

Project-583

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```
data <- read.csv("data_cleaned.csv", header = TRUE, check.names = FALSE)
dim(data) # 189, 31
data_no_date <- data[, -1]
head(data)
```

We aim to predict the price of ethylene glycol using the following 29 explanatory variables.
MEG refers to ethylene glycol.

y MEG spot price

Up-stream price of MEG

x1 WTI: West Texas Intermediate(Crude Oil)Price

x2 Brent: Brent Crude oil price

x3 Coal price

Up-stream(Ethylene)Profit

x4 domestic; x5 foreign

MEG Profit

x6 made of coal; x7 made of Ethylene

MEG operating rate

x8 domestic; x9 foreign

Downstream profits

x10 Recycled Bottle Chips; x11 polyester chips; x12 polyester bottle chip; x13 POY; x14 FDY; x15 DTY;

x16 Polyester **Downstream operating rate**

x17 Polyester; x18 filament; x19 Direct spinning; x20 chip spun filament; x21 texturing machine; x22 weaving machine

Downstream Inventory

x24 Polyester; x25 FDY; x26 DTY; x27 POY

MEG Inventory

x28 foreign Port; x29 domestic Port; x30 factory

1. A statistically descriptive analysis of the dataset.

The data structure is shown below. We can see that all the variables are continuous values.

```
str(data)

## 'data.frame':   189 obs. of  31 variables:
## $ date: chr  "2019/1/4" "2019/1/11" "2019/1/18" "2019/1/25" ...
## $ y : int  5160 5115 4985 5070 5110 5010 4980 5125 5230 5130 ...
## $ x1 : num  48 51.6 53.8 53.7 56.3 ...
## $ x2 : num  57.1 60.5 62.7 61.6 62.8 ...
## $ x3 : num  570 565 570 565 570 570 575 600 640 628 ...
## $ x4 : num  345.2 28.3 85.5 246.8 364 ...
## $ x5 : num  29.14 4.46 3.8 4.84 26.26 ...
## $ x6 : num  1300 1227 1078 1166 1198 ...
```

