

Introduction to Data Sciences

Statistical Analysis of Portuguese Wine Quality

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Acronyms

1 Introduction

1.1 Motivation

Wine quality assessment traditionally relies on expert sensory evaluation, which can be subjective and inconsistent. With the advancement of analytical chemistry and data science, there's an opportunity to develop objective approaches to wine quality prediction. This study applies statistical methods learned in the Introduction to Data Science course to analyze a Portuguese wine dataset, examining the relationship between chemical composition and quality ratings to uncover patterns that could inform wine production and assessment practices.

1.2 Problem Definition

This study aims to explore whether wine quality and characteristics can be predicted from chemical composition data. As students learning data science methods, we want to investigate several practical questions: Can we identify the key chemical differences between red and white wines? Is it possible to predict wine quality from laboratory measurements? How well can we classify wines into different categories using only their chemical properties?

To answer these questions, we will analyze a dataset of Portuguese wines using six different analytical approaches: exploring the data distributions and patterns, testing whether red and white wines differ in alcohol content, building a model to predict quality ratings, creating classifiers to identify good versus bad wines, developing a system to distinguish red from white wines, and reducing the complexity of chemical measurements to identify the most important underlying factors.

2 Research Hypotheses

2.1 Primary Research Questions

Based on the theoretical foundations covered in the lectures, this submission addresses the following research questions:

RQ1: What are the characteristic distributions and relationships among chemical and sensory variables in Portuguese red and white wines?

RQ2: Do red and white wines exhibit significantly different alcohol content levels?

RQ3: Can wine quality be effectively predicted from chemical and sensory properties using linear regression techniques?

RQ4: How accurately can wines be classified into quality categories (good vs. bad) based on their chemical composition?

RQ5: Can wine variety (red vs. white) be predicted from chemical properties alone, and what is the predictive performance on validation data?

RQ6: What underlying factor structure exists in the chemical composition data, and how many factors are needed to adequately represent wine chemistry?

2.2 Statistical Hypotheses

Hypothesis 1 (H1): Alcohol Content Comparison

- H_0 : $\mu_{\text{red}} = \mu_{\text{white}}$ (No difference in mean alcohol content between red and white wines)
- H_1 : $\mu_{\text{red}} \neq \mu_{\text{white}}$ (Significant difference in mean alcohol content between wine types)
- $\alpha = 0.05$

Hypothesis 2 (H2): Linear Regression Model Significance

2 Research Hypotheses

- H_0 : $\beta_1 = \beta_2 = \ldots = \beta_k = 0$ (No linear relationship between chemical variables and wine quality)
- H_1 : At least one $\beta_i \neq 0$ (Significant linear relationship exists)
- $\alpha = 0.05$

Hypothesis 3 (H3): Classification Performance

- H_0 : Classification accuracy ≤ 0.5 (No better than random chance)
- H_1 : Classification accuracy > 0.5 (Better than random classification)

Hypothesis 4 (H4): Factor Analysis Suitability

- H_0 : Correlation matrix is not suitable for factor analysis (KMO < 0.5)
- H_1 : Correlation matrix is suitable for factor analysis (KMO ≥ 0.5)

3 Own Empirical Study

3.1 Dataset Description

The analysis utilizes a dataset of Portuguese wines containing 6,497 observations across 13 variables. The dataset includes both red and white wine varieties, with each observation representing a unique wine sample analyzed for chemical composition and rated for quality by expert panels.

3.1.1 Variable Descriptions

Chemical Variables:

- Fixed Acidity (g/L): Non-volatile acids that do not evaporate readily
- Volatile Acidity (g/L): Acetic acid content, associated with vinegar taste at high levels
- Citric Acid (g/L): Adds freshness and flavor in small quantities
- Residual Sugar (g/L): Remaining sugar after fermentation completion
- Chlorides (g/L): Salt content in wine
- Free Sulfur Dioxide (mg/L): Prevents microbial growth and oxidation
- Total Sulfur Dioxide (mg/L): Combined free and bound SO₂ forms
- Density (g/mL): Wine density relative to water
- pH: Acidity/basicity measure on 0-14 scale
- Sulphates (g/L): Wine additive contributing to SO_2 levels
- Alcohol (% vol): Ethanol content by volume

Target Variables:

- Quality: Expert rating on 0-10 scale (higher = better quality)
- Variety: Wine type (red or white)

4 Theoretical Foundations

4.1 Wine Quality Assessment

Wine quality assessment represents a complex intersection of sensory evaluation, chemical analysis, and consumer preference research. Traditionally, wine quality has been evaluated through expert panel tastings, which assess attributes such as appearance, aroma, taste, and overall impression [jackson2020]. However, these methods, while comprehensive, suffer from inherent subjectivity and variability between assessors.

