

P8106_HW1

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2/16/2025

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Loading libraries

```
library(ISLR)
library(glmnet)
library(caret)
library(tidymodels)
library(corrplot)
library(ggplot2)
library(plotmo)
library(ggrepel)
library(pls)
library(knitr)
```

Question (a): Lasso Model

To start, we load the training and testing data and subsequently set a seed for reproducibility.

Next, we initialise 10-fold cross-validation to partition the training data into 10 equal subsets. This allows training the model on 9 folds while validating on the final fold. This ensures we evaluate the performance of the model, while avoiding overfitting.

```
# Load training and testing data

training_data <- read.csv("housing_training.csv")
testing_data <- read.csv("housing_test.csv")
```

```

set.seed(29) # Ensure results are reproducible

# Using 10 fold cross-validation

ctrl1 <- trainControl(method = "cv", number = 10)

```

Next, we proceed to fit a lasso regression model using the training data. Sale_Price is the outcome variable, with all other variables as predictors. The lasso model is tuned over a sequence of 100 lambda values ranging from $\exp(6)$ to $\exp(-5)$.

```

set.seed(29) # Ensure results are reproducible

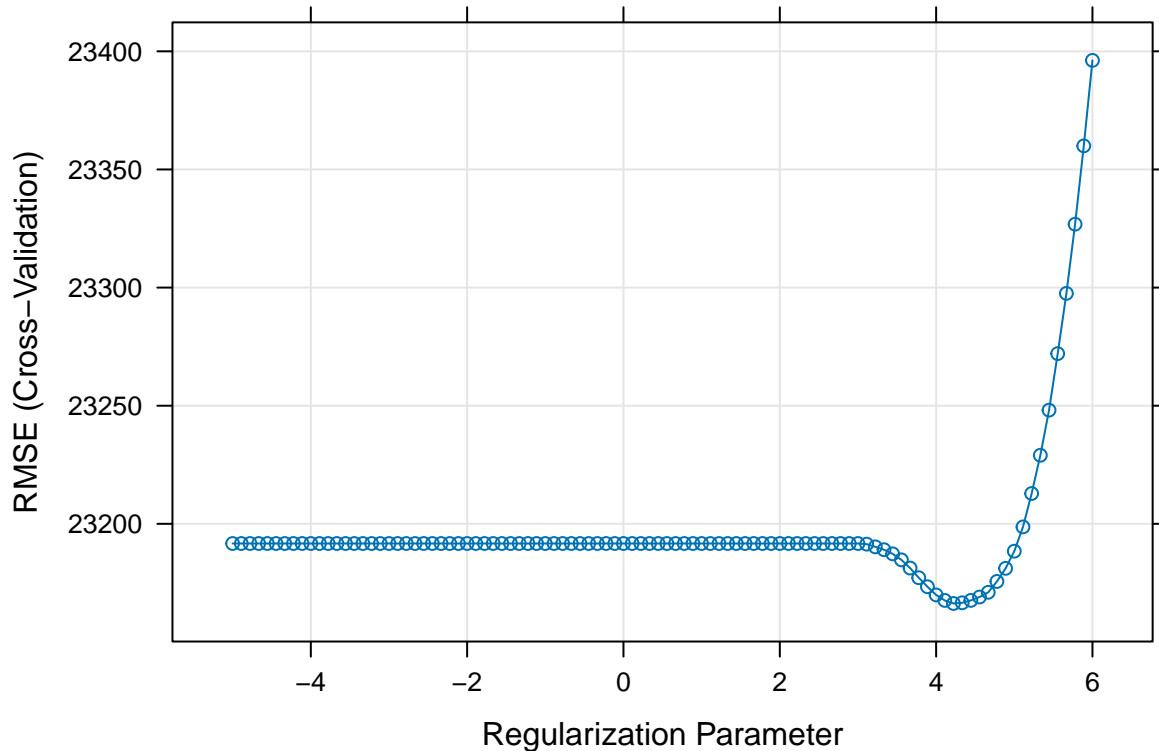
# Fit the Lasso model

lasso.fit <- train(
  Sale_Price ~ .,
  data = training_data,
  method = "glmnet",
  tuneGrid = expand.grid(alpha = 1,
                        lambda = exp(seq(6, -5, length = 100))),
  trControl = ctrl1
)

# Plot

plot(lasso.fit, xTrans = log)

