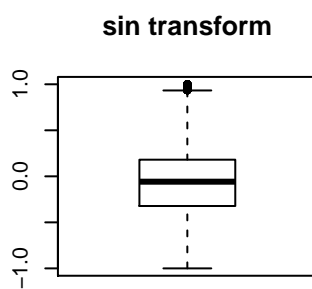
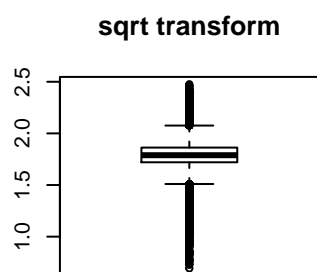
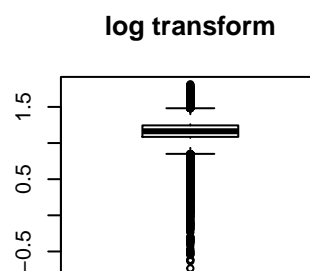
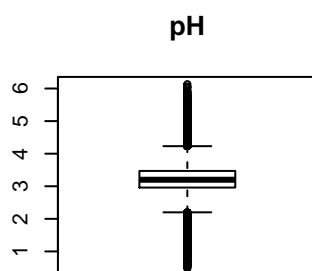
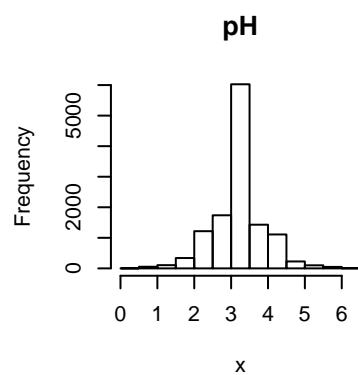


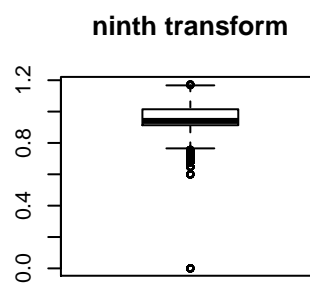
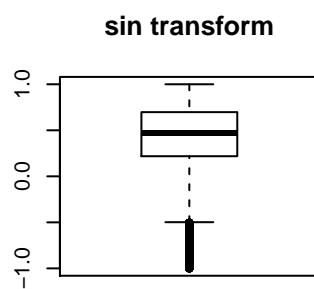
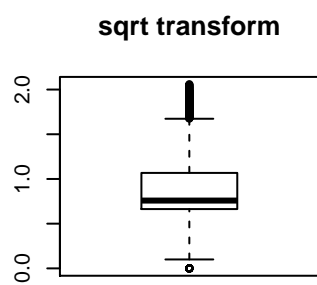
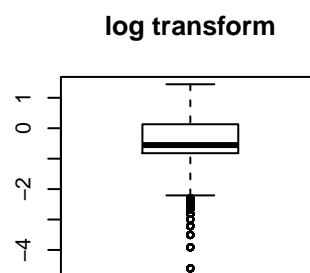
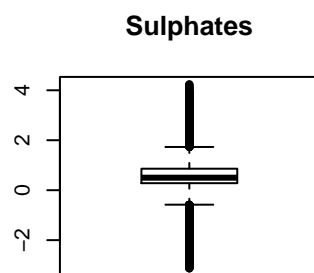
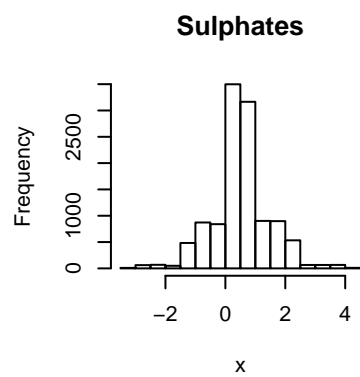


