

sommelieR

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Introduction

- ▶ Wine tasting can range from a casual pastime to a lucrative profession
- ▶ For professional sommeliers, considerable time and training is required to adequately rate wine quality
- ▶ Intuitively, we might expect expert ratings to reflect the underlying chemical composition of the wines
- ▶ We decided to more formally investigate this idea

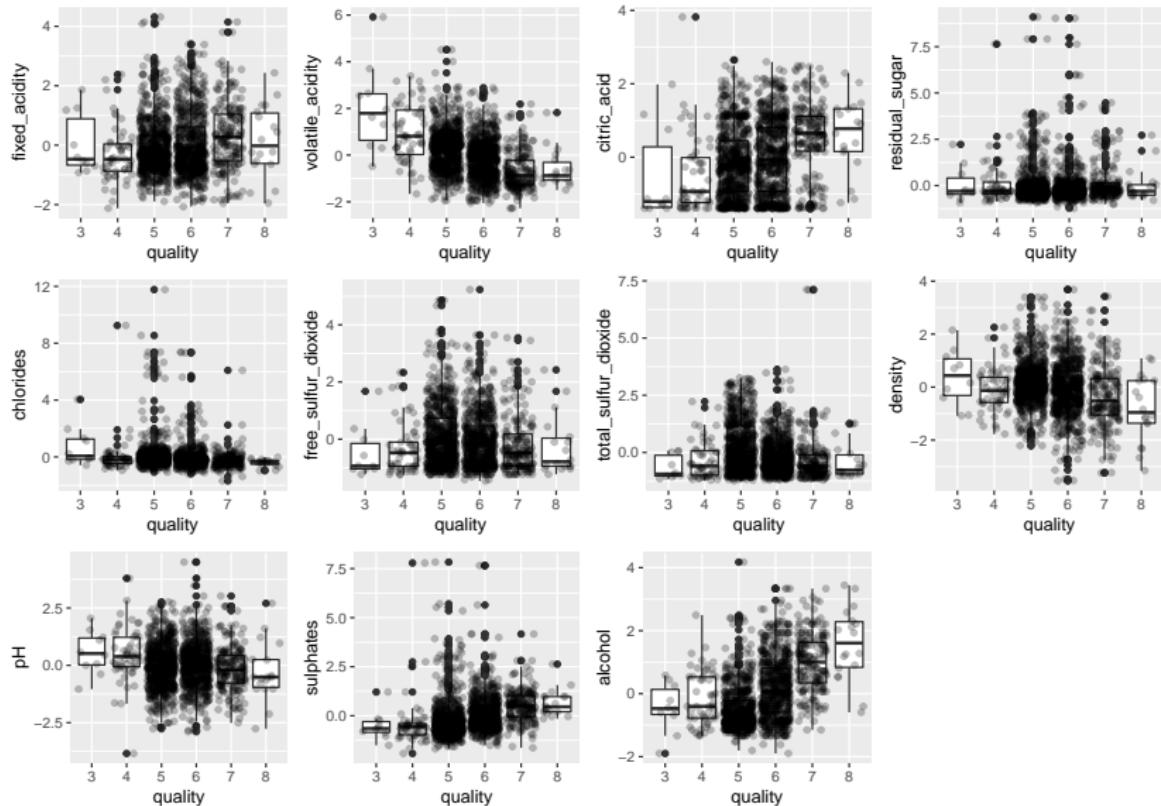
Project Aim

To determine how accurately expert wine quality ratings can be predicted using a set of easily measured chemical components.