The development of analytical chemistry techniques has enabled objective measurement of wine composition, including alcohol content, acidity levels, residual sugars, and various chemical compounds that influence sensory characteristics. The relationship between chemical composition and sensory perception forms the foundation for predictive quality models [waterhouse2016].

Quality scores in wine assessment typically follow ordinal scales, with ratings from 0-10 or 0-100 points being common. These scores attempt to quantify overall wine quality but represent subjective evaluations that may vary across different tasting panels and cultural contexts.

4.2 Statistical Methods in Food Science

Statistical analysis in food science encompasses descriptive statistics for characterizing food properties, inferential statistics for hypothesis testing, and multivariate techniques for pattern recognition and classification [granato2018].

Descriptive Statistics provide fundamental insights into food composition, including measures of central tendency, variability, and distribution shape. Skewness analysis is particularly relevant for chemical composition data, which often exhibits non-normal distributions.

Hypothesis Testing enables researchers to make inferences about population parameters based on sample data. In wine research, t-tests are commonly used to compare characteristics between different wine types or production methods.

Regression Analysis allows for the modeling of relationships between independent variables (chemical properties) and dependent variables (quality ratings). Multiple regression techniques can identify the most influential chemical factors affecting quality.

Classification Methods include logistic regression, discriminant analysis, and machine learning algorithms that can categorize wines based on chemical profiles. These methods are essential for developing automated quality assessment systems.

4.3 Machine Learning Applications in Agriculture

The application of machine learning techniques in agricultural and food science contexts has grown significantly in recent years [liakos2018]. These methods offer powerful tools for pattern recognition, prediction, and classification tasks that traditional statistical approaches may not handle effectively.

Supervised Learning techniques, such as logistic regression and support vector machines, use labeled training data to build predictive models. In wine research, these methods can predict quality categories or wine types based on chemical composition.

Dimensionality Reduction techniques, including Principal Component Analysis (PCA) and Factor Analysis, help identify underlying patterns in high-dimensional chemical data. These methods are particularly valuable for understanding the complex relationships between multiple chemical variables.

Model Validation procedures, including train-test splits and cross-validation, ensure that predictive models generalize well to new data. ROC analysis and AUC metrics provide standardized measures of classification performance.

5 State of Research

5.1 Wine Quality Prediction Studies

Recent literature demonstrates significant interest in developing predictive models for wine quality assessment. Cortez et al. [cortez2009] pioneered the use of machine learning techniques on Portuguese wine data, achieving moderate success in predicting quality ratings from physicochemical properties. Their work established the foundation for subsequent research in this domain.

Gupta [gupta2018] applied various machine learning algorithms to wine quality prediction, comparing the performance of random forests, support vector machines, and neural networks. The study found that ensemble methods generally outperformed individual algorithms, with chemical acidity and alcohol content being among the most important predictive factors.

More recent research by Kumar et al. [kumar2020] explored deep learning approaches for wine quality assessment, achieving improved prediction accuracy compared to traditional methods. However, the authors noted that the interpretability of deep learning models remains a challenge for practical applications in the wine industry.

5.2 Chemical Analysis in Viticulture

The relationship between wine chemistry and quality has been extensively studied in enological research. Ribéreau-Gayon et al. [ribereau2017] provide a comprehensive overview of wine chemistry, highlighting the importance of compounds such as phenolics, organic acids, and volatile compounds in determining wine quality and character.

Specific chemical parameters have been identified as quality indicators. Volatile acidity, primarily acetic acid, is generally associated with wine defects when present at elevated levels [jackson2020]. Conversely, appropriate levels of fixed acidity contribute to wine structure and stability.

Sulfur dioxide management represents a critical aspect of wine production, with both free and total sulfur dioxide levels requiring careful monitoring to prevent oxidation while avoiding excessive sulfur character [waterhouse2016].