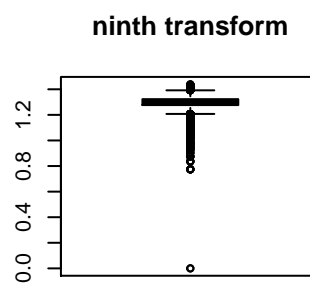
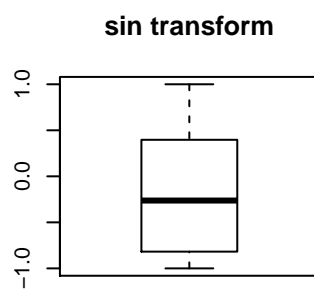
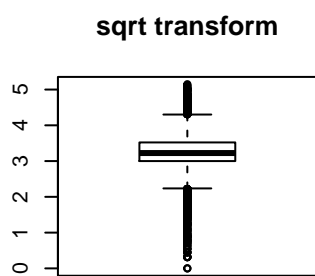
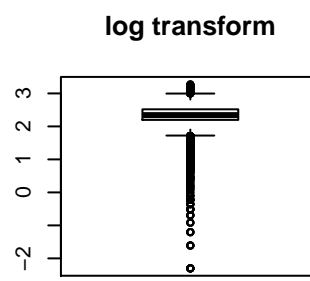
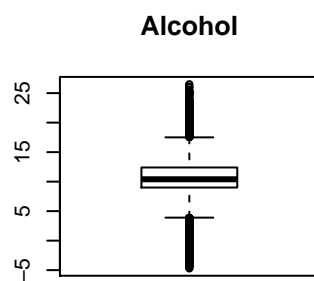
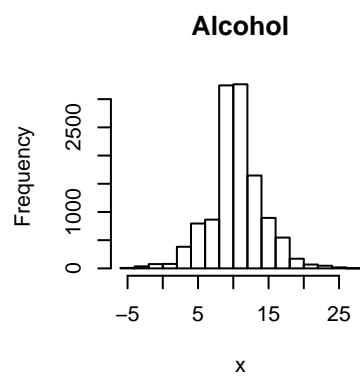


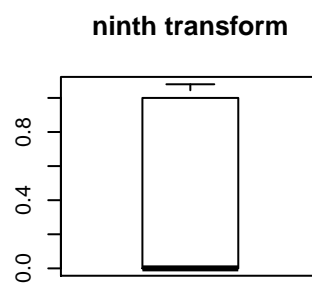
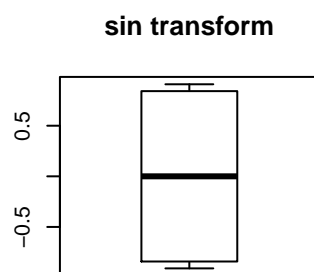
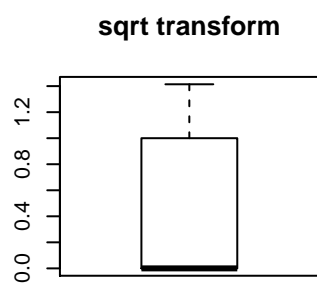
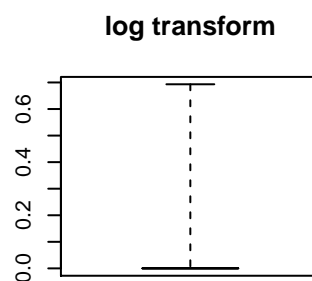
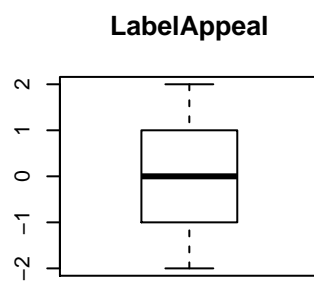
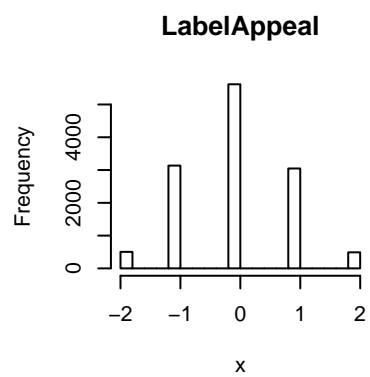


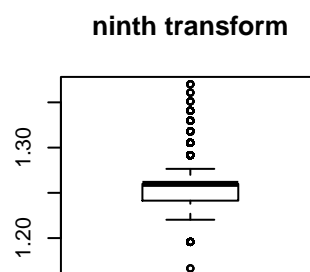
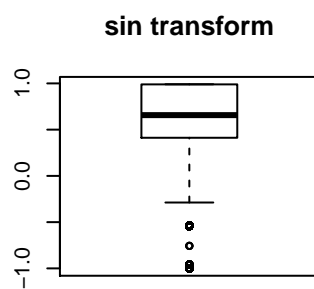
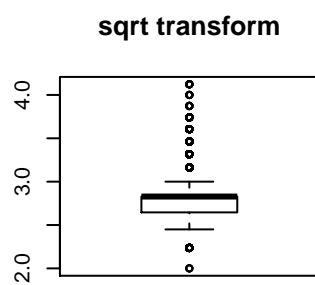
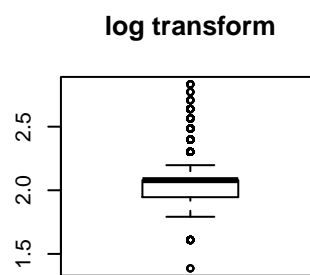
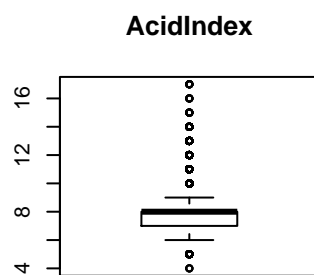
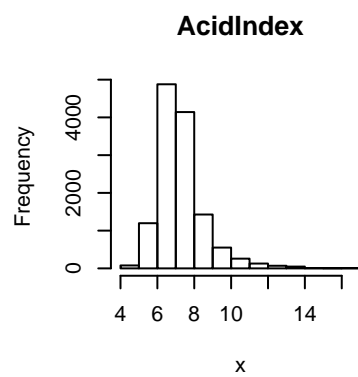