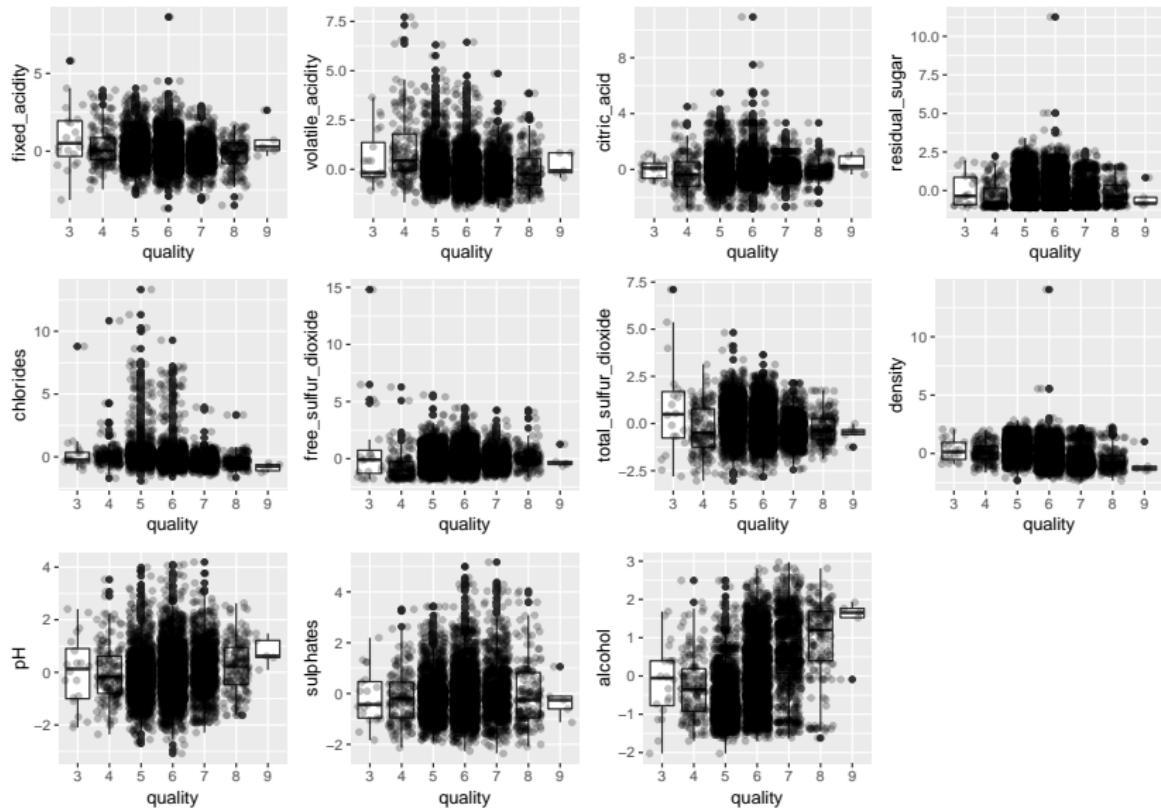
Data

- ▶ 2 datasets of expert quality ratings of red and white Vinho Verde wines
 - ▶ source: <http://archive.ics.uci.edu/ml/datasets/wine+quality>
 - ▶ 1599 red wines
 - ▶ 4898 white wines
- ▶ Outcome variable: wine quality
 - ▶ ordinal variable theoretically ranging from 0-10
 - ▶ Observed ratings range from 3-9
 - ▶ 0 is poor, 10 is excellent
 - ▶ classes are unbalanced
- ▶ Predictor variables: 11 physiochemical wine components:
 - ▶ fixed acidity, volatile acidity, citric acid, residual sugar, chlorides, free sulfur dioxide, total sulfur dioxide, density, pH, sulphates, alcohol