```
## $ x7 : num 241 234 -270 -573 -683 ...
## $ x8 : num 0.714 0.729 0.761 0.77 0.801 ...
## $ x9 : num 0.998 0.982 0.941 0.963 0.957 ...
## $ x10 : num -219.8 -155.2 -155.2 60.3 60.3 ...
## $ x11 : num 213.8 175.6 172.2 218.4 98.1 ...
## $ x12 : num 68.8 275.6 322.1 378.4 323.1 ...
## $ x13 : num -81.2 -154.4 -137.8 108.4 -11.9 ...
## $ x14 : num 479 396 442 628 508 ...
## $ x15 : int 490 480 470 460 460 470 450 590 215 250 ...
## $ x16 : num 689 681 677 698 578 ...
## $ x17 : num 84.8 83.9 81.9 78.3 74 78.6 85.4 86.6 88.5 90.5 ...
## $ x18 : num 75.7 75.1 71.7 62.9 62.1 62.1 71.7 75.7 80.1 83.6 ...
## $ x19 : num 82.3 81.8 78 71.7 67.5 73 82.7 84.5 86.7 88.9 ...
## $ x20 : num 48.8 48.8 47 29 29 29 29 84 54 60 ...
## $ x21 : int 76 72 66 17 10 20 50 81 87 89 ...
## $ x22 : int 65 57 50 9 5 34 45 75 84 85 ...
## $ x24 : num 4.7 0.2 1.4 0 2 5.9 7.6 6 5.2 6.4 ...
## $ x25 : num 15.6 12.5 11.1 9.2 11.1 14.2 16.8 18 11.5 14.2 ...
## $ x26 : num 16.7 13.5 13.8 13.8 18 27.2 27.8 28.6 22.9 24.2 ...
## $ x27 : num 13.1 6.7 6.5 3.8 5.5 11.6 14.8 15.4 9.8 10.3 ...
## $ x28 : int 620 624 606 618 627 608 610 629 650 641 ...
## $ x29 : num 78.5 83.4 90.2 89.3 92.4 ...
## $ x30 : num 15.1 14 13.5 14.8 17.8 16.5 15.4 15.4 14.8 14.6 ...
```

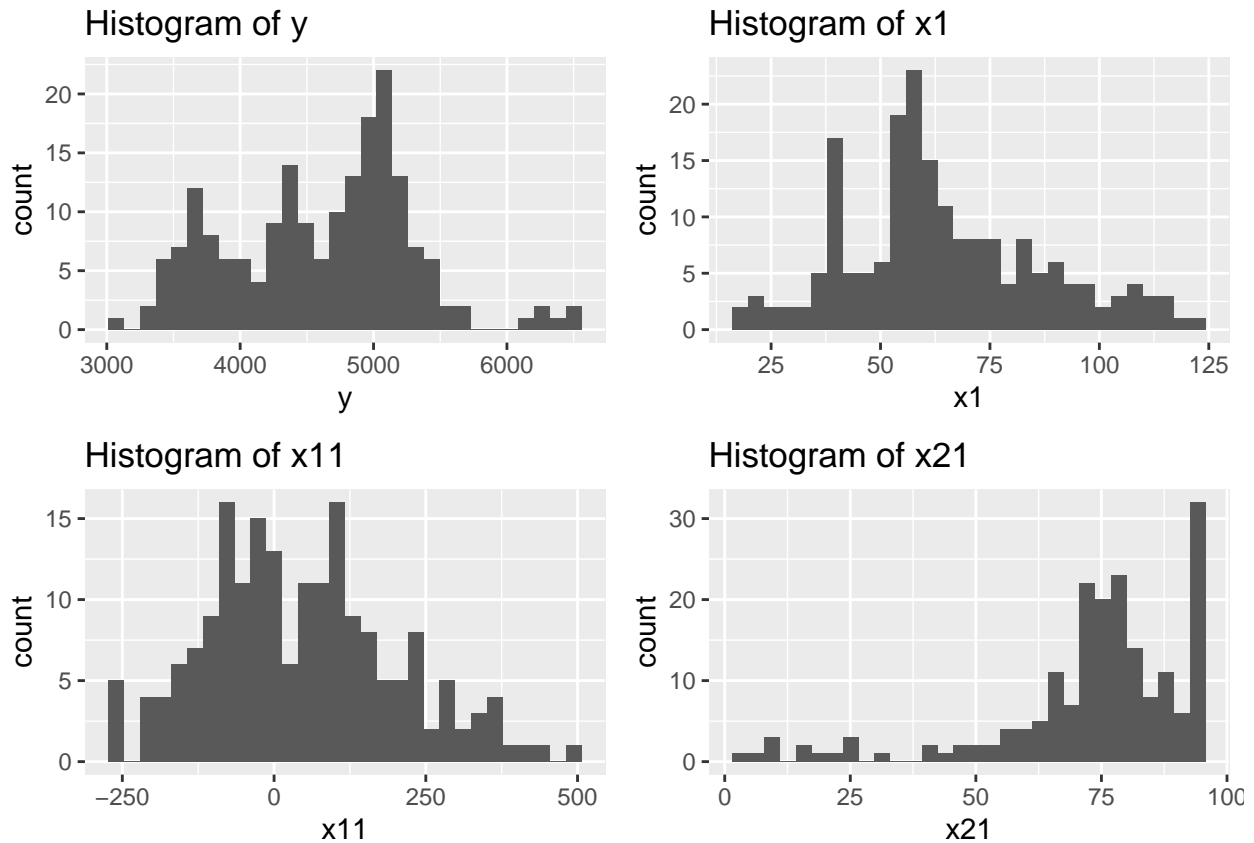
Here is the summary statistics of the response variable.