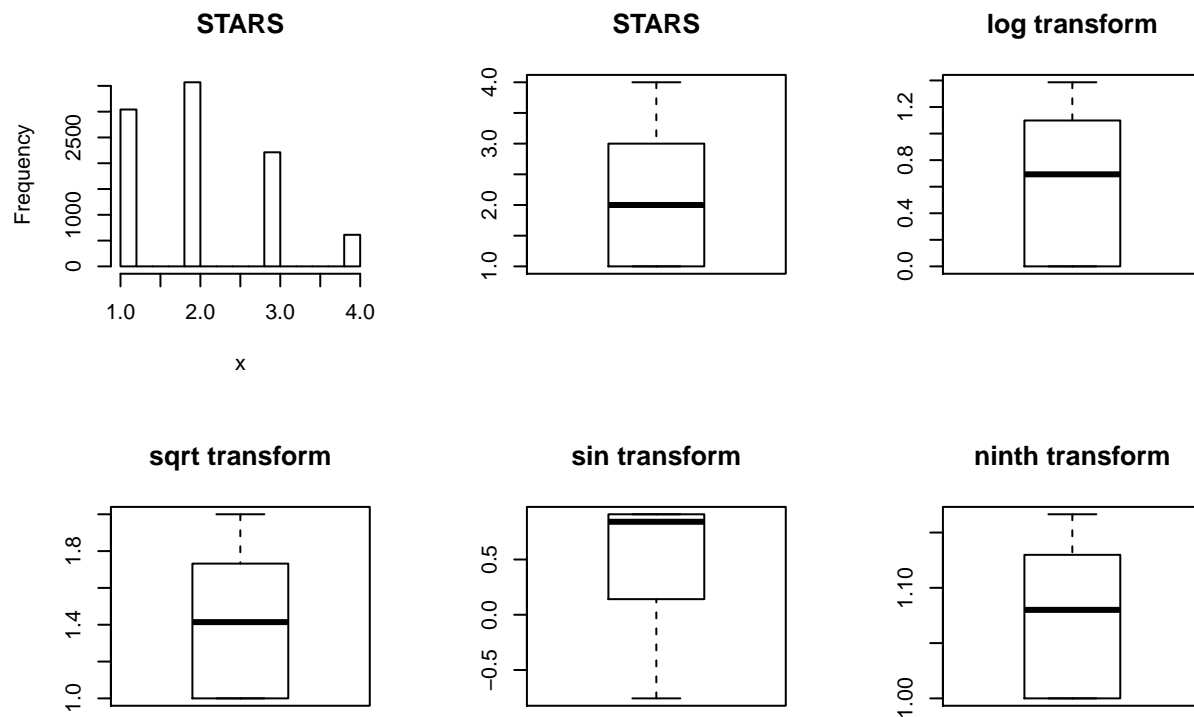












2. Data Preparation

Now that we have completed the preliminary analysis, we will be cleaning and consolidating data into one dataset for use in analysis and modeling. We will be puring the belwo steps as guidelines:

- Missing Flags
- Missing values treatment
- Outliers treatment
- Dummy Variables

2.1 Missing Flags

We create flag variables to indicate whether some of the fields are missing any values. If the value is missing, we code it with 1 and if the value is present we code it with 0. The following are the variables that are created:

- ResidualSugar_MISS
- Chlorides_MISS
- FreeSulfurDioxide_MISS
- TotalSulfurDioxide_MISS
- pH_MISS
- Sulphates_MISS
- Alcohol_MISS
- STARS_MISS

2.2 Missing values treatment

Next we impute missing values. We can go ahead and use the mean as impute values. We will replace the missing values in the original variables. However, for STARS, we will code the missing value as a '0' instead of a mean. The following are the variables that are impacted:

- ResidualSugar
- Chlorides
- FreeSulfurDioxide
- TotalSulfurDioxide
- pH
- Sulphates
- Alcohol
- STARS

2.3 Outliers treatment

For outliers, we will use the capping method. In this method, we will replace all outliers that lie outside the 1.5 times of IQR limits. We will cap it by replacing those observations less than the lower limit with the value of 5th %ile and those that lie above the upper limit with the value of 95th %ile.