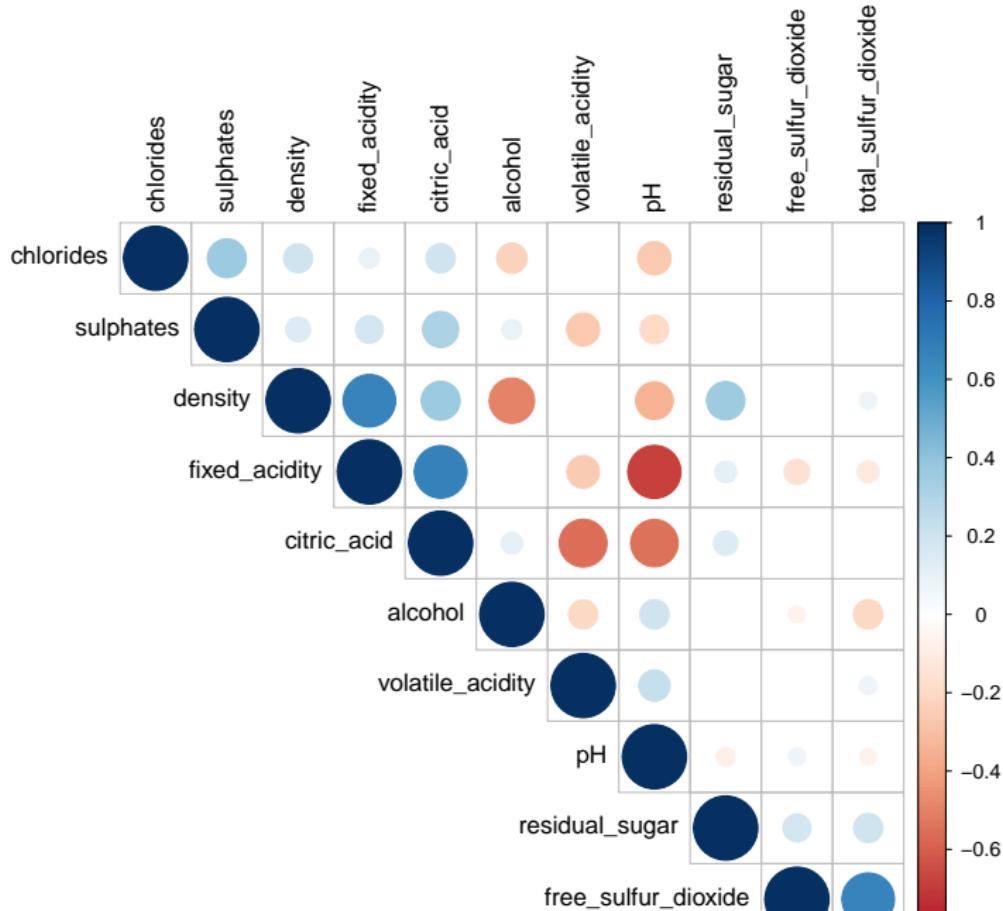
Looking At the Red Wine Data



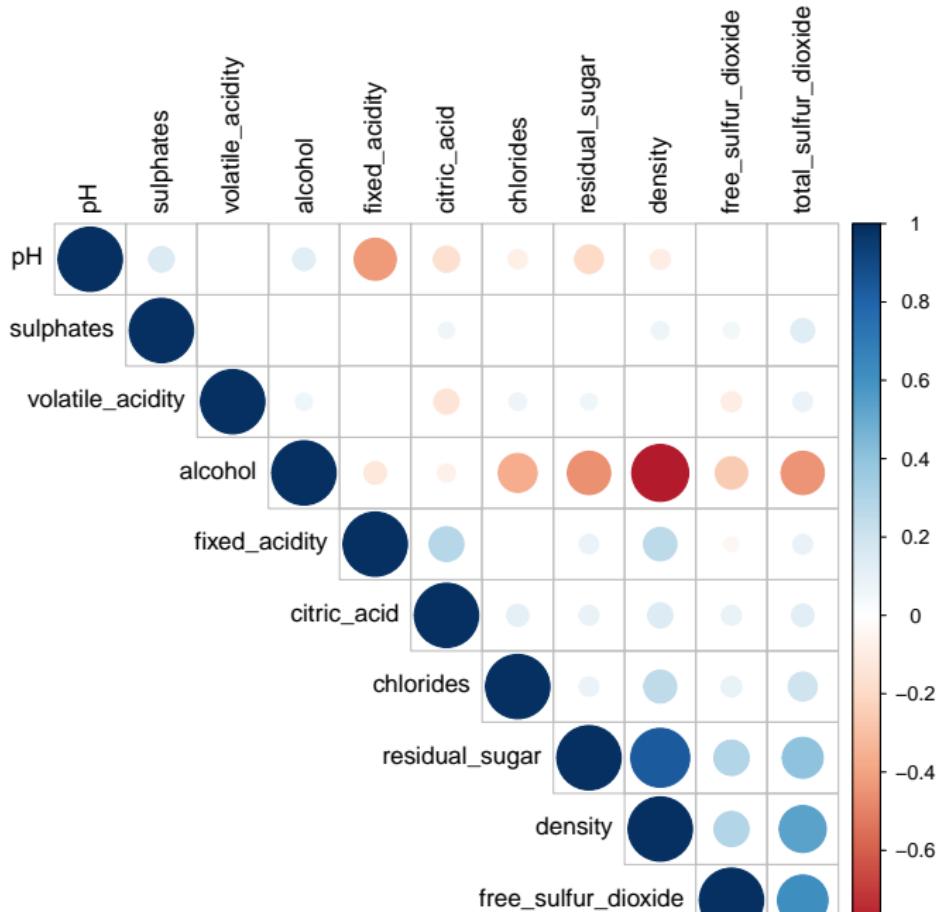
Looking At the White Wine Data



Correlations- Red Wine



Correlations- White Wine



Methods

- ▶ The red and white wine datasets were split into training and testing sets
 - ▶ 80% of available data sampled for training models
 - ▶ Remaining 20% used to evaluate accuracy
 - ▶ Relative frequencies of quality ratings in full dataset were preserved in training and testing data
- ▶ Modeling Approach
 - ▶ Machine learning
 - ▶ Random forest classification of quality ratings
 - ▶ Likelihood based
 - ▶ linear regression
 - ▶ partial proportional odds model
 - ▶ multinomial model

Methods: Linear Regression

- ▶ $Z = X^T \beta + \epsilon$, where $\epsilon \sim N(0, \sigma^2 I)$
- ▶ $\hat{\beta} = (X^T X)^{-1} X^T Z$
- ▶ Predicted classification: The nearest integer to the predicted mean rating

Methods: Partial Proportional Odds Models

Three different approaches were considered:

- ▶ Non-proportional odds: $\text{logit}(P(Z \leq j|x)) = \alpha_j + x^T \beta_j$
- ▶ Proportional odds: $\text{logit}(P(Z \leq j|x)) = \alpha_j + x^T \beta$
- ▶ Partial proportional odds:
$$\text{logit}(P(Z \leq j|x)) = \alpha_j + x_*^T \beta + x_{**}^T \beta_j$$
- ▶ $\alpha_j > \alpha_i$ for $j > i$
- ▶ Log likelihood maximized using BFGS (default) or any available method in package optimx
- ▶ Predicted classification: the rating with the highest predicted probability

Methods: Multinomial Regression

- ▶ Response takes values in K classes
- ▶ Model: $\log \frac{Pr(G=k|X=x)}{P(G=K|X=x)} = X\beta_k$ for $k = 1, \dots, (K - 1)$
- ▶ Predictions:
$$P(G = k|X = x) = \frac{\exp(X\beta_k)}{1 + \sum_{l=1}^{K-1} \exp(X\beta_l)}, k = 1, \dots, K - 1$$
- ▶ Considered 3 Models: All Predictors with Linear Terms, Reduced Model with Linear Terms, and Reduced Model with All Second Order Terms
- ▶ Computational Efficiency Challenges

Methods: Random Forest

- ▶ Something here describing the random forest approach

Variable Selection

- ▶ Random Forest: all physiochemical variables included
- ▶ Likelihood based models:
 - ▶ We examined the correlations between predictors and considered best subsets found in R based on linear regression.
 - ▶ The final predictor variables selected were:
 - ▶ red wine: volatile acidity, total sulfur dioxide, pH, alcohol, sulphates
 - ▶ white wine: pH, volatile_acidity, residual_sugar, alcohol

Model Evaluation

Models were compared on the following metrics:

- ▶ Accuracy
 - ▶ Correct Classifications/Total
- ▶ Kappa
- ▶ Weighted Kappa - used to select the final model
 - ▶ A more useful version of Kappa for data with inherent ordering that penalizes misclassifications proportional to the distance from the true category.
 - ▶ For instance, when the true quality rating is 4, a prediction of 7 will be penalized more severely than a prediction of 5.

