Applied Economic Forecasting

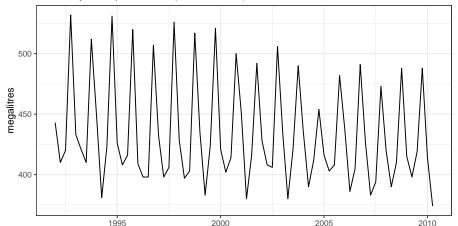
3. Evaluation of Basic Forecasting Models

- Some simple forecasting methods
- 2 Transformations & Adjustments
- 3 Residual diagnostics
- 4 Evaluating forecast accuracy
- **6** Prediction intervals

Section 1

Some simple forecasting methods

Quarterly beer production (1992 – 2010)

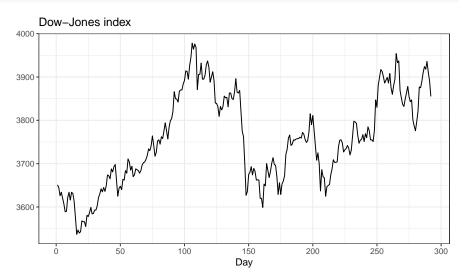


Year

Number of pigs slaughtered in Victoria



```
autoplot(dj) +
labs(title = "Dow-Jones index", x = "Day", y = "")
```



1. Average method

- Forecast of all future values is equal to mean of historical data $\{y_1, \ldots, y_T\}$.
- Forecasts: $\hat{y}_{T+h|T} = \bar{y} = (y_1 + \dots + y_T)/T$

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2. Naïve method

- Forecasts equal to last observed value.
- Forecasts: $\hat{y}_{T+h|T} = y_T$.
- Consequence of efficient market hypothesis.

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2. Naïve method

- Forecasts equal to last observed value.
- Forecasts: $\hat{y}_{T+h|T} = y_T$.
- Consequence of efficient market hypothesis.

3. Seasonal naïve method

- Forecasts equal to last value from same season.
- Forecasts: $\hat{y}_{T+h|T} = y_{T+h-m(k+1)}$, where m = seasonal period and k is the integer part of (h-1)/m.

4. Drift method

- Forecasts equal to last value plus average change.
- Forecasts:

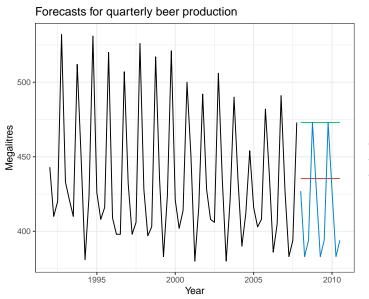
$$\hat{y}_{T+h|T} = y_T + \frac{h}{T-1} \sum_{t=2}^{T} (y_t - y_{t-1})$$
$$= y_T + \frac{h}{T-1} (y_T - y_1).$$

• Equivalent to extrapolating a line drawn between first and last observations.

Functions in R:

- Mean: meanf(y, h=__)
- Naïve: naive(y, h=__)
- Seasonal naïve: snaive(y, h=_)
- Drift: rwf(y, drift=TRUE, h=__)

where h is the forecasting horizon of interest.





Seasonal naïve

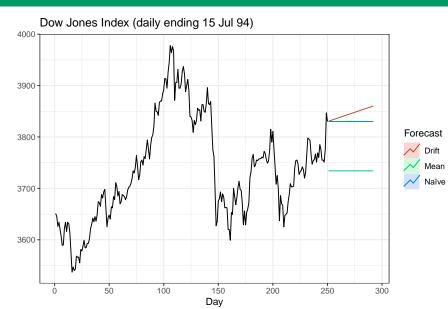
```
beer2 <- window(ausbeer, start=1992, end=c(2007,4))
# Plot some forecasts
autoplot(beer2) +
  autolayer(meanf(beer2, h=11), PI=FALSE, series="Mean") +
  autolayer(naive(beer2, h=11), PI=FALSE, series="Naïve") +</pre>
```

labs(title = "Forecasts for quarterly beer production",

autolayer(snaive(beer2, h=11), PI=FALSE,

x = "Year", y= "Megalitres") +
guides(colour=guide legend(title="Forecast"))