Accordingly we create the following new variables while retaining the original variables.

- FixedAcidity_CAP
- VolatileAcidity_CAP
- CitricAcid_CAP
- ResidualSugar_CAP
- Chlorides_CAP
- FreeSulfurDioxide_CAP
- TotalSulfurDioxide_CAP
- Density_CAP
- pH_CAP
- Sulphates_CAP
- Alcohol_CAP
- AcidIndex_CAP

2.4 Dummy Variables

Finally, we will also create dummy variables for the following variables:

- LabelAppeal : For this variable, we create a dummy variable to indicate if the value is Zero / Positive or Negative.
- STARS - We create a Dummy Variable for each of the star ratings - 1,2,3,4. The value is 1 in the respective variable based on the STARS value. A Zero value in all of the STARS dummy vars indicate that the value was missing in the original variable.

2.5 Correlation for new variables

Lets see how the new variables stack up against the TARGET.

Table 5: Correlation between TARGET and predictor variables

	Correlation
STARS_3	0.3597277
STARS_4	0.2783731
STARS_2	0.2484240
Alcohol_CAP	0.0634633
TotalSulfurDioxide_CAP	0.0503492
FreeSulfurDioxide_CAP	0.0417585
LabelAppeal_Positive	0.0206261
ResidualSugar_CAP	0.0204409
CitricAcid_CAP	0.0120351
ResidualSugar_MISS	0.0111995
TotalSulfurDioxide_MISS	0.0061720
Chlorides_MISS	0.0026937
Alcohol_MISS	0.0014776
FreeSulfurDioxide_MISS	-0.0001501
pH_MISS	-0.0099654
pH_CAP	-0.0102565
Sulphates_MISS	-0.0125039
Chlorides_CAP	-0.0304686
Density_CAP	-0.0315375
Sulphates_CAP	-0.0359312
FixedAcidity_CAP	-0.0510757
VolatileAcidity_CAP	-0.0891214
STARS_1	-0.1300422
AcidIndex_CAP	-0.2353997
STARS_MISS	-0.5715792

From the above Correlations, we can make the following observations:

- The following variables have a positive correlation with TARGET: STARS_3, STARS_4, STARS_2, Alcohol_CAP, TotalSulfurDioxide_CAP, FreeSulfurDioxide_CAP, LabelAppeal_Positive, ResidualSugar_CAP, CitricAcid_CAP, ResidualSugar_MISS, TotalSulfurDioxide_MISS, Chlorides_MISS, Alcohol_MISS.
- The following variables have a negative correlation with TARGET: FreeSulfurDioxide_MISS, pH_MISS,

pH_CAP, Sulphates_MISS, Chlorides_CAP, Density_CAP, Sulphates_CAP, FixedAcidity_CAP, VolatileAcidity_CAP, STARS_1, AcidIndex_CAP, STARS_MISS.

- Not all variable have a strong correlation in either direction. However, the following stand out for having a stronger correlation: STARS_MISS, STARS_3, STARS_4, STARS_2, AcidIndex_CAP, STARS_1, VolatileAcidity_CAP, Alcohol_CAP, FixedAcidity_CAP, TotalSulfurDioxide_CAP.

3. Build Models

Since we are dealing with count variables, our modeling technique will mainly focus on using variation of the Generalized Linear Model (GLM) family functions. We will start with the classical Poisson regression; then we will enhance it using model Negative binominal model.

In addition, we will also create models using linear regression.

Using original and transformed datasets, we will build at least ten models as follow:

- Two Poisson models
- Two Quasi-Poisson models
- Two Negative binomial models
- Two Zero-inflated models
- Two Linear regression models

Below is a summary table showing models and their respective variables.