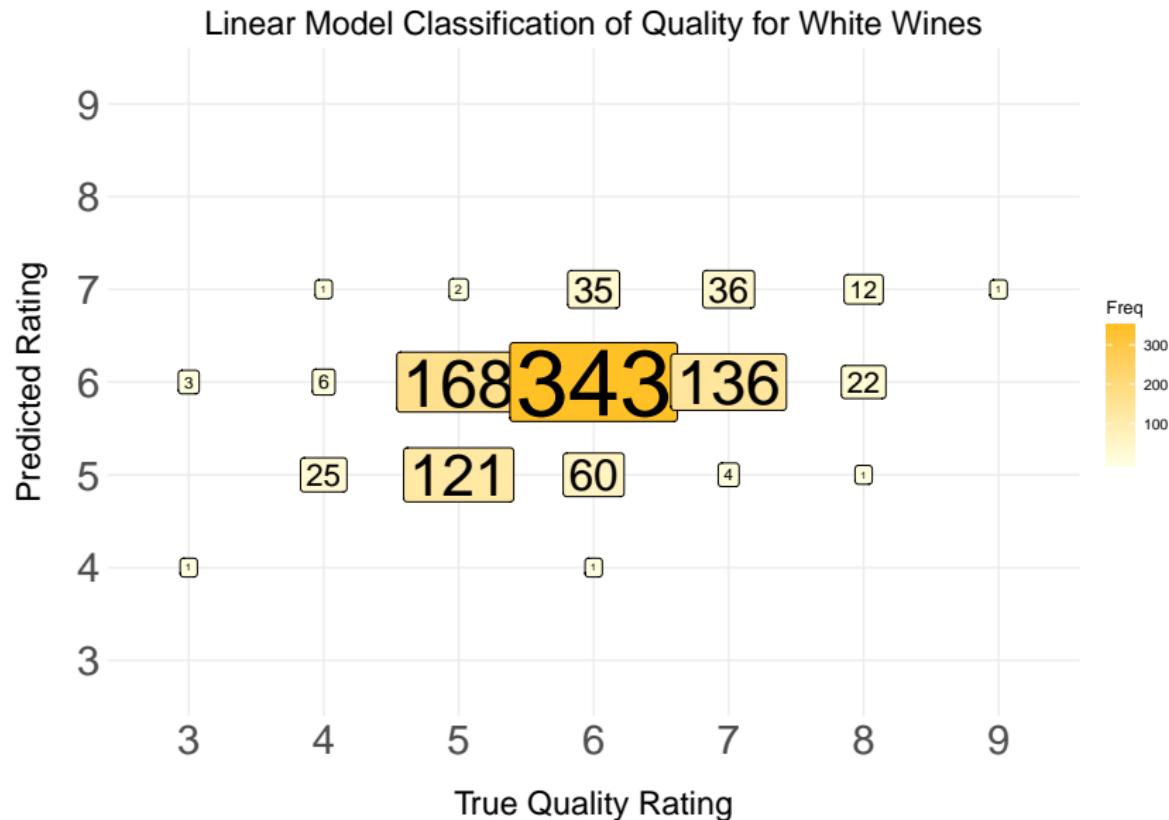
Results: Linear Model

- ▶ Red Wine
 - ▶ Accuracy: 56.15
 - ▶ Kappa: 0.2728
 - ▶ Weighted Kappa: 0.4545
- ▶ White Wine
 - ▶ Accuracy: 51.12
 - ▶ Kappa: 0.1969
 - ▶ Weighted Kappa: 0.4067

Results: Linear Model (Red Wine)



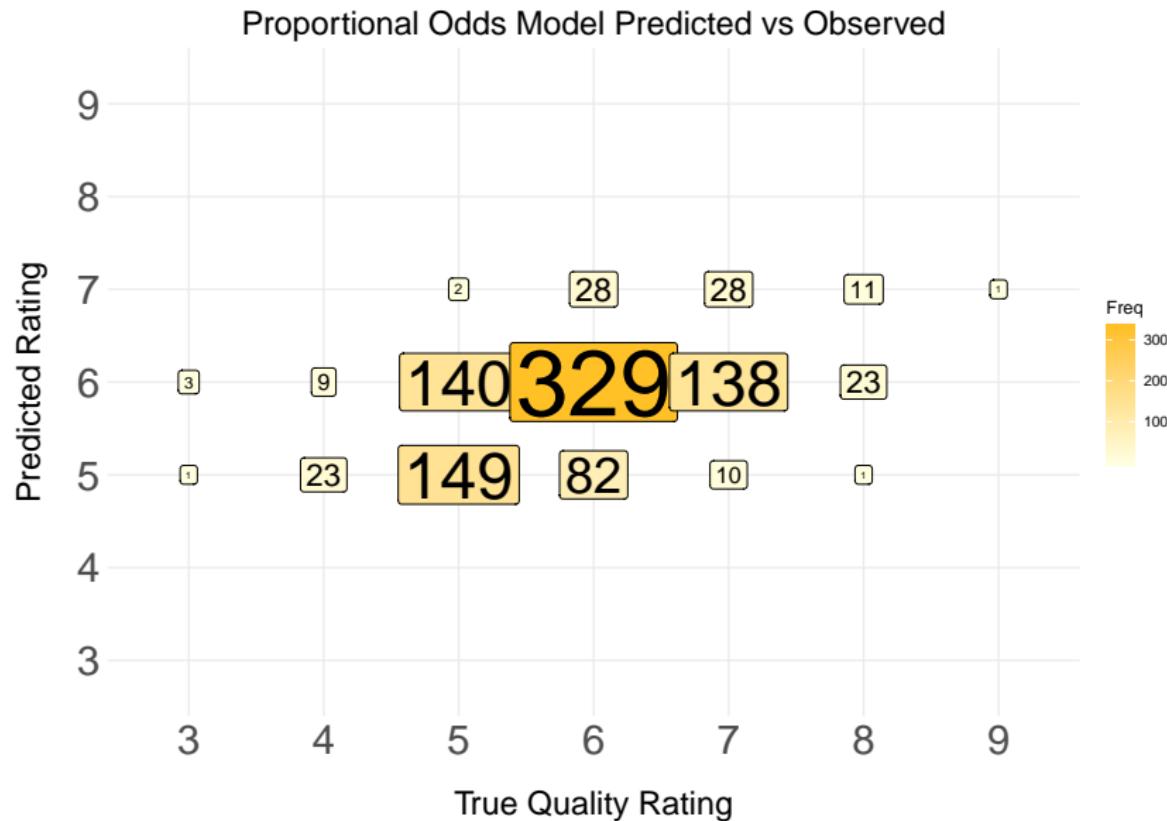
Results: Linear Model (White Wine)



Results: Partial Proportional Odds Model (White Wine)

- ▶ For white wine, only the proportional odds model converged
 - ▶ Overall Accuracy: 51.74%
 - ▶ Kappa: 0.2108
 - ▶ Weighted Kappa: 0.3993

Results: Proportional Odds Model (White Wine)



Results: Partial Proportional Odds Model (Red Wine)