```
summary(data$y)
```

```
##      Min. 1st Qu.  Median    Mean 3rd Qu.    Max.
##      3130   4065    4705    4612   5110    6557
```

Let's explore the histogram of the response variable y, and some explanatory variables.

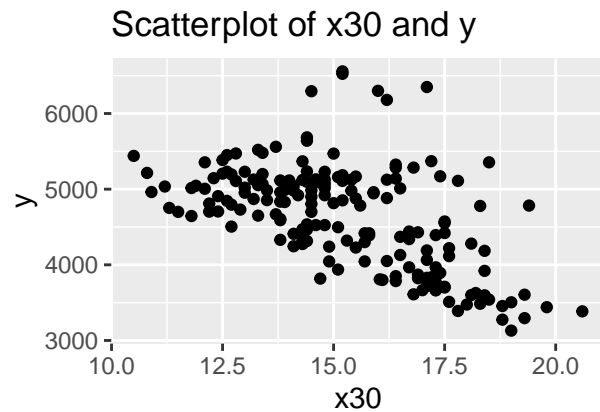
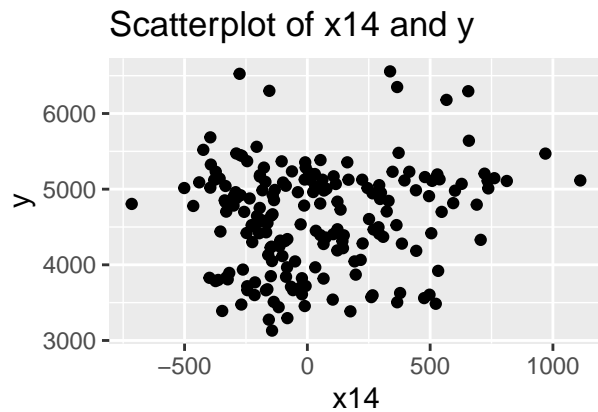
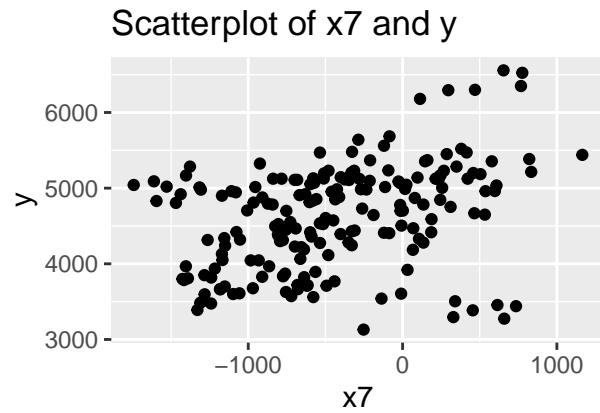
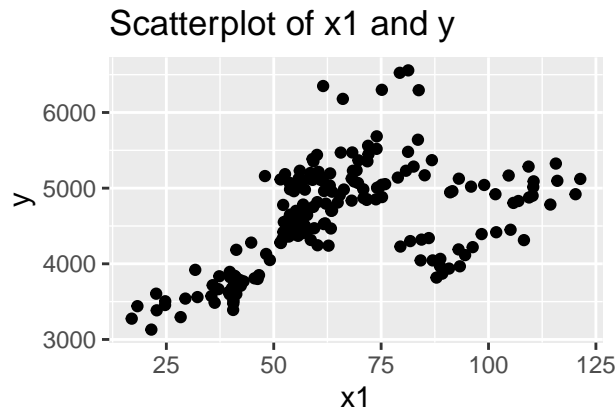
```
hist_y <- ggplot(data, aes(x = y)) + geom_histogram() + labs(title = "Histogram of y")
hist_x1 <- ggplot(data, aes(x = x1)) + geom_histogram() + labs(title = "Histogram of x1")
hist_x11 <- ggplot(data, aes(x = x11)) + geom_histogram() + labs(title = "Histogram of x11")
hist_x21 <- ggplot(data, aes(x = x21)) + geom_histogram() + labs(title = "Histogram of x21")
grid.arrange(hist_y, hist_x1, hist_x11, hist_x21, nrow = 2, ncol = 2)
```



The histogram of the response variable y seems to have 3 peaks. The histogram of the explanatory variable x_1 and x_{11} are right skewed. The histogram of the explanatory variable x_{21} is left skewed.

Let's explore the scatter plots of the response variable y and some explanatory variables.