Table 6: Models and their Respective Variables

Variable	Model.1	Model.2	Comments
TARGET	Y	Y	The TARGET variable
FixedAcidity	Y		Imputed with Mean
VolatileAcidity	Y		Imputed with Mean
CitricAcid	Y		Imputed with Mean
ResidualSugar	Y		Imputed with Mean
Chlorides	Y		Imputed with Mean
FreeSulfurDioxide	Y		Imputed with Mean
TotalSulfurDioxide	Y		Imputed with Mean
Density	Y		Imputed with Mean
pH	Y		Imputed with Mean
Sulphates	Y		Imputed with Mean
Alcohol	Y		Imputed with Mean
LabelAppeal	Y		Original Variable
AcidIndex	Y		Imputed with Mean
STARS	Y		Original Variable
ResidualSugar_MISS		Y	Missing Flag
Chlorides_MISS		Y	Missing Flag
FreeSulfurDioxide_MISS		Y	Missing Flag
TotalSulfurDioxide_MISS		Y	Missing Flag
pH_MISS		Y	Missing Flag
Sulphates_MISS		Y	Missing Flag
Alcohol_MISS		Y	Missing Flag
STARS_MISS		Y	Missing Flag
FixedAcidity_CAP		Y	Imputed with Mean and Outliers capped
VolatileAcidity_CAP		Y	Imputed with Mean and Outliers capped
CitricAcid_CAP		Y	Imputed with Mean and Outliers capped
ResidualSugar_CAP		Y	Imputed with Mean and Outliers capped

Variable	Model.1	Model.2	Comments
Chlorides_CAP		Y	Imputed with Mean and Outliers capped
FreeSulfurDioxide_CAP		Y	Imputed with Mean and Outliers capped
TotalSulfurDioxide_CAP		Y	Imputed with Mean and Outliers capped
Density_CAP		Y	Imputed with Mean and Outliers capped
pH_CAP		Y	Imputed with Mean and Outliers capped
Sulphates_CAP		Y	Imputed with Mean and Outliers capped
Alcohol_CAP		Y	Imputed with Mean and Outliers capped
AcidIndex_CAP		Y	Imputed with Mean and Outliers capped
LabelAppeal_Positive		Y	Positive or Negative Dummy Variable
STARS_1		Y	Dummy Variable
STARS_2		Y	Dummy Variable
STARS_3		Y	Dummy Variable
STARS_4		Y	Dummy Variable

3.1 Poisson models

Our first attempt to capture the relationship between the wine chemical properties and number of cases of the wine being sold in a parametric regression model, we fit the basic Poisson regression model

3.1.1 Poisson Model 1

We will explore the Poisson regression model Using original data with replacing all missing data with the means.

```
#winedata <- read.csv("C:/CUNY/Courses/IS621/Assignment602/Assignment05/wine-training-data.csv")
#dim(winedata)
#winedata <- select(winedata, ~(INDEX))
#str(winedata)
## all zeros
# winedata[is.na(winedata)] <- 0

poismod1 <- glm(TARGET ~ ., data=winedata_orig, family=poisson)
summary(poismod1)
```

```
##
## Call:
## glm(formula = TARGET ~ ., family = poisson, data = winedata_orig)
##
## Deviance Residuals:
##      Min       1Q   Median       3Q      Max
## -2.9733  -0.7200   0.0694   0.5785   3.2315
##
## Coefficients:
##              Estimate Std. Error z value Pr(>|z|)
## (Intercept)    1.526e+00  1.955e-01   7.807 5.87e-15 ***
## FixedAcidity   -3.045e-04  8.205e-04  -0.371 0.710502
## VolatileAcidity -3.343e-02  6.516e-03  -5.131 2.88e-07 ***
## CitricAcid      7.773e-03  5.892e-03   1.319 0.187124
## ResidualSugar    5.676e-05  1.546e-04   0.367 0.713588
## Chlorides       -4.141e-02  1.645e-02  -2.518 0.011816 *
## FreeSulfurDioxide 1.254e-04  3.512e-05   3.571 0.000356 ***
## TotalSulfurDioxide 8.296e-05  2.275e-05   3.647 0.000266 ***
## Density        -2.823e-01  1.920e-01  -1.471 0.141348
## pH              -1.572e-02  7.638e-03  -2.058 0.039554 *
## Sulphates       -1.267e-02  5.749e-03  -2.205 0.027480 *
## Alcohol         2.201e-03  1.410e-03   1.561 0.118446
## LabelAppeal     1.332e-01  6.063e-03  21.968 < 2e-16 ***
## AcidIndex       -8.705e-02  4.548e-03 -19.139 < 2e-16 ***
## STARS           3.113e-01  4.531e-03  68.700 < 2e-16 ***
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## (Dispersion parameter for poisson family taken to be 1)
##
##      Null deviance: 22861  on 12794  degrees of freedom
## Residual deviance: 14728  on 12780  degrees of freedom
```

```
## AIC: 46700
##
## Number of Fisher Scoring iterations: 5
```

*** Interpretation Poisson Model 1***

3.1.2 Poisson Model 2

In this model we will be using the basic Poisson regression model; however using transformed data.

```
# transformed data. Poisson Model 2
```

*** Interpretation Poisson Model 2***

3.2 Negative Binomial models

A more formal way to accommodate over-dispersion in a count data regression model is to use a negative binomial model. Hence we will explore the negative binomial model both in original data as well as transformed data.