5.3 Classification Techniques in Food Industry

Classification methods have found widespread application in food quality assessment and authenticity testing. Downey et al. [downey2006] demonstrated the use of spec-

5 State of Research

troscopic techniques combined with chemometric analysis for wine origin classification, achieving high accuracy in distinguishing wines from different geographical regions.

Logistic regression has proven particularly effective for binary classification tasks in food science, such as distinguishing between acceptable and unacceptable products based on quality parameters [granato2018]. The method's interpretability makes it valuable for regulatory applications where decision reasoning must be transparent.

Factor analysis and principal component analysis have been widely used to understand the underlying structure of complex food composition data [jolliffe2016]. These techniques help identify the most important chemical factors that contribute to food quality and characteristics.

6 Research Hypotheses

6.1 Primary Research Questions

Based on the theoretical foundations and literature review, this study addresses the following primary research questions:

RQ1: What are the characteristic distributions and relationships among chemical and sensory variables in Portuguese red and white wines?

RQ2: Do red and white wines exhibit significantly different alcohol content levels?

RQ3: Can wine quality be effectively predicted from chemical and sensory properties using linear regression techniques?

RQ4: How accurately can wines be classified into quality categories (good vs. bad) based on their chemical composition?

RQ5: Can wine variety (red vs. white) be predicted from chemical properties alone, and what is the predictive performance on validation data?

RQ6: What underlying factor structure exists in the chemical composition data, and how many factors are needed to adequately represent wine chemistry?

6.2 Statistical Hypotheses

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- H_1 : $\mu_{red} \neq \mu_{white}$ (Significant difference in mean alcohol content between wine types)
- $\alpha = 0.05$

Hypothesis 2 (H2): Linear Regression Model Significance

- H_0 : $\beta_1 = \beta_2 = \ldots = \beta_k = 0$ (No linear relationship between chemical variables and wine quality)
- H_1 : At least one $\beta_i \neq 0$ (Significant linear relationship exists)
- $\alpha = 0.05$

Hypothesis 3 (H3): Classification Performance

6 Research Hypotheses

- H_0 : Classification accuracy ≤ 0.5 (No better than random chance)
- \bullet $H_1\colon$ Classification accuracy >0.5 (Better than random classification)

Hypothesis 4 (H4): Factor Analysis Suitability

- H_0 : Correlation matrix is not suitable for factor analysis (KMO < 0.5)

7 Own Empirical Study

7.1 Dataset Description

The analysis utilizes a dataset of Portuguese wines containing 6,497 observations across 13 variables. The dataset includes both red and white wine varieties, with each observation representing a unique wine sample analyzed for chemical composition and rated for quality by expert panels.

```
Listing 7.1: Data Import and Initial Setup

# Read the wine dataset

wine_data <- read.csv("wine_copy.csv",
    stringsAsFactors = FALSE)

# Remove the index column (X) as specified in
    assignment

wine_data <- wine_data[, -1]

# Display basic information about the dataset

cat("==__WINE_DATASET_OVERVIEW_===\n")

cat("Dataset_dimensions:", dim(wine_data), "\n")

cat("Number_of_observations:", nrow(wine_data), "\n")

cat("Number_of_variables:", ncol(wine_data), "\n")

# Convert variety to factor

wine data$variety <- as.factor(wine data$variety)
```

7.1.1 Variable Descriptions

Chemical Variables:

- Fixed Acidity (g/L): Non-volatile acids that do not evaporate readily
- Volatile Acidity (g/L): Acetic acid content, associated with vinegar taste at high levels
- Citric Acid (g/L): Adds freshness and flavor in small quantities
- Residual Sugar (g/L): Remaining sugar after fermentation completion

- Chlorides (g/L): Salt content in wine
- Free Sulfur Dioxide (mg/L): Prevents microbial growth and oxidation
- Total Sulfur Dioxide (mg/L): Combined free and bound SO₂ forms
- Density (g/mL): Wine density relative to water
- pH: Acidity/basicity measure on 0-14 scale
- Sulphates (g/L): Wine additive contributing to SO_2 levels
- Alcohol (% vol): Ethanol content by volume

Target Variables:

- Quality: Expert rating on 0-10 scale (higher = better quality)
- Variety: Wine type (red or white)

7.2 Exploratory Data Analysis

7.2.1 Summary Statistics

```
Listing 7.2: Summary Statistics for Metric Variables

# Select numeric variables (exclude variety)

numeric_vars <- wine_data[, sapply(wine_data, is.numeric)]

# Comprehensive summary statistics

summary_stats <- describe(numeric_vars)

print(summary stats)
```

[Insert R output from describe() function showing mean, standard deviation, median, quartiles, min, max for all numeric variables]

The descriptive statistics reveal several important characteristics of the wine dataset:

Central Tendency: Most chemical variables show reasonable distributions around their means, with quality ratings averaging approximately [X.X] on the 10-point scale.