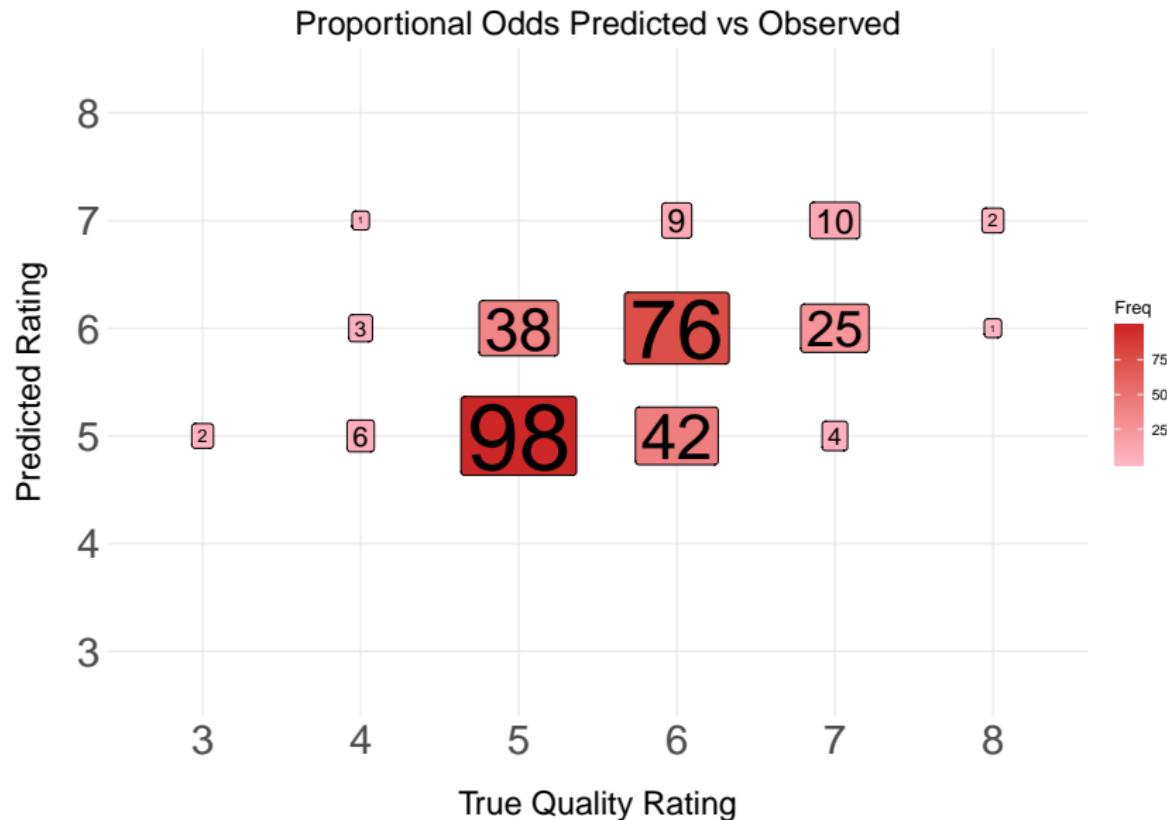
- ▶ Proportional, partial proportional, and non-proportional odds models all converged
- ▶ Results presented for partial proportional model with coefficient for total sulfur dioxide allowed to vary with the level of wine quality

Results: Comparison of Partial Proportional Odds Models (Red Wine)

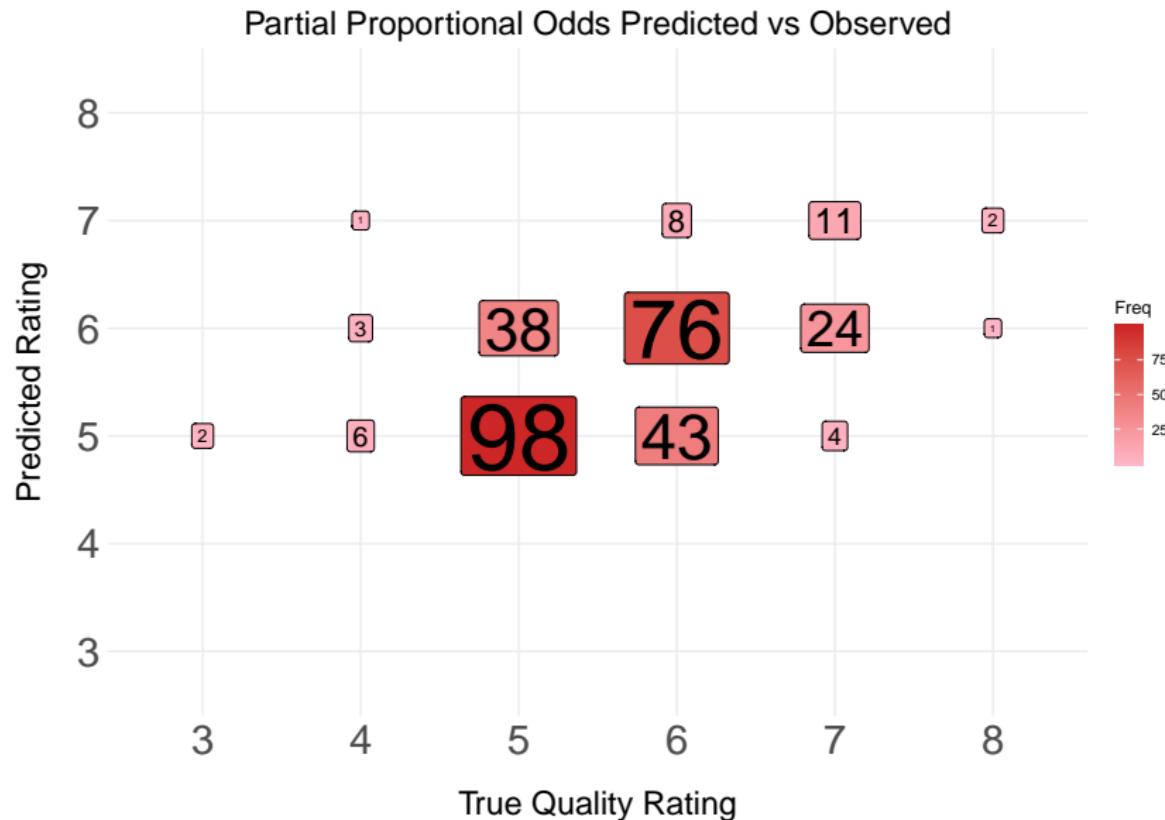
Table 1: Comparison of Partial Proportional Odds Models for Red Wine Quality

Model	Accuracy	Kappa	Weighted Kappa
Proportional Odds	58.0442	0.3065	0.4707
Partial Proportional Odds	58.3596	0.3117	0.4742
Non-Proportional Odds	57.7287	0.3080	0.4357