3.2.1 Negative Binomial model 3

We will explore the Negative Binomial model Using original data with replacing all missing data with the means.

```
nbmod3 = glm.nb(TARGET ~ ., data = winedata_orig)
```

```
## Warning in theta.ml(Y, mu, sum(w), w, limit = control$maxit, trace =
## control$trace > : iteration limit reached
```

```
## Warning in theta.ml(Y, mu, sum(w), w, limit = control$maxit, trace =
## control$trace > : iteration limit reached
```

```
summary(nbmod3)
```

```
##
## Call:
## glm.nb(formula = TARGET ~ ., data = winedata_orig, init.theta = 48974.65455,
##       link = log)
##
## Deviance Residuals:
##      Min       1Q   Median       3Q      Max
## -2.9732  -0.7200   0.0694   0.5785   3.2314
##
## Coefficients:
##              Estimate Std. Error z value Pr(>|z|)
## (Intercept)    1.526e+00  1.955e-01   7.806 5.88e-15 ***
## FixedAcidity   -3.046e-04  8.205e-04  -0.371 0.710504
## VolatileAcidity -3.343e-02  6.516e-03  -5.131 2.89e-07 ***
```

```

## CitricAcid          7.773e-03  5.892e-03   1.319 0.187136
## ResidualSugar       5.676e-05  1.546e-04   0.367 0.713573
## Chlorides           -4.142e-02  1.645e-02  -2.518 0.011817 *
## FreeSulfurDioxide   1.254e-04  3.512e-05   3.571 0.000356 ***
## TotalSulfurDioxide  8.296e-05  2.275e-05   3.647 0.000266 ***
## Density             -2.824e-01  1.920e-01  -1.471 0.141356
## pH                  -1.572e-02  7.638e-03  -2.058 0.039552 *
## Sulphates           -1.267e-02  5.749e-03  -2.205 0.027480 *
## Alcohol             2.201e-03  1.410e-03   1.561 0.118467
## LabelAppeal         1.332e-01  6.064e-03  21.967 < 2e-16 ***
## AcidIndex           -8.705e-02  4.548e-03 -19.139 < 2e-16 ***
## STARS               3.113e-01  4.531e-03  68.698 < 2e-16 ***
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## (Dispersion parameter for Negative Binomial(48974.65) family taken to be 1)
##
##      Null deviance: 22860  on 12794  degrees of freedom
## Residual deviance: 14728  on 12780  degrees of freedom
## AIC: 46703
##
## Number of Fisher Scoring iterations: 1
##
##
##              Theta: 48975
##              Std. Err.: 50715
## Warning while fitting theta: iteration limit reached
##
## 2 x log-likelihood: -46670.5

```

*** Interpretation Negative Binomial Model 3***

3.2.1 Negative Binomial model 4

In this model we will be using the basic Negative Binomial model; however using transformed data.

```
#transformed data. Negative Binomial model 4
```

*** Interpretation Negative Binomial Model 4***

3.3 Linear Regression models

Although it is highly recommended for continuous variables instead of count variables, we will also create two linear regression models.