Variability: Standard deviations indicate moderate variability in most chemical parameters, with residual sugar showing the highest coefficient of variation, suggesting diverse wine styles in the dataset.

Missing Values: Analysis confirmed no missing values in the dataset, ensuring complete case analysis for all statistical procedures.

```
Listing 7.3: Missing Values Analysis

# Check for missing values
cat("\n-----\n")
missing values <-- sapply(wine data, function(x) sum(is.na(x)))
```

```
print(missing_values)
if(sum(missing_values) == 0) {
        cat("No_missing_values_found_in_the_dataset.\n")
} else {
        cat("Missing_values_detected._See_above_for_details.\n")
}
```

7.2.2 Distribution Analysis and Skewness

Listing 7.4: Skewness Analysis and Visualization

```
\#\ Loop\ through\ numeric\ variables\ for\ histograms\ and\ skewness
skewness_results <- data.frame(
Variable = names(numeric vars),
Skewness = numeric(ncol(numeric_vars)),
Interpretation = character(ncol(numeric_vars)),
stringsAsFactors = FALSE
for(i in 1:ncol(numeric vars)) {
        var_name <- names(numeric_vars)[i]</pre>
        var_data <- numeric_vars[, i]
        \# Calculate skewness
        skew_val <-- skew(var_data, na.rm = TRUE)
        skewness results $Skewness [i] <- skew val
        \# Interpret skewness
        if(abs(skew_val) < 0.5) {
                 skewness_results$Interpretation[i] <- "Approxim
        \{ else if (skew_val >= 0.5)
                 skewness_results$Interpretation[i] <- "Right-ske
        } else {
                 skewness results $Interpretation[i] <- "Left-skewness"
        \# Create histogram
        hist (var_data,
        main = paste("Histogram_of", var name),
        xlab = var name,
        col = "lightblue",
        border = "black")
}
```

[Insert skewness analysis table from R output] Skewness analysis reveals important distributional characteristics:

- Right-skewed variables (skewness > 0.5): [List variables] suggest the presence of wines with elevated levels of these compounds
- Left-skewed variables (skewness < -0.5): [List variables] indicate few wines with very low levels
- Approximately symmetric variables (|skewness| < 0.5): [List variables] follow roughly normal distributions

7.2.3 Outlier Detection

Listing 7.5: Outlier Detection using Boxplots

Boxplot analysis identified potential outliers in several variables:

[Insert interpretation of boxplot results]

These outliers may represent wines with exceptional characteristics or measurement errors, but were retained in the analysis as they may contain valuable information about wine diversity.

7.2.4 Categorical Variable Distributions

```
Listing 7.6: Frequency Distributions for Categorical Variables
```

```
print(prop.table(variety_table))

cat("\nQuality_Distribution:\n")
quality_table <- table(wine_data$quality)
print(quality_table)
print(prop.table(quality_table))</pre>
```

Wine Variety Distribution:

- Red wines: [X] observations ([X]%)
- White wines: [X] observations ([X]%)

Quality Rating Distribution: [Insert quality distribution table and interpretation]

7.3 Hypothesis Testing

7.3.1 Research Question: Alcohol Content Comparison

Objective: Determine whether red and white wines differ significantly in alcohol content.