```



Based on the plot, it appears as though the optimal lambda value is around $\exp(4)$, as this is where the RMSE is minimised. Higher lambda values (i.e., greater penalisation) appear to result in poorer model performance, likely due to excessive shrinkage forcing too many coefficients to zero, leading to underfitting.

```
set.seed(29) # Ensure results are reproducible
```

```
# Find optimal tuning parameter
```

```
lasso.fit$bestTune
```

```
##      alpha      lambda
```

```
## 84      1 68.18484
```

```
# Extracting coefficients for each predictor, at the optimal lambda
```

```
coef(lasso.fit$finalModel, lasso.fit$bestTune$lambda)
```

```
## 40 x 1 sparse Matrix of class "dgCMatrix"
```

```
##              s1
## (Intercept) -4.820791e+06
## Gr_Liv_Area  6.534680e+01
## First_Flr_SF 8.043483e-01
## Second_Flr_SF .
## Total_Bsmt_SF 3.542591e+01
## Low_Qual_Fin_SF -4.089879e+01
```

```
## Wood_Deck_SF          1.161853e+01
## Open_Porch_SF        1.539927e+01
## Bsmt_Unf_SF          -2.088675e+01
## Mas_Vnr_Area         1.091770e+01
## Garage_Cars          4.078354e+03
## Garage_Area          8.182394e+00
## Year_Built           3.232484e+02
## TotRms_AbvGrd        -3.607362e+03
## Full_Bath            -3.820746e+03
## Overall_QualAverage   -4.845814e+03
## Overall_QualBelow_Average -1.244202e+04
## Overall_QualExcellent 7.559703e+04
## Overall_QualFair      -1.073410e+04
## Overall_QualGood      1.211373e+04
## Overall_QualVery_Excellent 1.358907e+05
## Overall_QualVery_Good 3.788544e+04
## Kitchen_QualFair      -2.476713e+04
## Kitchen_QualGood      -1.713660e+04
## Kitchen_QualTypical   -2.525278e+04
## Fireplaces           1.051146e+04
## Fireplace_QuFair      -7.657866e+03
## Fireplace_QuGood      .
## Fireplace_QuNo_Fireplace 1.385656e+03
## Fireplace_QuPoor      -5.632703e+03
## Fireplace_QuTypical   -7.010013e+03
## Exter_QualFair        -3.316061e+04
## Exter_QualGood        -1.492745e+04
## Exter_QualTypical     -1.936658e+04
## Lot_Frontage         9.952901e+01
## Lot_Area             6.042265e-01
## Longitude            -3.285809e+04
## Latitude             5.492284e+04
## Misc_Val             8.240622e-01
## Year_Sold            -5.568142e+02
```

Note that at the optimal lambda value, most of the predictors remain in the model. However, some are shrunk to zero (i.e., Second_Flr_SF, Fireplace_QuGood) during the variable selection process, and removed from the model. Therefore, this final model includes **37 predictors**.

```
set.seed(29) # Ensure results are reproducible

# Finding RMSE

lasso_preds <- predict(lasso.fit, newdata = testing_data)

lasso_rmse <- sqrt(mean((lasso_preds - testing_data$Sale_Price)^2))

print(lasso_rmse)
```

```
## [1] 20969.2
```