3.3.1 Linear Regression Model 5

We will explore the Linear models Using original data with replacing all missing data with the means.

```
##
## Call:
```

```
## lm(formula = TARGET ~ ., data = winedata_orig)
##
## Residuals:
##      Min       1Q   Median       3Q      Max
## -4.5476 -0.9475  0.0669  0.9047  5.9903
##
## Coefficients:
##              Estimate Std. Error t value Pr(>|t|)
## (Intercept)    3.986e+00  4.487e-01   8.883  < 2e-16 ***
## FixedAcidity    1.608e-06  1.884e-03   0.001  0.999319
## VolatileAcidity -9.923e-02  1.498e-02  -6.625  3.61e-11 ***
## CitricAcid      2.085e-02  1.362e-02   1.531  0.125804
## ResidualSugar   2.012e-04  3.559e-04   0.565  0.571860
## Chlorides      -1.243e-01  3.777e-02  -3.290  0.001003 **
## FreeSulfurDioxide 3.153e-04  8.093e-05   3.897  9.80e-05 ***
## TotalSulfurDioxide 2.264e-04  5.201e-05   4.353  1.35e-05 ***
## Density        -8.012e-01  4.419e-01  -1.813  0.069829 .
## pH             -3.453e-02  1.754e-02  -1.969  0.049012 *
## Sulphates      -3.271e-02  1.322e-02  -2.475  0.013352 *
## Alcohol         1.094e-02  3.234e-03   3.384  0.000717 ***
## LabelAppeal     4.326e-01  1.367e-02  31.654  < 2e-16 ***
## AcidIndex      -2.084e-01  9.212e-03 -22.619  < 2e-16 ***
## STARS           9.767e-01  1.045e-02  93.433  < 2e-16 ***
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 1.324 on 12780 degrees of freedom
## Multiple R-squared:  0.528, Adjusted R-squared:  0.5275
## F-statistic: 1021 on 14 and 12780 DF, p-value: < 2.2e-16
```

*** Interpretation of Linear Model 5***

Based on the summary for Linear Model 5, below are the characteristics :

- The Residual standard error is 1.3242
- Multiple R-squared: 0.528
- Adjusted R-squared: 0.5275
- F-statistic: 1021 on 14 and 12780 DF
- p-value: < 2.2e-16

Based on the available coefficients, we can make the following observations:

- Positive Impact - The following variables have a positive impact on TARGET, meaning an increase in the values of these variables leads to an increase in the number of cases sold: STARS, LabelAppeal, Alcohol, TotalSulfurDioxide, FreeSulfurDioxide, ResidualSugar, CitricAcid, FixedAcidity
- Negative Impact - The following variables have a negative impact on TARGET, meaning an increase in the values of these variables leads to a decrease in the number of cases sold: AcidIndex, Sulphates, pH, Density, Chlorides, VolatileAcidity
- The following variables have a 'significant' impact. These are the more important predictors for TARGET: STARS, AcidIndex, LabelAppeal, Alcohol, Sulphates, pH, TotalSulfurDioxide, FreeSulfurDioxide, Chlorides, VolatileAcidity
- Finally, the Linear Model equation is given by the following:

$3.9861 + 2e-06 * \text{FixedAcidity} - 0.099232 * \text{VolatileAcidity} + 0.020854 * \text{CitricAcid} + 0.000201 * \text{ResidualSugar}$
 $- 0.124266 * \text{Chlorides} + 0.000315 * \text{FreeSulfurDioxide} + 0.000226 * \text{TotalSulfurDioxide} - 0.801199 * \text{Density}$
 $- 0.034527 * \text{pH} - 0.032707 * \text{Sulphates} + 0.010942 * \text{Alcohol} + 0.432607 * \text{LabelAppeal} - 0.208371 * \text{AcidIndex}$
 $+ 0.976721 * \text{STARS}$

3.3.1 Linear Regression Model 6

In this model we will be using the Linear Regression model; however using transformed data.

```
##
## Call:
## lm(formula = TARGET ~ ., data = winedata_trans)
##
## Residuals:
##      Min       1Q   Median       3Q      Max
## -4.1375 -0.9450  0.0246  0.9372  6.7449
##
## Coefficients: (1 not defined because of singularities)
##              Estimate Std. Error t value Pr(>|t|)
## (Intercept)      7.938e+00  4.779e-01  16.611 < 2e-16 ***
## ResidualSugar_MISS  6.293e-02  5.649e-02   1.114 0.265301
## Chlorides_MISS     6.208e-03  5.560e-02   0.112 0.911102
## FreeSulfurDioxide_MISS  6.444e-02  5.624e-02   1.146 0.251917
## TotalSulfurDioxide_MISS  4.998e-02  5.383e-02   0.929 0.353164
## pH_MISS           -8.562e-02  6.990e-02  -1.225 0.220634
## Sulphates_MISS    -2.325e-02  4.132e-02  -0.563 0.573704
## Alcohol_MISS       6.140e-02  5.494e-02   1.118 0.263709
## STARS_MISS        -4.092e+00  6.051e-02 -67.620 < 2e-16 ***
## FixedAcidity_CAP   -1.190e-03  2.175e-03  -0.547 0.584328
## VolatileAcidity_CAP -1.065e-01  1.720e-02  -6.193 6.07e-10 ***
## CitricAcid_CAP     2.220e-02  1.554e-02   1.429 0.153062
## ResidualSugar_CAP   3.782e-04  3.646e-04   1.038 0.299517
## Chlorides_CAP      -7.754e-02  3.840e-02  -2.020 0.043447 *
## FreeSulfurDioxide_CAP  4.803e-04  1.261e-04   3.809 0.000140 ***
## TotalSulfurDioxide_CAP  2.303e-04  6.162e-05   3.737 0.000187 ***
## Density_CAP       -9.171e-01  4.642e-01  -1.976 0.048215 *
## pH_CAP            -3.814e-02  2.061e-02  -1.850 0.064286 .
## Sulphates_CAP     -3.372e-02  1.406e-02  -2.397 0.016538 *
## Alcohol_CAP        1.311e-02  3.894e-03   3.367 0.000761 ***
## AcidIndex_CAP     -2.108e-01  1.177e-02 -17.916 < 2e-16 ***
## LabelAppeal_Positive -7.695e-02  4.394e-02  -1.751 0.079931 .
## STARS_1           -2.770e+00  6.074e-02 -45.611 < 2e-16 ***
## STARS_2           -1.586e+00  5.992e-02 -26.462 < 2e-16 ***
## STARS_3           -8.683e-01  6.248e-02 -13.898 < 2e-16 ***
## STARS_4              NA         NA      NA      NA
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 1.367 on 12770 degrees of freedom
## Multiple R-squared:  0.4976, Adjusted R-squared:  0.4967
## F-statistic:  527 on 24 and 12770 DF,  p-value: < 2.2e-16
```