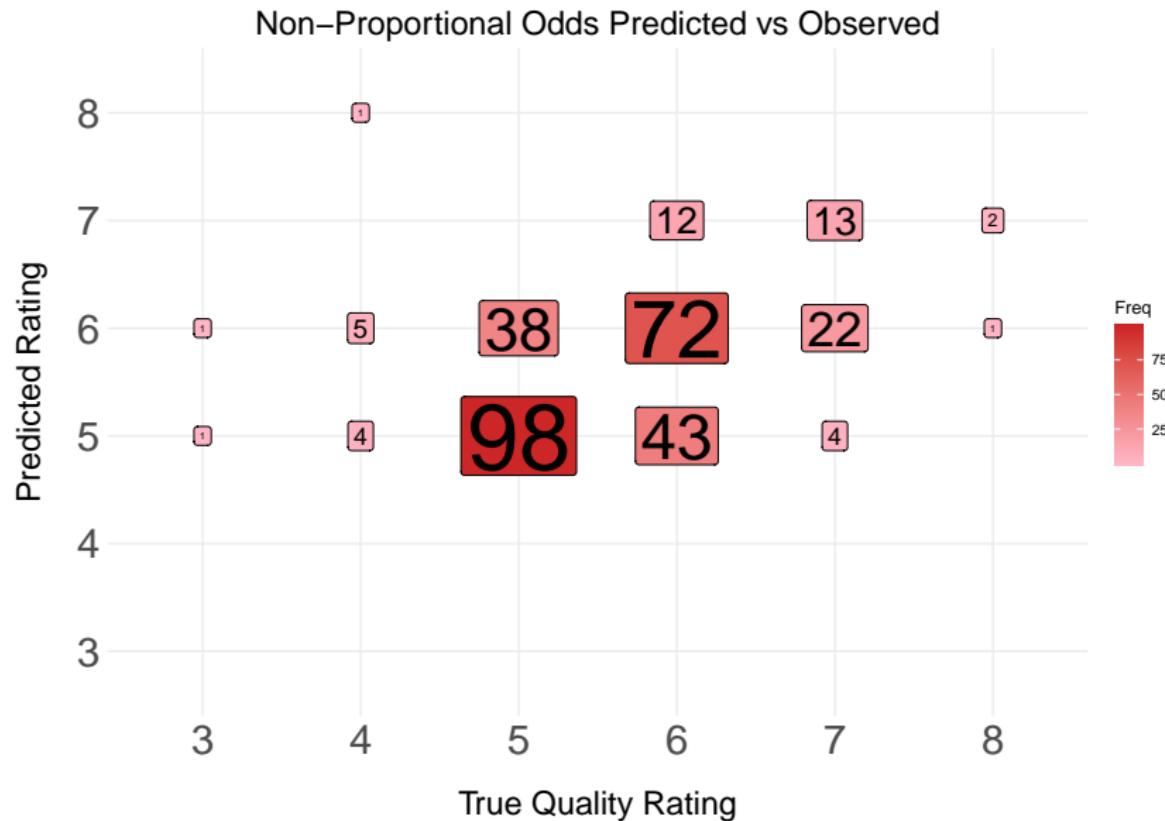
Results: Proportional Odds Model (Red Wine)



Results: Partial Proportional Odds Model (Red Wine)



Results: Non-Proportional Odds Model (Red Wine)



Results: Multinomial Regression (Red Wine Quality Classification)

Table 2: Comparison of Multinomial Regression Models for Red Wine Quality

Model	Accuracy	Kappa	Weighted Kappa	3	4	5	6	7	8
Full Model (Linear Terms)	58.3596	0.3158	0.5227	0	20	74.2647	55.9055	28.2051	0
Reduced Model (Linear Terms)	58.3596	0.3193	0.4952	0	10	72.7941	56.6929	33.3333	0
Reduced Model (Second Order Terms)	55.2050	0.2702	0.4789	0	0	67.6471	55.1181	33.3333	0

Results: Multinomial Regression - Full Model Confusion Matrix (Red Wine)