For the lasso model, the optimal tuning parameter lambda is **68.18484**, representing where RMSE is minimised. The test error (RMSE) at this lambda is **20969.2**.

```

set.seed(29) # Ensure results are reproducible

# Using 1se cross-validation.
# Code from: https://www.rdocumentation.org/packages/caret/versions/6.0-92/topics/oneSE

ctrl_1se <- trainControl(
  method = "cv",
  selectionFunction = "oneSE"
)

# Fit the lasso model using 1se

lasso_1se_fit <- train(
  Sale_Price ~ .,
  data = training_data,
  method = "glmnet",
  tuneGrid = expand.grid(
    alpha = 1,
    lambda = exp(seq(6, -5, length = 100))
  ),
  trControl = ctrl_1se
)

# Optimal lambda using 1SE

lasso_lambda_1se <- lasso_1se_fit$bestTune$lambda
print(lasso_lambda_1se)

## [1] 403.4288

# Extracting coefficients for each predictor, at the optimal lambda

coef(lasso_1se_fit$finalModel, s = lasso_lambda_1se)

## 40 x 1 sparse Matrix of class "dgCMatrix"
##                               s1
## (Intercept)                -3.919159e+06
## Gr_Liv_Area                  6.099153e+01
## First_Flr_SF                 9.477449e-01
## Second_Flr_SF                .
## Total_Bsmt_SF                3.627699e+01
## Low_Qual_Fin_SF              -3.523480e+01
## Wood_Deck_SF                 1.000632e+01
## Open_Porch_SF                1.203918e+01
## Bsmt_Unf_SF                  -2.059528e+01
## Mas_Vnr_Area                 1.297320e+01
## Garage_Cars                  3.491107e+03
## Garage_Area                  9.740129e+00
## Year_Built                   3.150805e+02
## TotRms_AbvGrd                -2.518326e+03
## Full_Bath                    -1.415647e+03
## Overall_QualAverage           -4.006722e+03
## Overall_QualBelow_Average     -1.084480e+04

```

```
## Overall_QualExcellent      8.719850e+04
## Overall_QualFair          -8.763236e+03
## Overall_QualGood           1.111389e+04
## Overall_QualVery_Excellent 1.559964e+05
## Overall_QualVery_Good      3.730815e+04
## Kitchen_QualFair          -1.420320e+04
## Kitchen_QualGood          -7.639239e+03
## Kitchen_QualTypical        -1.644274e+04
## Fireplaces                 8.190689e+03
## Fireplace_QuFair          -3.809974e+03
## Fireplace_QuGood           2.196176e+03
## Fireplace_QuNo_Fireplace    .
## Fireplace_QuPoor          -1.484163e+03
## Fireplace_QuTypical        -4.125304e+03
## Exter_QualFair            -1.695505e+04
## Exter_QualGood            .
## Exter_QualTypical          -4.790664e+03
## Lot_Frontage               8.663344e+01
## Lot_Area                   5.915806e-01
## Longitude                  -2.246220e+04
## Latitude                   3.767830e+04
## Misc_Val                   3.093854e-01
## Year_Sold                  -1.654627e+02
```

```
# Lasso 1SE RMSE
```

```
lasso_1SE_preds <- predict(lasso_1se_fit, newdata = testing_data)

lasso_1SE_rmse <- sqrt(mean((lasso_1SE_preds - testing_data$Sale_Price)^2))

print(lasso_1SE_rmse)
```

```
## [1] 20511.62
```

Using the 1SE rule, the optimal lambda is **403.4288**. During the variable selection process, variables `Second_Flr_SF`, `Fireplace_QuNo_Fireplace`, and `Exter_QualGood` are removed from the model. When the 1SE rule is applied, there are **36 predictors** included in the model, which is 1 fewer than the original lasso model.