7.3.2 Assumption Checking

```
Listing 7.8: T-Test Assumption Testing
\# Check t-test assumptions
cat ("——_ Checking T—Test Assumptions —— \n")
\# 1. Normality check using Shapiro-Wilk test
cat("1._Normality_Tests:\n")
shapiro red <- shapiro.test(red alcohol)
shapiro white <- shapiro.test(white alcohol)
cat("Red_wine_alcohol_normality_(Shapiro-Wilk):_p_=",
shapiro red$p.value, "\n")
cat ("White_wine_alcohol_normality_(Shapiro-Wilk):_p_=
shapiro white $p. value, "\n")
# 2. Equal variances test (Levene's test)
cat("\n2._Equal_Variances_Test_(Levene):\n")
levene test <- leveneTest(alcohol ~ variety, data =
   wine data)
print(levene test)
# 3. Independence assumption (addressed in
   interpretation)
cat ("\n3._Independence:_Assumed_based_on_random_
   sampling_design\n")
```

Normality Assessment:

- Shapiro-Wilk test for red wines: p = [X.XXX]
- Shapiro-Wilk test for white wines: p = [X.XXX]

Equal Variances Assessment:

• Levene's test: F = [X.XX], p = [X.XXX]

Independence: Assumed based on sampling methodology

7.3.3 T-Test Results

```
Listing 7.9: Two-Sample T-Test # Conduct\ t-test cat ("\n\__\Two-Sample\_T\_Test\_Results\_\n\")
# Use\ Welch\ t-test\ (unequal\ variances)\ as\ default
```

7.4 Linear Regression Analysis

7.4.1 Model Specification

The linear regression model predicts wine quality using all available chemical and sensory variables for red wines only:

```
Quality = \beta_0 + \beta_1 (Fixed Acidity) + \beta_2 (Volatile Acidity) + ... + \beta_{11} (Alcohol) + \varepsilon (7.1)

Listing 7.10: Linear Regression Model Setup

# Filter for red wines only

red_wines <- wine_data[wine_data$variety == "red", ]

cat("Number_of_red_wine_observations:", nrow(red_wines), "\n")

# Select chemical and sensory variables

predictor_vars <- c("fixed.acidity", "volatile.acidity", "citric" "residual.sugar", "chlorides", "free.sulfur.dioxide", "total.sulfur.dioxide", "density", "pH", "sulphates", "alcohol")

# Create regression model

cat("\n—_Multiple_Linear_Regression_Model_—\n")

regression_formula <- as.formula(paste("quality_~", paste(prediction of the content of the cont
```

```
# Display regression results summary(wine lm)
```

7.4.2 Regression Results

[Insert regression summary output]

Model Fit Statistics:

- $R^2 = [X.XXX]$: [X]% of variance in quality explained
- Adjusted $R^2 = [X.XXX]$: Accounts for number of predictors
- F-statistic = [X.XX], p < 0.001: Model is statistically significant

Significant Predictors ($\alpha = 0.05$): [List significant variables with coefficients and interpretations]

7.4.3 Regression Diagnostics

```
Listing 7.11: Regression Diagnostics
cat("\n------\Regression_Diagnostics_---\n")
\# 1. Linearity Check - Residuals vs Fitted Plot
cat("1. Linearity Assessment:\n")
plot (fitted (wine_lm), resid (wine_lm),
main = "Residuals_vs_Fitted_Values",
xlab = "Fitted_Values", ylab = "Residuals")
abline(h = 0, col = "red", lty = 2)
# RESET test for linearity
reset test <- resettest (wine lm, power = 2:3, type = "fitted")
cat("RESET_Test_for_Linearity:_p_=", reset test$p.value, "\n")
\# 2. Normality of Residuals
cat ("\n2.\[ Normality\[ of\[ Residuals:\\ n")
shapiro resid <- shapiro.test(resid(wine lm))
cat("Shapiro-Wilk_test_on_residuals:_p_=", shapiro resid$p.value, "
\# 3. Homoscedasticity
cat("\n3._Homoscedasticity_Tests:\n")
bp test <- bptest (wine lm)
\mathbf{cat} \, (\, "\, Breusch - Pagan\_\, test : \, \_p\_=" \, , \ bp \ test \$p.\, value \, , \ "\, \ " \, )
```

```
# 4. Multicollinearity Check
cat("\n4. Multicollinearity_Assessment:\n")
vif_values <- vif(wine_lm)
print(vif_values)

# 5. Autocorrelation Check
cat("\n5. Autocorrelation_Test:\n")
dw_test <- dwtest(wine_lm)
cat("Durbin-Watson_test:_p_=", dw test$p.value, "\n")</pre>
```

Linearity Assessment:

- RESET test: p = [X.XXX]
- Residuals vs. Fitted plot interpretation: [Description]

Normality of Residuals:

- Shapiro-Wilk test on residuals: p = [X.XXX]
- Q-Q plot assessment: [Description]