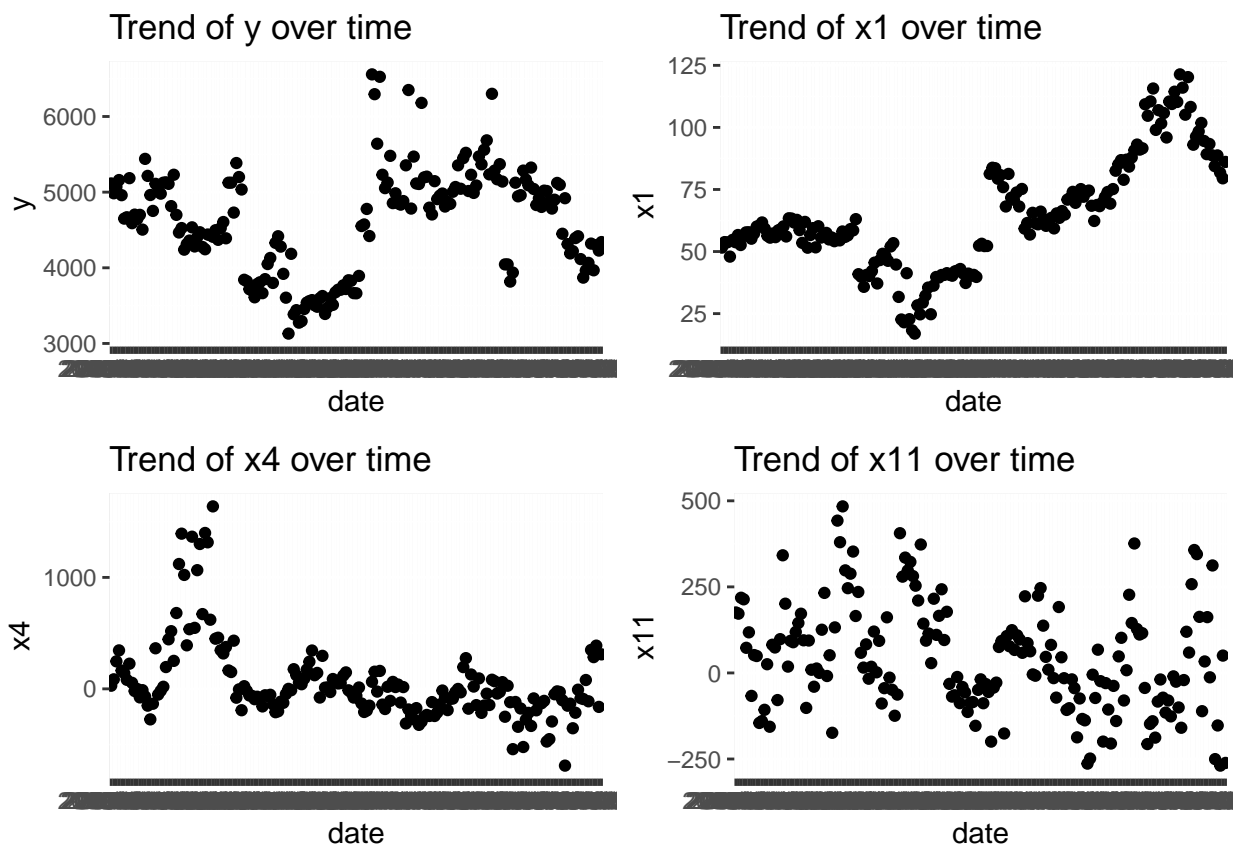
```
# create scatterplots of the variables x1, x2, x3, x4, x5, x7, x10, x11, x12
scatter_x1 <- ggplot(data = data, aes(x = x1, y = y)) + geom_point() + labs(title = "Scatterplot of x1 & y")
scatter_x7 <- ggplot(data = data, aes(x = x7, y = y)) + geom_point() + labs(title = "Scatterplot of x7 & y")
scatter_x14 <- ggplot(data = data, aes(x = x14, y = y)) + geom_point() + labs(title = "Scatterplot of x14 & y")
scatter_x30 <- ggplot(data = data, aes(x = x30, y = y)) + geom_point() + labs(title = "Scatterplot of x30 & y")
grid.arrange(scatter_x1, scatter_x7, scatter_x14, scatter_x30, ncol = 2, nrow = 2)
```



There seems to have a positive linear relationship between the response variable y and the explanatory variables x1, x7. And a negative relationship between y and x30. It seems reasonable because x1 represent Crude Oil Price, which is positively related to the price of y(MEG), while x30 represents the factory inventory, the higher the inventory, the lower the price. There is no obvious relationship between y and x14.

Let's explore the trend of the response variable y and some explanatory variables over time.