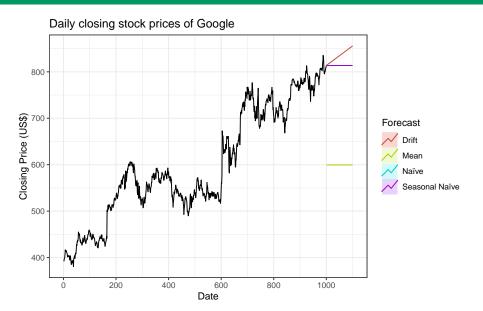
series="Seasonal naïve") +

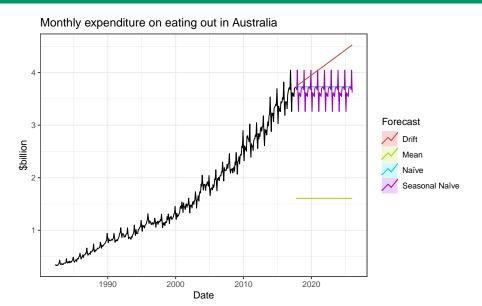


- Mean: meanf(y, h=__)
- Naïve: naive(y, h=__)
- Seasonal naïve: snaive(y, h=__)
- Drift: rwf(y, drift=TRUE, h=__)

Your turn

- Use these four functions to produce forecasts for goog and auscafe. Set h=100.
- Plot the results using autoplot().





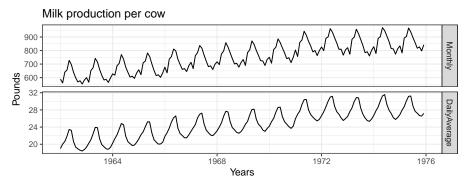
Section 2

Transformations & Adjustments

Calendar Adjustments

Some of the observed variation in seasonal data may be due to simple issues such as the number of calendar days. Usually much easier to remove the variation before fitting a forecasting model. The ${\tt monthdays}()$ function will compute the number of days in each month or quarter.

```
dframe <- cbind(Monthly = milk, DailyAverage = milk/monthdays(milk))
autoplot(dframe, facet = TRUE) + labs(x = "Years",
    y = "Pounds") + ggtitle("Milk production per cow")</pre>
```



Population adjustments

Data affected by population change (or size) are better expressed on a per-capita basis.

For example, if you are studying the relationship between income and literacy across two regions, rather than compare the total GDP, it might be more constructive to look at income per person. With such a large population, China's total GDP will definitely be larger than a Luxembourg for example. After removing the effect of population size however, Luxembourg would be the leader (IMF 2020).

Inflation adjustments

- Data which are affected by the value of money are best adjusted before modelling. For example, the average cost of a new house will have increased over the last few decades due to inflation.
- A \$1 this year is not the same as a \$1 twenty years ago. For this reason, financial time series are usually adjusted so that all values are stated in dollar values from a particular year (real prices).
- Price indexes are often constructed by government agencies. For consumer goods, a common price index is the Consumer Price Index (or CPI).
- To convert current prices (y_t) to real prices (in 2000 dollars x_{2000}) we can use the transformation:

$$x_{2000} = \frac{y_{2020}}{CPI_{\text{based in 2000}}}$$

Box-Cox transformations

Variance stabilization

If the data show different variation at different levels of the series, then a transformation can be useful.

Denote original observations as y_1, \ldots, y_n and transformed observations as w_1, \ldots, w_n .

Box-Cox transformations

Variance stabilization

If the data show different variation at different levels of the series, then a transformation can be useful.

Denote original observations as y_1, \ldots, y_n and transformed observations as w_1, \ldots, w_n .