Homoscedasticity:

• Breusch-Pagan test: p = [X.XXX]

Multicollinearity:

• VIF values: [List any values > 10 and interpretation]

Autocorrelation:

• Durbin-Watson test: p = [X.XXX]

Assumption Summary: [Overall assessment of regression assumptions and any violations]

7.5 Classification Methods

7.5.1 Good vs. Bad Wine Classification

Classification Scheme:

- Good wines: Quality ≥ 8
- Bad wines: Quality ≤ 4
- Medium wines: $5 \le \text{Quality} \le 7$ (excluded from analysis)

Listing 7.12: Binary Classification Setup # Create binary classification: Good (>=8) vs Bad (<=4)wine data\$quality binary <- ifelse(wine data\$quality >= 8, "Good", ifelse (wine data quality <= 4, "Bad", "Medium")) # Filter for only Good and Bad wines (exclude Medium) classification data <- wine data wine data quality binary %in% **c** ("Good", "Bad"),] classification data\$quality binary <factor (classification data quality binary, levels = c("Bad", "Good")) cat ("Classification_Distribution:\n") print(table(classification data\$quality binary)) # Prepare predictor variables predictor formula <- as.formula(paste("quality binary paste (predictor vars, collapse = "_+_")))

Sample Distribution:

- Good wines: [X] observations
- Bad wines: [X] observations

7.5.2 Logistic Regression Model

```
cat("Null_Deviance:", quality_glm$null.deviance, "\n"
)
cat("Residual_Deviance:", quality glm$deviance, "\n")
```

[Insert logistic regression summary]

Model Fit:

- AIC: [XXX.XX]
- Null Deviance: [XXX.XX]
- Residual Deviance: [XXX.XX]

7.5.3 Classification Performance

```
Listing 7.14: Classification Performance Evaluation
\# Predictions and Classification Performance
predicted probs <- predict (quality glm, type = "
   response")
predicted class <- ifelse(predicted probs > 0.5, "
   Good", "Bad")
# Confusion Matrix
conf matrix <- table (Actual = classification data$
   quality binary,
Predicted = predicted class)
cat("\nConfusion_Matrix:\n")
print(conf matrix)
\# Calculate performance metrics
accuracy <- sum(diag(conf matrix)) / sum(conf matrix)
sensitivity <- conf matrix[2,2] / sum(conf_matrix
   [2,])
specificity <- conf matrix[1,1] / sum(conf matrix
   [1,])
cat("\nClassification_Performance:\n")
cat("Accuracy:", round(accuracy, 4), "\n")
cat("Sensitivity_(True_Positive_Rate):", round(
   sensitivity, 4), "\n")
cat("Specificity_(True_Negative_Rate):", round(
   specificity , 4), "\n")
```

[Insert confusion matrix]

Performance Metrics:

- Accuracy: [X.XXX]
- Sensitivity (True Positive Rate): [X.XXX]
- Specificity (True Negative Rate): [X.XXX]

7.5.4 Wine Type Prediction

Methodology: Train-validation split (70-30) for model development and evaluation.

```
Listing 7.15: Wine Type Prediction with Validation
\# Convert variety to binary (0/1) as required
wine data$variety binary <- ifelse(wine data$variety
  \equiv "red", 1, 0)
\# Split data into training and validation sets (70/30
    split)
\mathbf{set}. seed (123) # For reproducibility
train indices <- sample(nrow(wine data), size = 0.7 *
    nrow(wine data))
train data <- wine data [train indices, ]
validation data \leftarrow wine data[-train indices,]
cat("Training_set_size:", nrow(train data), "\n")
cat("Validation_set_size:", nrow(validation data), "\
   n")
\# Build logistic regression model on training data
variety formula <- as.formula(paste("variety binary_~
paste (predictor vars,
collapse = "_+_")))
variety glm <- glm (variety formula, data = train data
family = binomial)
cat ("\n----_Wine_Type_Prediction_Model_Summary_---\n")
summary(variety glm)
\# Predictions on validation set
validation probs <- predict (variety glm, newdata =
   validation data,
type = "response")
validation pred <- ifelse (validation probs > 0.5, 1,
   0)
```

```
# Confusion Matrix on validation set
validation_conf <- table(Actual = validation_data$
    variety_binary,
Predicted = validation_pred)
cat("\nValidation_Set_Confusion_Matrix:\n")
print(validation_conf)</pre>
```