```
# plot the variables versus date
trend_y <- ggplot(data = data, aes(x = date, y = y)) + geom_point() + labs(title = "Trend of y over time")
trend_x1 <- ggplot(data = data, aes(x = date, y = x1)) + geom_point() + labs(title = "Trend of x1 over time")
trend_x4 <- ggplot(data = data, aes(x = date, y = x4)) + geom_point() + labs(title = "Trend of x4 over time")
trend_x11 <- ggplot(data = data, aes(x = date, y = x11)) + geom_point() + labs(title = "Trend of x11 over time")
grid.arrange(trend_y, trend_x1, trend_x4, trend_x11, ncol = 2, nrow = 2)
```



The trend plots of y and x1 seem to have some similarities. The trend plots of x4 and x11 have no obvious trend.

2. Applications of statistical analysis techniques

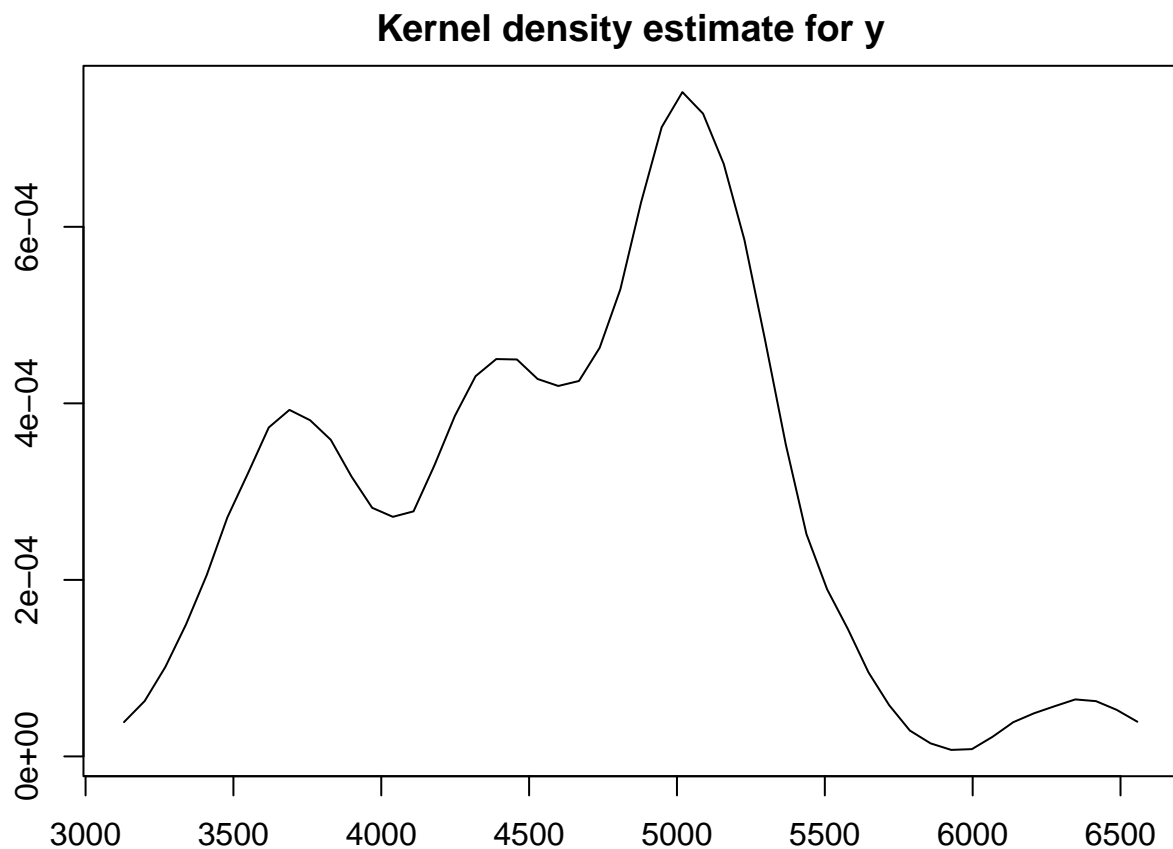
kernel density

Let's try to plot a kernel density estimate for y using an Epanechnikov kernel.