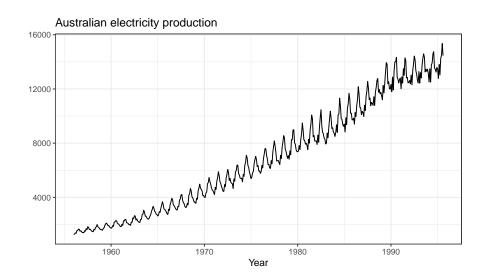
Mathematical transformations for stabilizing variation

Square root
$$w_t = \sqrt{y_t}$$

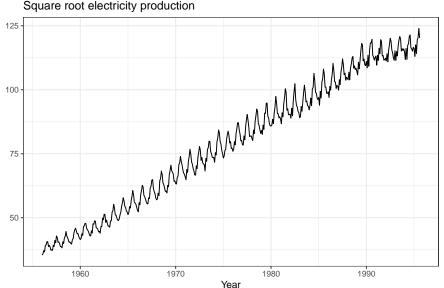
Cube root $w_t = \sqrt[3]{y_t}$ Increasing

Logarithm $w_t = \log(y_t)$ strength

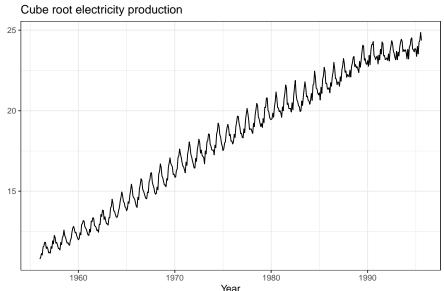
Logarithms, in particular, are useful because they are more interpretable: changes in a log value are relative (percent) changes on the original scale.

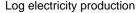


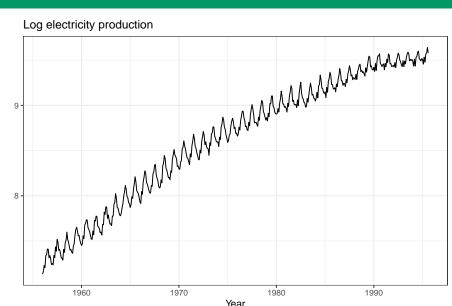


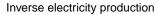


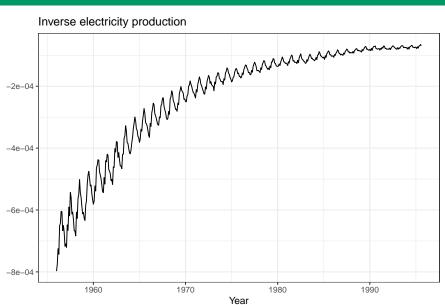












Box-Cox transformations

Each of these transformations is close to a member of the family of **Box-Cox transformations**:

$$w_t = \begin{cases} \log(y_t), & \lambda = 0; \\ (y_t^{\lambda} - 1)/\lambda, & \lambda \neq 0. \end{cases}$$

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- $\lambda = 1$: (No substantive transformation)
- $\lambda = \frac{1}{2}$: (Square root plus linear transformation)
- $\lambda = 0$: (Natural logarithm)
- $\lambda = -1$: (Inverse plus 1)

Section 3

Residual diagnostics

Fitted values

- $\hat{y}_{t|t-1}$ is the forecast of y_t based on observations y_1, \ldots, y_t .
- We call these "fitted values".
- Sometimes drop the subscript: $\hat{y}_t \equiv \hat{y}_{t|t-1}$.
- Often not true forecasts since parameters are estimated on all data.

For example:

- $\hat{y}_t = \bar{y}$ for average method.
- $\hat{y}_t = y_{t-1} + (y_T y_1)/(T-1)$ for drift method.

Forecasting residuals

Residuals in forecasting: difference between observed value and its fitted value: $e_t = y_t - \hat{y}_{t|t-1}$.