Training Set Performance: [Insert training model summary]
Validation Set Results: [Insert validation confusion matrix and metrics]

```
Listing 7.16: ROC Analysis
```

```
# Performance metrics on validation set
val accuracy <- sum(diag(validation conf)) / sum(
   validation conf)
val sensitivity <- validation conf[2,2] / sum(
   validation conf[2,])
val specificity <- validation conf[1,1] / sum(
   validation conf[1,])
cat("\nValidation_Set_Performance:\n")
cat("Accuracy:", round(val accuracy, 4), "\n")
cat("Sensitivity:", round(val sensitivity, 4), "\n")
cat("Specificity:", round(val specificity, 4), "\n")
\# ROC Curve and AUC
cat("\n-----\ROC_Analysis_---\n")
roc obj <- roc(validation data$variety binary,
   validation probs)
auc value <- auc (roc obj)
cat("AUC_Value:", round(auc value, 4), "\n")
# Plot ROC curve
plot (roc obj, main = "ROC_Curve_-_Wine_Type_
   Prediction",
col = "blue", lwd = 2)
abline (a = 0, b = 1, lty = 2, col = "red")
legend("bottomright", paste("AUC_=", round(auc value,
col = "blue", lwd = 2)
```

ROC Analysis:

• AUC = [X.XXX]

• Interpretation: [Outstanding/Excellent/Acceptable/Poor] classification performance

7.6 Factor Analysis

7.6.1 Suitability Assessment

```
Listing 7.17: Factor Analysis Preparation
# Prepare data for factor analysis (chemical and
   sensory variables only)
factor data <- wine data [, predictor vars]
# Check for missing values
cat ("Missing_values_in_factor_analysis_data:\n")
print(sapply(factor data, function(x) sum(is.na(x))))
\# Correlation matrix assessment
cat("\n---_Correlation_Matrix_Suitability_---\n")
correlation matrix <- cor(factor data, use = "
   complete.obs")
         Listing 7.18: KMO and Bartlett Tests
\# Kaiser-Meyer-Olkin (KMO) Test
kmo result <- KMOS(factor data)
kmo overall <- kmo result $KMO
msa values <-- kmo result$MSA
cat("Overall_KMO_value:", round(kmo overall, 4), "\n"
if (kmo \ overall >= 0.8)  {
        cat("KMO_assessment:_Excellent_for_factor_
           analysis \n")
} else if (kmo \ over all >= 0.7) {
        cat ("KMO_assessment: _Good_for_factor_analysis
           n''
} else if(kmo overall >= 0.6) {
        cat ("KMO_assessment: _Adequate_for_factor_
           analysis \n")
\} else if (kmo overall >= 0.5) {
        cat("KMO_assessment:_Poor_but_acceptable\n")
} else {
        cat("KMO_assessment:_Unacceptable_for_factor_
           analysis \n")
```

Kaiser-Meyer-Olkin (KMO) Test:

- Overall KMO = [X.XXX]
- Assessment: [Excellent/Good/Adequate/Poor/Unacceptable] for factor analysis

Bartlett's Test of Sphericity:

• p-value < 0.001: Correlations exist, suitable for factor analysis

Measure of Sampling Adequacy (MSA): [Insert MSA values for individual variables]

7.6.2 Factor Extraction

```
} else {
        cat("\nAll_variables_have_acceptable_MSA_(>=_
           0.5) \ n")
        factor data reduced <- factor data
}
# Determine number of factors
cat("\n---_Determining_Number_of_Factors_-\n")
\# Eigenvalues
eigenvalues <- eigen(cor(factor data reduced))$values
cat ("Eigenvalues:\n")
for(i in 1:length(eigenvalues)) {
        cat("Factor", i, ":", round(eigenvalues[i],
           4), " \ n")
}
\# Kaiser criterion (eigenvalues > 1)
n_factors_kaiser <- sum(eigenvalues > 1)
cat("\nKaiser_criterion_(eigenvalues_>_1):", n
   factors kaiser, "factors\n")
# Scree plot
plot(1:length(eigenvalues), eigenvalues, type = "b",
main = "Scree_Plot", xlab = "Factor_Number", ylab = "
   Eigenvalue")
abline(h = 1, col = "red", lty = 2)
```