```
par(mar = c(2, 2, 2, 2))
bw <- npudensbw( ~ y, data = data, ckertype = "epanechnikov", bwmethod = "cv.ml")
```

```
## Multistart 1 of 1 |Multistart 1 of 1 |Multistart 1 of 1 |Multistart 1 of 1 /Multistart 1 of 1 |Multi
```

```
fhat <- npudens(bws = bw)
plot(fhat, main = "Kernel density estimate for y")
```



As stated in the histogram part, the distribution of y is not normal, and it seems to have 3 peaks.

Let's conduct Pearson tests for normality.

```
pearson.test(data$y)
```

```
##
##  Pearson chi-square normality test
##
## data:  data$y
## P = 44.772, p-value = 4.441e-05
```

From the pearson chi-square normality tests, we can see that the p-value is very small for y , so we have evidence that y is not normally distributed

Linear regression

```
model_ls <- lm(y ~ ., data = data_no_date)
summary(model_ls)
```

```
##
## Call:
## lm(formula = y ~ ., data = data_no_date)
##
## Residuals:
##      Min       1Q   Median       3Q      Max
## -165.392  -52.794   -1.755    44.694   171.121
##
## Coefficients:
```

```
##           Estimate Std. Error t value Pr(>|t|)
## (Intercept)  1.981e+03  3.006e+02   6.590 6.14e-10 ***
## x1          -7.209e+00  4.750e+00  -1.518 0.131105
## x2           9.303e+00  4.693e+00   1.982 0.049149 *
## x3           1.185e+00  1.305e-01   9.076 4.13e-16 ***
## x4           3.944e-02  5.211e-02   0.757 0.450279
## x5           7.938e-02  5.962e-01   0.133 0.894244
## x6           4.855e-01  5.288e-02   9.181 < 2e-16 ***
## x7           9.900e-02  1.654e-02   5.985 1.38e-08 ***
## x8          -1.227e+02  1.491e+02  -0.823 0.411734
## x9           2.112e+02  1.421e+02   1.486 0.139360
## x10          -9.404e-02  4.265e-02  -2.205 0.028886 *
## x11           1.536e-01  8.178e-02   1.878 0.062259 .
## x12           8.932e-03  2.857e-02   0.313 0.754965
## x13          -1.344e-01  4.317e-02  -3.115 0.002185 **
## x14           1.451e-01  4.062e-02   3.571 0.000471 ***
## x15          -6.469e-02  5.319e-02  -1.216 0.225689
## x16          -1.069e-01  5.485e-02  -1.950 0.052972 .
## x17          -3.266e+00  4.925e+00  -0.663 0.508218
## x18           1.187e-01  1.250e-01   0.950 0.343681
## x19           6.022e+00  4.419e+00   1.363 0.174922
## x20          -3.370e+00  1.397e+00  -2.412 0.016986 *
## x21          -2.405e+00  1.211e+00  -1.986 0.048775 *
## x22           3.065e+00  1.164e+00   2.634 0.009275 **
## x24           5.010e+00  3.501e+00   1.431 0.154350
## x25          -1.115e+01  4.879e+00  -2.285 0.023624 *
## x26          -3.468e+00  3.101e+00  -1.118 0.265201
## x27           1.080e+01  3.528e+00   3.060 0.002597 **
## x28           3.228e+00  4.014e-01   8.043 1.90e-13 ***
## x29          -5.339e-01  6.139e-01  -0.870 0.385811
## x30          -2.251e+01  5.826e+00  -3.864 0.000162 ***
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 74.85 on 159 degrees of freedom
## Multiple R-squared:  0.99, Adjusted R-squared:  0.9882
## F-statistic: 542.8 on 29 and 159 DF, p-value: < 2.2e-16
```

From the summary output, we can see that only variables x3, x6, x7, x13, x14, x20, x22, x25, x27, x28 and x30 are significant. The R-squared and adjusted R-square are high, which is expected as we have lots of variables. Variable selection is needed to reduce the number of variables.

Let's try to do variable selection using stepwise regression.

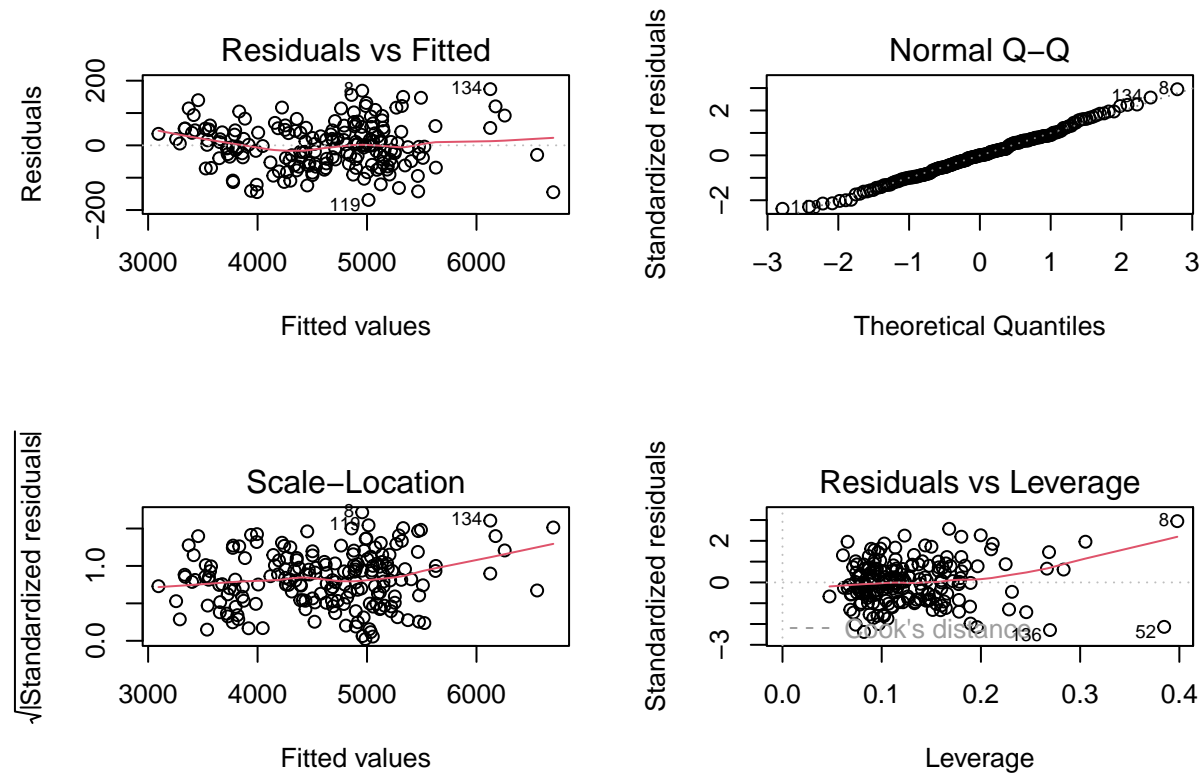
```
model_step <- stepAIC(model_ls, direction = "both", criterion = "bic", trace = FALSE)
model_step
```