Question (b): Elastic Net

To fit the elastic net model, I began with a wide lambda range.

```
# Set seed to ensure reproducibility

set.seed(16)

# Fit elastic net model
# Tuning the different lambda ranges

enet.fit <- train(Sale_Price ~ .,
                  data = training_data,
```

```

method = "glmnet",
tuneGrid = expand.grid(alpha = seq(0, 1, length = 21),
                      lambda = exp(seq(6, -5, length = 100))),
trControl = ctrl1)

# Results

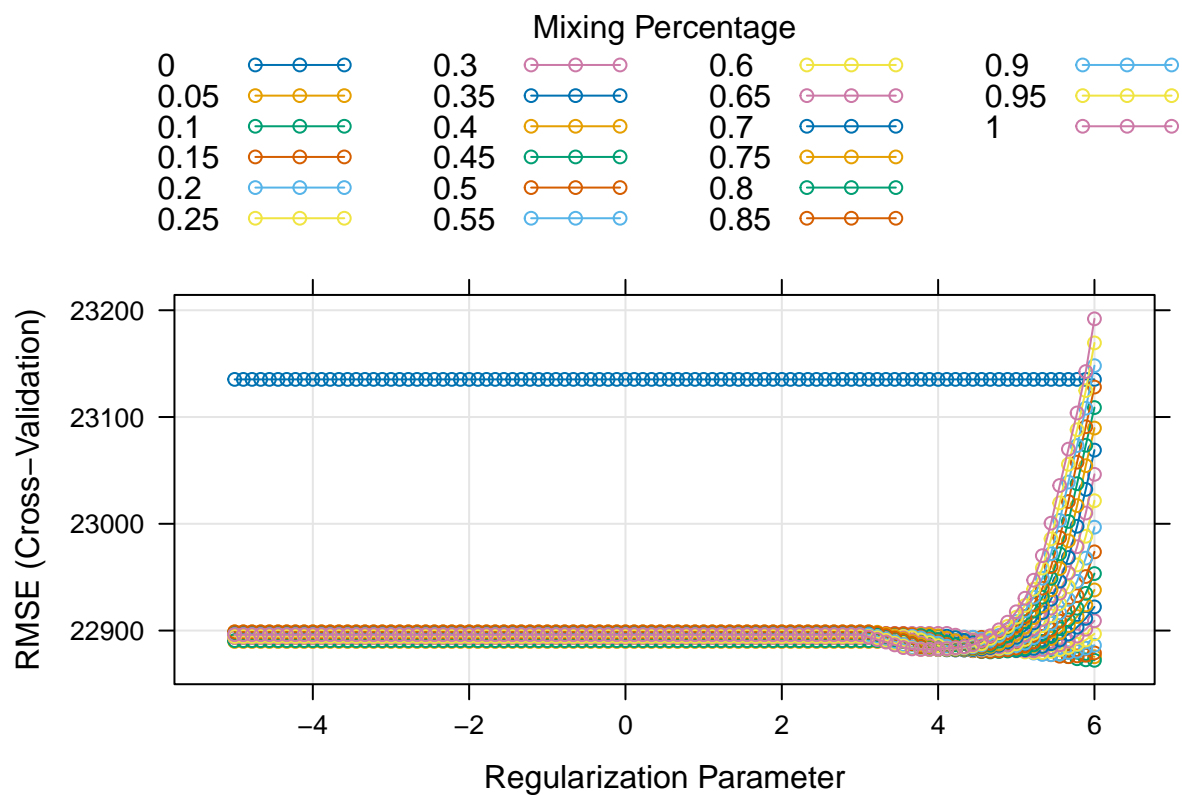
print(enet.fit$bestTune)

##      alpha  lambda
## 300    0.1 403.4288

# Cross validation plot

plot(enet.fit, xTrans = log)

```



After reviewing the cross-validation plot, I refined the lambda range.

```

# Set seed to ensure reproducibility

set.seed(16)

# Adjusting

enet.fit <- train(Sale_Price ~ .,

```

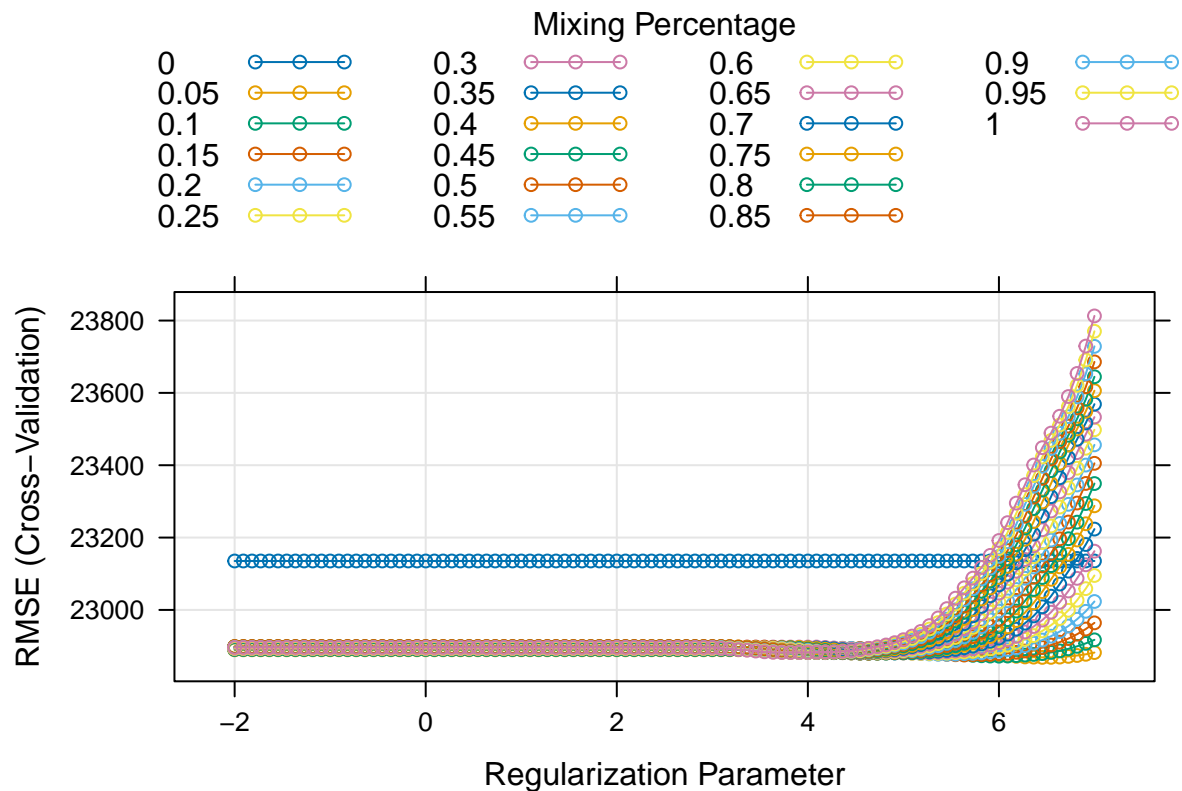
```

data = training_data,
method = "glmnet",
tuneGrid = expand.grid(alpha = seq(0, 1, length = 21),
                      lambda = exp(seq(7, -2, length = 100))),
trControl = ctrl1)

# Cross validation plot

plot(enet.fit, xTrans = log)