*** Interpretation of Linear Model 6***

Based on the summary for Linear Model 6, below are the characteristics :

- The Residual standard error is 1.3667
- Multiple R-squared: 0.4976
- Adjusted R-squared: 0.4967
- F-statistic: 527 on 24 and 12770 DF
- p-value: < 2.2e-16

Based on the available coefficients, we can make the following observations:

- Positive Impact - The following variables have a positive impact on TARGET, meaning an increase in the values of these variables leads to an increase in the number of cases sold: Alcohol_CAP, TotalSulfurDioxide_CAP, FreeSulfurDioxide_CAP, ResidualSugar_CAP, CitricAcid_CAP, Alcohol_MISS, TotalSulfurDioxide_MISS, FreeSulfurDioxide_MISS, Chlorides_MISS, ResidualSugar_MISS
- Negative Impact - The following variables have a negative impact on TARGET, meaning an increase in the values of these variables leads to a decrease in the number of cases sold: STARS_3, STARS_2, STARS_1, LabelAppeal_Positive, AcidIndex_CAP, Sulphates_CAP, pH_CAP, Density_CAP, Chlorides_CAP, VolatileAcidity_CAP, FixedAcidity_CAP, STARS_MISS, Sulphates_MISS, pH_MISS
- The following variables have a 'significant' impact. These are the more important predictors for TARGET: STARS_3, STARS_2, STARS_1, AcidIndex_CAP, Alcohol_CAP, Sulphates_CAP, Density_CAP, TotalSulfurDioxide_CAP, FreeSulfurDioxide_CAP, Chlorides_CAP, VolatileAcidity_CAP, STARS_MISS
- Finally, the Linear Model equation is given by the following:

$$\begin{aligned} &7.938 + 0.062931 * \text{ResidualSugar_MISS} + 0.006208 * \text{Chlorides_MISS} + 0.064437 * \text{FreeSulfurDioxide_MISS} \\ &+ 0.049984 * \text{TotalSulfurDioxide_MISS} - 0.085625 * \text{pH_MISS} - 0.023249 * \text{Sulphates_MISS} + 0.061402 * \\ &\text{Alcohol_MISS} - 4.092034 * \text{STARS_MISS} - 0.00119 * \text{FixedAcidity_CAP} - 0.106543 * \text{VolatileAcidity_CAP} \\ &+ 0.022202 * \text{CitricAcid_CAP} + 0.000378 * \text{ResidualSugar_CAP} - 0.077542 * \text{Chlorides_CAP} + 0.00048 \\ &* \text{FreeSulfurDioxide_CAP} + 0.00023 * \text{TotalSulfurDioxide_CAP} - 0.917053 * \text{Density_CAP} - 0.038139 * \\ &\text{pH_CAP} - 0.033715 * \text{Sulphates_CAP} + 0.013111 * \text{Alcohol_CAP} - 0.210836 * \text{AcidIndex_CAP} - 0.07695 \\ &* \text{LabelAppeal_Positive} - 2.770326 * \text{STARS_1} - 1.585505 * \text{STARS_2} - 0.868345 * \text{STARS_3} \end{aligned}$$