```
##
## Call:
## lm(formula = y ~ x1 + x2 + x3 + x4 + x6 + x7 + x9 + x10 + x11 +
##       x13 + x14 + x15 + x16 + x19 + x20 + x21 + x22 + x24 + x25 +
##       x26 + x27 + x28 + x30, data = data_no_date)
##
## Coefficients:
## (Intercept)          x1          x2          x3          x4          x6
## 1858.80540      -9.16272     11.29471     1.22914     0.04681     0.49736
```

##	x7	x9	x10	x11	x13	x14
##	0.10436	217.80130	-0.11073	0.18016	-0.11633	0.12868
##	x15	x16	x19	x20	x21	x22
##	-0.06754	-0.14431	3.00727	-3.38212	-1.94951	2.42204
##	x24	x25	x26	x27	x28	x30
##	4.01386	-8.11560	-5.01568	10.32603	3.15784	-24.67567

From the stepwise model selection, 6 variables are removed from the model.

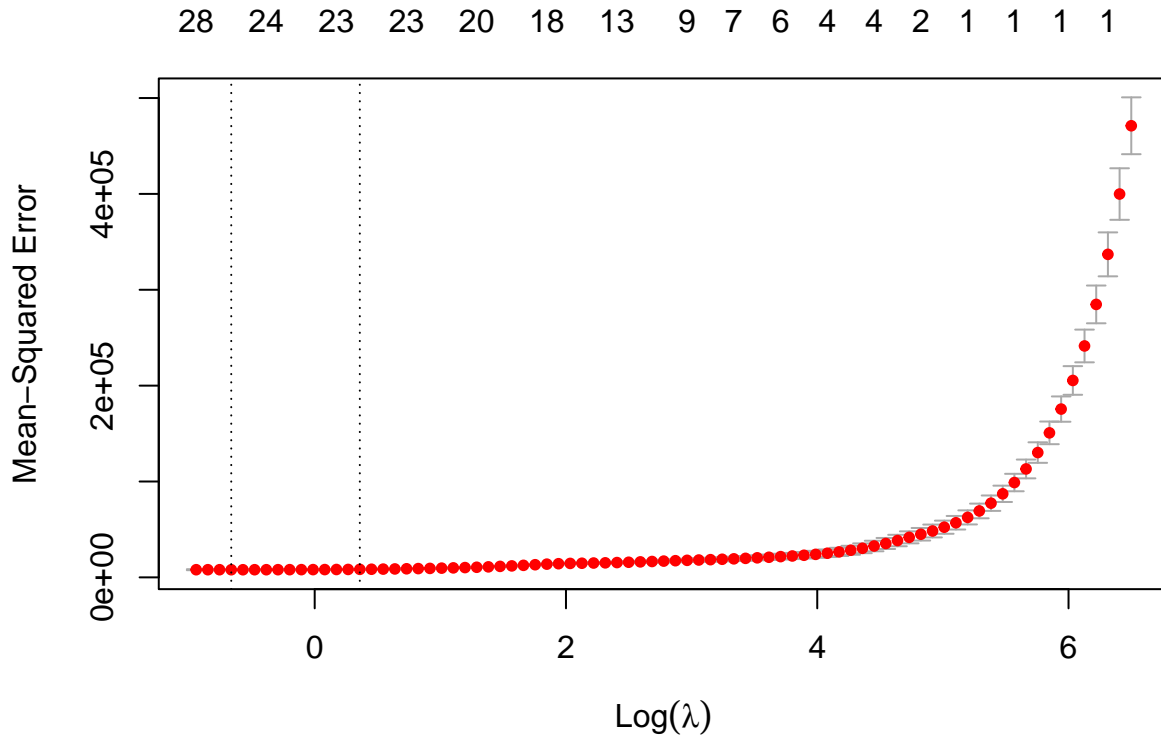
```
par(mfrow=c(2,2))
plot(model_step)
```



The residual plot shows no obvious trend. The scale-location plot shows a slightly increasing trend. The normal Q-Q plot shows that the residuals are roughly normally distributed. The leverage plot shows that there are couple potential outliers.

Lasso