```



```

# Optimal lambda

print(enet.fit$bestTune)

```

```

##      alpha  lambda
## 194  0.05 635.5848

```

The cross validation plot shows the RMSE values were fairly stable at lower regularisation values, but increasing steeply when $\log(\lambda) \geq 6$. Therefore, the selected tuning parameters are **alpha = 0.05** and **lambda = 635.5848**.

```

# Set seed to ensure reproducibility

set.seed(16)

```



```

# Predictions using testing dataset

enet.pred <- predict(enet.fit, newdata = testing_data)

# Test error

enet_test_mse <- mean((enet.pred - testing_data$Sale_Price)^2)

# Results

print(enet_test_mse)

```

```
## [1] 438041526
```

From this, the test error of the model is **438041526**.

```

# Set seed to ensure reproducibility

set.seed(16)

# Applying 1SE rule to elastic net model

enet_1se_fit <- train(
  Sale_Price ~ .,
  data = training_data,
  method = "glmnet",
  tuneGrid = expand.grid(
    alpha = seq(0, 1, length = 21),
    lambda = exp(seq(7, -2, length = 100))
  ),
  trControl = ctrl_1se
)

enet_1se_fit$bestTune$lambda

```

```
## [1] 1096.633
```

```
enet_1se_fit$bestTune$alpha
```

```
## [1] 0
```

Yes, it is possible to apply the 1SE rule to selecting tuning parameters for elastic net. The elastic net method includes penalties from both ridge regression and lasso (the mixing parameter alpha that determines the balance between ridge and lasso penalties, and the overall regularisation strength lambda). The 1SE rule is defined as the most regularised model such that error is within one standard error of the minimum.

Therefore, using the 1SE rule, it is possible to select the most regularised model (i.e., the largest lambda) for each alpha value that has error within one standard error of the minimum, then compare across different alpha values to give the effective regularisation via the ridge-type penalty and feature selection via the lasso penalty, as determined by cross-validation.

Based on our data, the 1SE rule model parameters are alpha = **0** and lambda = **1096.633**. Given that alpha = 0, this indicates ridge regression was the optimal model.

I proceeded to find the test error of the 1SE model.

```
# Set seed to ensure reproducibility

set.seed(16)

# Predictions using testing dataset for 1SE model

enet_1se_pred <- predict(enet_1se_fit, newdata = testing_data)

# Test error

enet_1se_test_mse <- mean((enet_1se_pred - testing_data$Sale_Price)^2)

# Results

print(enet_1se_test_mse)
```

```
## [1] 426357707
```

The test error for the 1SE rule elastic net model is **426357707**.