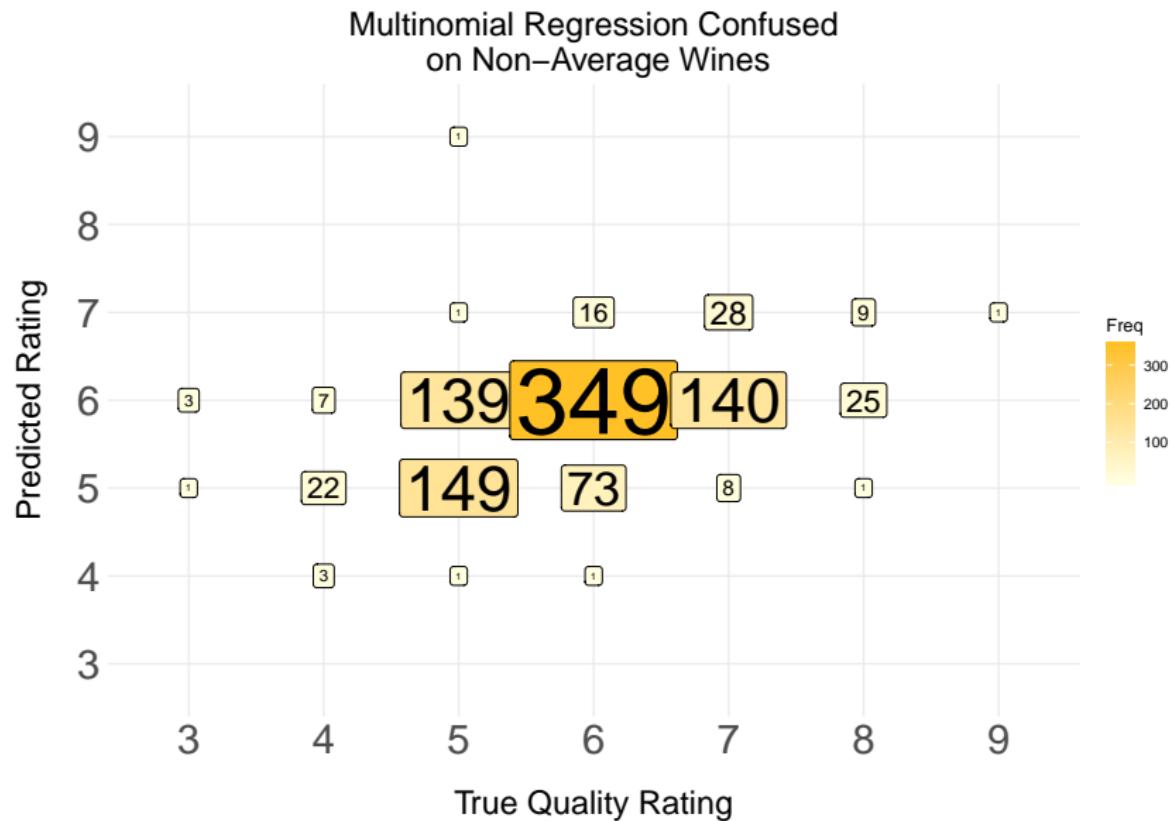


Results: Multinomial Regression (White Wine Quality Classification)

Table 3: Comparison of Multinomial Regression Models for Red Wine Quality

Model	Accuracy	Kappa	Weighted Kappa	3	4	5	6	7	8	9
Full Model (Linear Terms)	54.0900	0.2451	0.4121	0	9.375	51.2027	79.4989	15.9091	0	0
Reduced Model (Linear Terms)	51.9427	0.2201	0.4093	0	3.125	54.6392	72.6651	16.4773	0	0
Reduced Model (Second Order Terms)	53.2720	0.2396	0.4010	0	3.125	54.2955	74.4875	19.8864	0	0

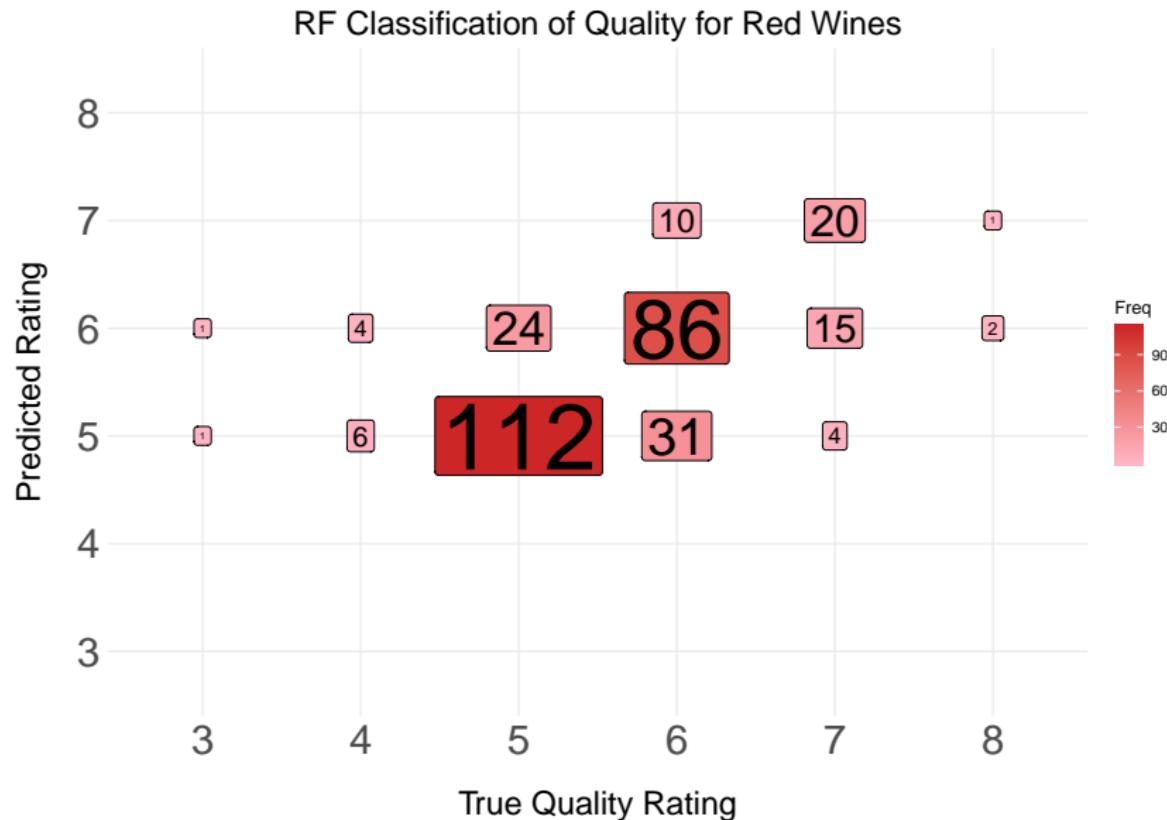
Results: Multinomial Regression - Full Model Confusion Matrix (White Wine)



Results: Random Forests (Red Wine)