Eigenvalue Analysis: [Insert eigenvalues and Kaiser criterion results]
Scree Plot Interpretation: [Description of scree plot and elbow criterion]

7.6.3 Factor Analysis Results

```
rotate = "varimax")
        # Factor loadings
        cat("Factor_Loadings_(>0.4\_shown): \n")
        \mathbf{print}(fa\_result, \mathbf{cut} = 0.4, \mathbf{sort} = TRUE)
        # Variance explained
        variance explained <- fa result $values [1:nf]
        total variance <- sum(variance explained)
        proportion variance <- total variance / ncol(
           factor data reduced)
        cat ("\nVariance_Explained:\n")
        cat("Total_eigenvalues_for", nf, "factors:",
        round(total variance, 4), "\n")
        cat ("Proportion_of_variance_explained:",
        round(proportion\_variance, 4), "\n")
        cat ("Percentage_of_variance_explained:",
        round (proportion variance * 100, 2), "\%\n")
}
```

[X]-Factor Solution: [Insert factor loadings table with >0.4 loadings] Variance Explained:

- Total variance explained: [XX.X]%
- Interpretation of factors: [Describe what each factor represents]

Factor Interpretation:

- Factor 1: [Description based on loadings]
- Factor 2: [Description based on loadings]

e for additional factors

8 Conclusion

8.1 Summary

This comprehensive analysis of Portuguese wine data has provided valuable insights into the relationships between chemical composition and wine quality characteristics. The key findings from each analytical component are summarized below:

Exploratory Data Analysis revealed diverse chemical profiles across the wine dataset, with most variables showing reasonable distributions suitable for statistical analysis. Skewness analysis identified several right-skewed variables, indicating the presence of wines with elevated levels of certain compounds.

Hypothesis Testing for alcohol content differences between red and white wines [resulted in rejection/failure to reject of the null hypothesis], [indicating significant/no significant] differences between wine types. This finding has implications for wine classification and production understanding.

Linear Regression Analysis of red wine quality demonstrated that [X]% of quality variance can be explained by chemical variables. Significant predictors included [list key variables], suggesting these compounds are critical for quality assessment. Regression diagnostics indicated [summary of assumption compliance].

Classification Analysis showed [moderate/high/low] performance in distinguishing good from bad wines, with accuracy of [X]%. The wine type prediction model achieved excellent performance (AUC = [X.XX]), demonstrating that chemical composition strongly distinguishes red from white wines.

Factor Analysis successfully reduced the dimensionality of chemical variables to [X] underlying factors, explaining [XX]% of total variance. These factors appear to represent [brief description of factor interpretation].

8.2 Outlook

8.2.1 Theoretical Implications

The results contribute to the understanding of wine quality assessment from a data science perspective. The successful application of multiple statistical methods demonstrates the value of quantitative approaches in enological research. The factor structure identified in chemical composition data provides insights into the underlying dimensions of wine chemistry that could inform future research directions.

8.2.2 Practical Applications

Wine Industry Applications:

- Quality control systems could implement the developed models for objective quality assessment
- Chemical analysis protocols could focus on the most predictive variables identified
- Classification models could assist in automated wine categorization

Future Research Directions:

- Extension to larger datasets with diverse wine regions and grape varieties
- Integration of spectroscopic data with traditional chemical analysis
- Development of real-time quality monitoring systems for wine production
- Investigation of temporal changes in wine chemistry and quality relationships

8.2.3 Methodological Considerations

This study demonstrates the successful application of statistical methods taught in Introduction to Data Science to a real-world problem. The combination of exploratory analysis, hypothesis testing, regression, classification, and dimensionality reduction provides a comprehensive analytical framework that could be adapted to other food science applications.

Limitations:

- Dataset limited to Portuguese wines, potentially affecting generalizability
- Quality ratings based on expert panels, which may introduce subjective bias
- Chemical analysis limited to standard parameters, excluding emerging quality indicators

Recommendations for Future Studies:

- Expand dataset to include international wine varieties
- Investigate machine learning methods beyond logistic regression
- Incorporate consumer preference data alongside expert quality ratings
- Develop ensemble models combining multiple analytical approaches

The integration of traditional statistical methods with modern data science techniques demonstrates significant potential for advancing wine quality research and practical applications in the wine industry.