Question (c): Partial least squares

I proceeded with fitting the partial least squares model.

```
# Set seed for reproducibility

set.seed(29)

# Fitting the model

pls_mod <- plsr(Sale_Price ~ .,
                data = training_data,
                scale = TRUE,
                validation = "CV")

summary(pls_mod)
```

```
## Data:      X dimension: 1440 39
## Y dimension: 1440 1
## Fit method: kernelpls
## Number of components considered: 39
##
## VALIDATION: RMSEP
## Cross-validated using 10 random segments.
##      (Intercept)  1 comps  2 comps  3 comps  4 comps  5 comps  6 comps
## CV           73685   33422   27955   25141   24106   23504   23315
## adjCV        73685   33415   27918   25071   24032   23431   23248
##      7 comps  8 comps  9 comps 10 comps 11 comps 12 comps 13 comps
## CV       23193   23167   23191   23193   23182   23178   23185
## adjCV    23131   23106   23126   23126   23115   23111   23117
```

```
##      14 comps 15 comps 16 comps 17 comps 18 comps 19 comps 20 comps
## CV      23187   23198   23201   23207   23211   23231   23237
## adjCV    23119   23129   23132   23138   23142   23160   23165
##      21 comps 22 comps 23 comps 24 comps 25 comps 26 comps 27 comps
## CV      23239   23243   23243   23244   23244   23247   23250
## adjCV    23167   23171   23171   23171   23171   23174   23176
##      28 comps 29 comps 30 comps 31 comps 32 comps 33 comps 34 comps
## CV      23250   23250   23250   23250   23250   23250   23250
## adjCV    23176   23177   23177   23177   23177   23177   23177
##      35 comps 36 comps 37 comps 38 comps 39 comps
## CV      23250   23250   23250   23250   23521
## adjCV    23177   23177   23177   23177   23373
##
## TRAINING: % variance explained
##      1 comps 2 comps 3 comps 4 comps 5 comps 6 comps 7 comps
## X      20.02   25.93   29.67   33.59   37.01   40.03   42.49
## Sale_Price 79.73   86.35   89.36   90.37   90.87   90.99   91.06
##      8 comps 9 comps 10 comps 11 comps 12 comps 13 comps 14 comps
## X      45.53   47.97   50.15   52.01   53.69   55.35   56.86
## Sale_Price 91.08   91.10   91.13   91.15   91.15   91.16   91.16
##      15 comps 16 comps 17 comps 18 comps 19 comps 20 comps
## X      58.64   60.01   62.18   63.87   65.26   67.10
## Sale_Price 91.16   91.16   91.16   91.16   91.16   91.16
##      21 comps 22 comps 23 comps 24 comps 25 comps 26 comps
## X      68.44   70.12   71.72   73.35   75.20   77.27
## Sale_Price 91.16   91.16   91.16   91.16   91.16   91.16
##      27 comps 28 comps 29 comps 30 comps 31 comps 32 comps
## X      78.97   80.10   81.83   83.55   84.39   86.34
## Sale_Price 91.16   91.16   91.16   91.16   91.16   91.16
##      33 comps 34 comps 35 comps 36 comps 37 comps 38 comps
## X      88.63   90.79   92.79   95.45   97.49   100.00
## Sale_Price 91.16   91.16   91.16   91.16   91.16   91.16
##      39 comps
## X      100.64
## Sale_Price 91.04
```

```
# Determine the optimal number of components
```

```
cv_mse <- RMSEP(pls_mod)
```

```
ncomp_cv <- which.min(cv_mse$val[1,]) - 1
```

```
# Optimal number of components
```

```
print(ncomp_cv)
```

```
## 8 comps
```

```
##      8
```

Based on the computation above, the optimal number of components is **8**.

```
# Set seed for reproducibility
```

```

set.seed(29)

# Calculate Test MSE

y2 <- testing_data$Sale_Price

predy2_pls <- predict(pls_mod, newdata = testing_data,
                      ncomp = ncomp_cv)

pls_test_mse <- mean((y2 - predy2_pls)^2)

print(pls_test_mse)

```

```
## [1] 440217938
```

The test error for this model is **440217938**.