```
knitr::opts_chunk$set(fig.width=5, fig.height=4)
model_lasso <- cv.glmnet(x = as.matrix(data_no_date[, -1]), y = data_no_date[, 1], alpha = 1)
plot(model_lasso)
```

Both minimum and 1se lines suggest to keep 23 variables, which is the same as the stepwise model selection.

3. Scientific questions.

4. Statistical analysis techniques I will use to answer those questions.

Q1. Based on our dataset, which models can be used to forecast the price of ethylene glycol?

From the above analysis, linear regression model seems to be good in interpretation. R-square is 0.99 and many variables are significant. However, there is a potential over-fitting problem with so many variables included in the model. Also, as response variable is not normal distributed, it may violate the assumptions of linear regression. Therefore, we can use other methods such as **nonparametric local linear regression model or regression trees or random forest**.

We also can consider using the **logistic regression**, but we need to convert the continuous response variable into a binary outcome by applying a threshold value. For example, if the response variable increases compared with last day or last week, then we assign it to 1, otherwise, we assign it to 0.

Nonparametric local linear regression model

```
library(np)
bw <- npregbw(y ~ x1 + x2 + x3 + x4 + x6 + x7 + x9 + x10 + x11 + x13 + x14 + x15 + x16 + x19 + x20 + x21)

## Multistart 1 of 5 |Multistart 1 of 5 |Multistart 1 of 5 |Multistart 1 of 5 /Multistart 1 of 5 |Multistart 1 of 5 |Multistart 1 of 5
model.ll <- npreg(bws=bw)
summary(model.ll)

##
## Regression Data: 189 training points, in 23 variable(s)
##           x1      x2      x3      x4      x6      x7      x9
## Bandwidth(s): 127.5724 10.17717 322.5403 2537.315 504.7473 336.0572 0.07111343
##           x10      x11      x13      x14      x15      x16      x19
```

```
## Bandwidth(s): 787.6651 310.6759 2438.71 4392.377 252.3407 329.8475 4.534213
##           x20      x21      x22      x24      x25      x26      x27
## Bandwidth(s): 17.95935 27.15539 177.2888 42.59719 45.24559 21.4809 28.94301
##           x28      x30
## Bandwidth(s): 492.8228 0.6055561
##
## Kernel Regression Estimator: Local-Linear
## Bandwidth Type: Fixed
## Residual standard error: 0.5214158
## R-squared: 0.9999994
##
## Continuous Kernel Type: Second-Order Gaussian
## No. Continuous Explanatory Vars.: 23
```

In nonparametric local linear regression model, R square is quite close to 1 (larger than R square in linear regression), which indicates that the model is fitting the data better than linear regression.

Logistic

```
data_logistic <- data_no_date[-1,]
data_logistic$newy <- ifelse(diff(data_no_date$y) > 0, 1, 0)
data_logistic$newy <- as.factor(data_logistic$newy)
model_Log <- glm(newy ~., data=data_logistic[,-1], family = binomial)
model_Log_step <- stepAIC(model_Log, direction = "both", criterion = "bic", trace = FALSE)
```

The variable selection helped us to reduce the number of variables (from 29 to 14), and we have a lower BIC value (from 229.3 to 206.2); Due to space limitations, we do not show results.

Q2. There are 29 variables in our dataset, how can we reduce the dimension and avoid multicollinearity? There are 2 methods we can consider, the first method is lasso, as used above, it can be used to reduce high-dimensional data in a model by shrinking the coefficients of irrelevant variables to zero. The other method is PCA, which aims to reduce the dimensionality of the data while retaining as much of its variance as possible.

PCA

```
pca <- prcomp(data_no_date[, -c(1)], scale.=TRUE)
summary(pca)

var_explained <- cumsum(pca $sdev^2 / sum(pca $sdev^2))
# Determine the number of components needed to retain at least 90% of the variance
which.max(var_explained >= 0.9)

## [1] 11
```

We reduce 29 variables to 11 principal components by PCA (those can explain 90% of variance).

Q3. which model is the most appropriate and accurate?

First of all, it depends on our purpose. If our priority is to interpret, linear regression, logistic regression or trees will be better. This involves examining the coefficients of the model to determine the direction and strength of the relationships between the variables. This information can be used to identify the most important variables and to generate hypotheses about the underlying mechanisms driving the relationship between the variables.

If we focus more on predication, we can divide our data into training and testing datasets, using cross validation to test the model with the smallest test MSE. Usually, random forest or boosting would be better in this case. As the data is time series, we can also use ARIMA model to do prediction.