- ▶ Accuracy: 68.77%
- ▶ Kappa: 0.4903
- ▶ Weighted Kappa: 0.5873

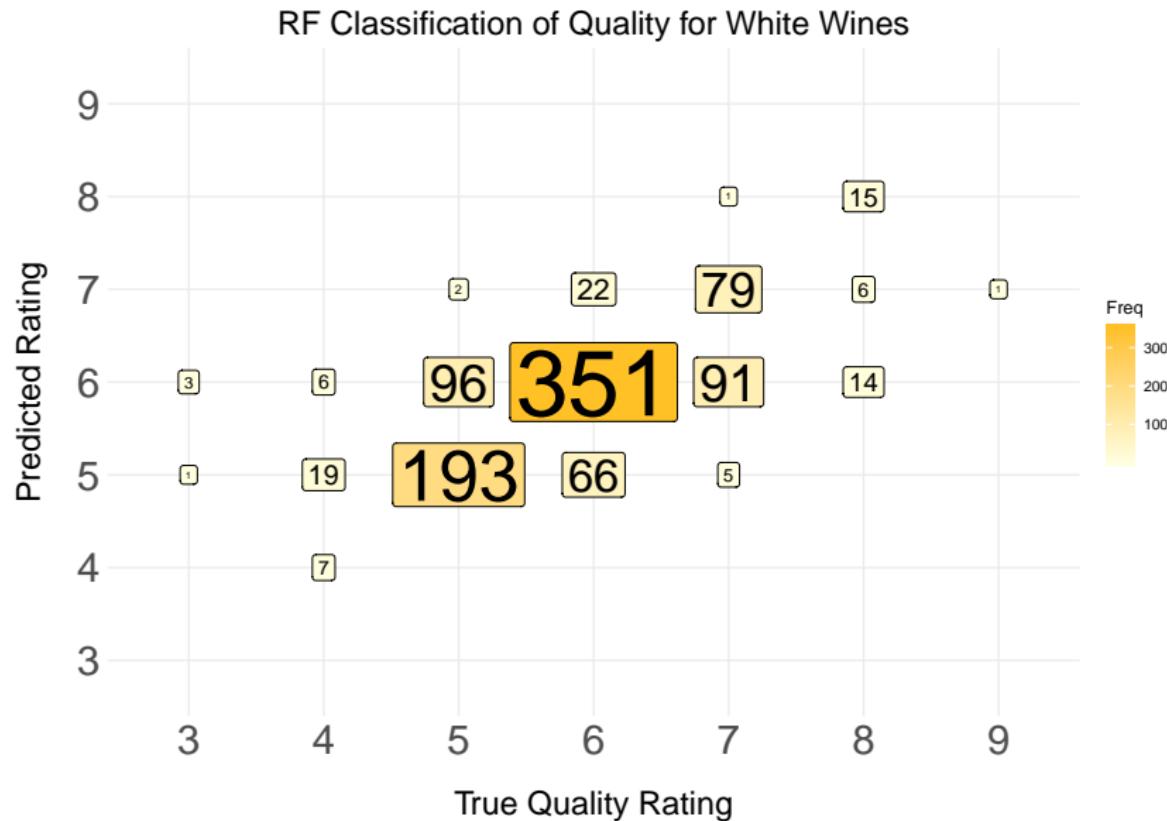
Results: Random Forests (Red Wine)



Results: Random Forests (White Wine)

- ▶ Accuracy: 65.95%
- ▶ Kappa: 0.4638
- ▶ Weighted Kappa: 0.6385

Results: Random Forests (White Wine)



Comparison of Results: Red Wine

Table 4: Comparison of Results for Red Wine

	Overall Results			Percent Correct by Category					
	Prediction Accuracy	Kappa	Weighted Kappa	3	4	5	6	7	8
Random Forest	68.7700	0.4903	0.0000	0	82.35	67.7200	51.2800	0.0000	0.5873
Proportional Odds	58.0442	0.3065	0.4707	0	0.00	72.0588	59.8425	25.6410	0.0000
Partial Proportional Odds	58.3596	0.3117	0.4742	0	0.00	72.0588	59.8425	28.2051	0.0000
Non-Proportional Odds	57.7287	0.3080	0.4357	0	0.00	72.0588	56.6929	33.3333	0.0000
Multinomial	58.3596	0.3158	0.5227	0	20.00	74.2647	55.9055	28.2051	0.0000
Linear Regression	56.1500	0.2728	0.4545	0	0.00	65.4400	63.7800	20.5100	0.0000

Comparison of Results: White Wine

Table 5: Comparison of Results for White Wine

	Overall Results			Percent Correct by Category						
	Prediction Accuracy	Kappa	Weighted Kappa	3	4	5	6	7	8	9
Random Forest	65.9500	0.4638	0.0000	21.88	66.320	79.9500	44.8900	42.8600	0	0.6385
Proportional Odds	51.7382	0.2108	0.3993	0.00	0.000	51.2027	74.9431	15.9091	0	0.0000
Multinomial	54.0900	0.2451	0.4121	0.00	9.375	51.2027	79.4989	15.9091	0	0.0000
Linear Regression	51.1200	0.1969	0.4067	0.00	0.000	41.5800	78.1300	20.4500	0	0.0000

Discussion: Random Forest

- ▶ Using random forests, we were able to predict expert wine quality ratings fairly accurately using measurements of a small number of chemical components of wine
 - ▶ White wine
 - ▶ Accuracy: 65.95%
 - ▶ Weighted Kappa: 0.6385
 - ▶ Red wine
 - ▶ Accuracy: 68.77%
 - ▶ Weighted Kappa: 0.5873
- ▶ Using the proposed cutoffs from Landis and Koch, these weighted Kappa values suggest moderate to substantial agreement with the expert ratings

Discussion: Likelihood Based Approaches

- ▶ The likelihood based approaches performed more poorly, with accuracies 10-15% lower than random forest and weighted kappas 0.1 to 0.2 lower
- ▶ Even so, the weighted kappas all suggested moderate agreement with the expert ratings and were clearly better than random guessing
- ▶ There was not a clear winner among the likelihood based approaches, accuracies and kappas were very similar across multinomial, ordinal, and linear regression models
 - ▶ Accounting for the ordered nature of the ratings did not make a difference in predictions
 - ▶ In some cases, doing so actually resulted in (slightly) worse predictive performance

Discussion: Limitations and Future Directions

- ▶ We only had data on a small number of reasonably easy to measure chemicals, we speculate some misclassification was due to failure to measure other important chemicals.
- ▶ Our goal was simply to predict quality ratings, in the future, it would be interesting characterize how individual chemicals relate to wine quality.
- ▶ There were only a small number of truly excellent wines (rating of 8 or higher) making it particularly difficult to predict these ratings. It would be worthwhile to examine whether various outlier detection algorithms are more suited to characterizing these wines than the approaches we took.

Bottom Line

Expert wine quality ratings can be predicted reasonable well using chemical components, but true wine connoisseurs are still better off consulting a sommelier.

References

Landis JR, Koch GG. The measurement of observer agreement for categorical data. Biometrics. 1977 Mar;33(1):159-74.