Question (d): Choose the best model for predicting the response and explain your choice.

```

# Comparison table of models
# Code from: https://bookdown.org/yihui/rmarkdown-cookbook/kable.html

# Convert MSE to RMSE for comparison
enet_test_rmse <- sqrt(enet_test_mse)
pls_test_rmse <- sqrt(pls_test_mse)
enet_1se_test_rmse <- sqrt(enet_1se_test_mse)

comparison_table <- tibble(
  Model = c("Lasso", "Lasso 1SE", "Elastic Net", "Elastic Net 1SE", "Partial Least Square Regression"),
  Test_Error = c(lasso_rmse, lasso_1se_rmse, enet_test_rmse, enet_1se_test_rmse, pls_test_rmse)
)

# Using kable to present table

knitr::kable(comparison_table)

```

| Model | Test_Error |
|---------------------------------|------------|
| Lasso | 20969.20 |
| Lasso 1SE | 20511.62 |
| Elastic Net | 20929.44 |
| Elastic Net 1SE | 20648.43 |
| Partial Least Square Regression | 20981.37 |

Based on the comparison table, the lasso model with the 1SE rule applied appears to perform the best as it has the lowest test error (RMSE = 20511.62) while undertaking feature selection. As discussed in class, the 1SE rule chooses the most regularised model whose error is within one standard error of the minimum,

which allows for a parsimonious model that may generalise better to new data while maintaining predictive performance.

In the case of predicting the housing sale price, the lasso with 1SE model removed various less important predictors during variable selection, which reduced the overall complexity of the model while keeping good predictive performance. This aligns with the bias-variance tradeoff, in which with increasing regularisation (λ), bias tends to increase but variance decreases, leading to improved generalisation.

The better performance of the lasso with 1SE compared to methods like elastic net and partial least squares regression suggests that this method is suitable as it provided the optimal balance in the bias-variance tradeoff for this housing price prediction task. This would be plausible since the lasso performs particularly well when there is a subset of true coefficients that are small or exactly zero, as appears to be the case with some of the housing price predictors.

Question (e): Retrain model using glmnet

```
# Set seed for reproducibility

set.seed(29)

# Matrix of training data predictors for glmnet

x.train <- model.matrix(Sale_Price ~ ., training_data)[,-1]

y.train <- training_data$Sale_Price

# Matrix of predictors for test data

x.test <- model.matrix(Sale_Price ~ ., testing_data)[,-1]

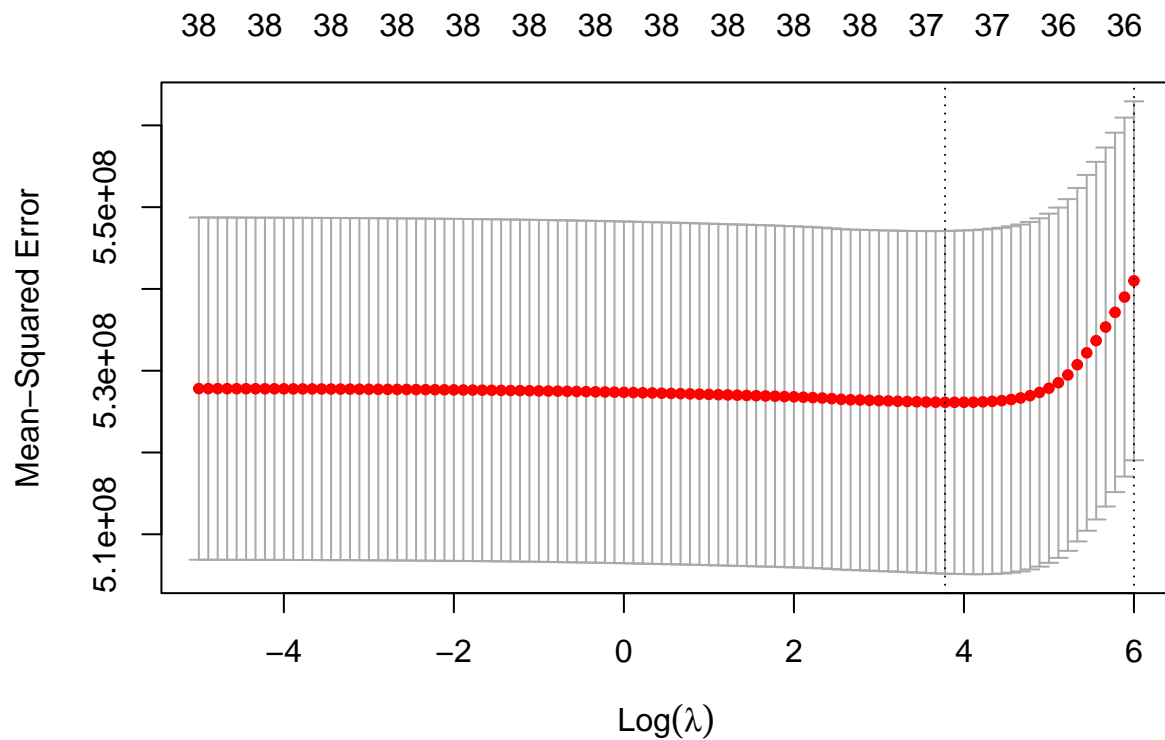
# Fit lasso

lasso_glmnet <- glmnet(x.train, y.train,
                      alpha = 1,
                      lambda = exp(seq(6, -5, length = 100)))

# Next, cross-validation for optimal lambda

cv.lasso <- cv.glmnet(x.train, y.train,
                    alpha = 1,
                    lambda = exp(seq(6, -5, length = 100)))

plot(cv.lasso)
```



```
# Best lambda that minimises RMSE
```

```
print(cv.lasso$lambda.min)
```

```
## [1] 43.71878
```

```
# Coefficients at optimal lambda
```

```
predict(cv.lasso, s = "lambda.min", type = "coefficients")
```

```
## 40 x 1 sparse Matrix of class "dgCMatrix"
##               lambda.min
## (Intercept)   -4.895849e+06
## Gr_Liv_Area    6.563920e+01
## First_Flr_SF   7.866606e-01
## Second_Flr_SF      .
## Total_Bsmt_SF   3.534723e+01
## Low_Qual_Fin_SF -4.125144e+01
## Wood_Deck_SF    1.175170e+01
## Open_Porch_SF   1.563393e+01
## Bsmt_Unf_SF     -2.089478e+01
## Mas_Vnr_Area    1.075014e+01
## Garage_Cars     4.109831e+03
## Garage_Area     8.092846e+00
## Year_Built      3.235015e+02
```

```

## TotRms_AbvGrd          -3.679742e+03
## Full_Bath              -3.994482e+03
## Overall_QualAverage    -4.909818e+03
## Overall_QualBelow_Average -1.257695e+04
## Overall_QualExcellent   7.464381e+04
## Overall_QualFair        -1.091948e+04
## Overall_QualGood         1.216881e+04
## Overall_QualVery_Excellent 1.341508e+05
## Overall_QualVery_Good    3.792195e+04
## Kitchen_QualFair        -2.550193e+04
## Kitchen_QualGood        -1.779096e+04
## Kitchen_QualTypical     -2.585294e+04
## Fireplaces              1.082636e+04
## Fireplace_QuFair        -7.714837e+03
## Fireplace_QuGood         .
## Fireplace_QuNo_Fireplace 1.883694e+03
## Fireplace_QuPoor        -5.696096e+03
## Fireplace_QuTypical     -7.002341e+03
## Exter_QualFair          -3.478720e+04
## Exter_QualGood          -1.640951e+04
## Exter_QualTypical       -2.086944e+04
## Lot_Frontage            1.005252e+02
## Lot_Area                 6.046880e-01
## Longitude               -3.364932e+04
## Latitude                 5.645067e+04
## Misc_Val                 8.584413e-01
## Year_Sold                -5.875701e+02

```

The final model includes 37 predictors, which is the same as the number of predictors identified in (a) using lasso (caret method).

Compared to the caret method implemented in Question (a), the tuning parameter lambda is notably different (glmnet = 43.71878, and caret = 68.18484). Both methods do use 10-fold cross validation to find the optimal lambda.

However, in glmnet, the built in cross validation function `cv.glmnet()` performs 10-fold cross validation once, computing RMSE for each lambda based on the 10 validation sets (Hastie et al., 2024). Whereas the caret package's `trainControl(method = "cv", number = 10)` also applies 10-fold cross validation, but averages the RMSE obtained across multiple resampling instances (Kuhn, 2020).

References

- Kuhn, M. (2020). Caret package documentation.
- Hastie, T., Tibshirani, R., & Friedman, J. (2024). An Introduction to glmnet.