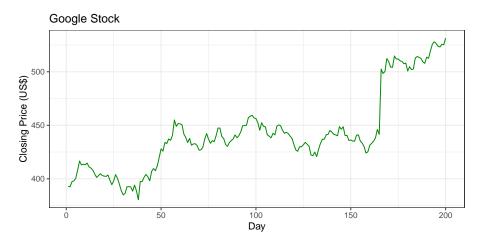
Assumptions

- \bullet $\{e_t\}$ uncorrelated.
 - If they aren't, then information left in residuals that should be used in computing forecasts.
- - If they don't, then forecasts are biased.

Useful properties (for prediction intervals)

- \bullet $\{e_t\}$ have constant variance.
- \bullet $\{e_t\}$ are normally distributed.

Example: Google stock price



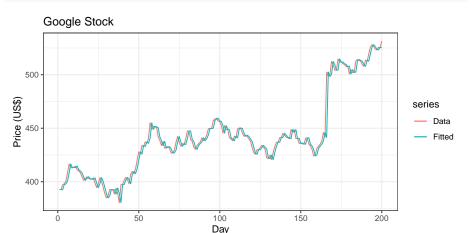
Example: Google stock price

Naïve forecast:

$$\hat{y}_{t|t-1} = y_{t-1} \\ e_t = y_t - y_{t-1}$$

Note: e_t are one-step-forecast residuals

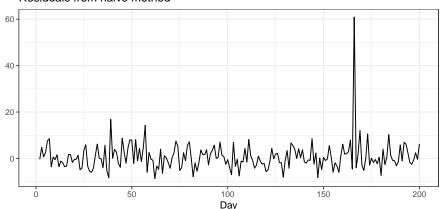
Example: Google stock price



Example: Google stock price

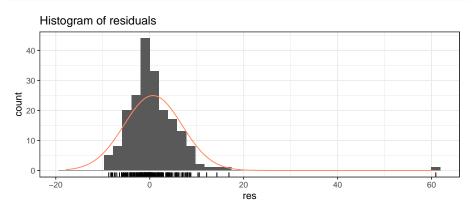
```
res <- residuals(naive(goog200))
autoplot(res) + labs(x = "Day", y = "") + ggtitle("Residuals from name of the state of the state
```

Residuals from naive method



Example: Google stock price

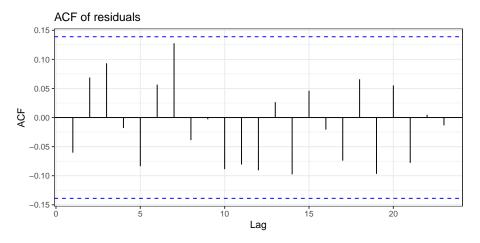
```
gghistogram(res, add.normal = TRUE) +
    ggtitle("Histogram of residuals")
```



Forecast with this model will be good but the non-normality of the residuals will prove problematic when calculating the prediction intervals.

Example: Google stock price

```
ggAcf(res) + ggtitle("ACF of residuals")
```



ACF of residuals

- We assume that the residuals are white noise (uncorrelated, mean zero, constant variance). If they aren't, then there is information left in the residuals that should be used in computing forecasts.
- So a standard residual diagnostic is to check the ACF of the residuals of a forecasting method.
- We *expect* these to look like white noise.

Portmanteau tests for autocorrelation

Consider a **whole set** of r_k values, and develop a test to see whether the **set** is (jointly) significantly different from a zero set.

1. Box-Pierce test

$$Q = T \sum_{k=1}^{h} r_k^2$$

where h is max lag being considered and T is number of observations.

- If each r_k close to zero, Q will be small.
- If some r_k values large (positive or negative), Q will be large.

Portmanteau tests for autocorrelation

Consider a **whole set** of r_k values, and develop a test to see whether the **set** is (jointly) significantly different from a zero set.

2. Ljung-Box test

$$Q^* = T(T+2) \sum_{k=1}^{h} (T-k)^{-1} r_k^2$$

where h is max lag being considered and T is number of observations.

- My preferences: h = 10 for non-seasonal data, h = 2m for seasonal data.
- Better performance, especially in small samples.

Portmanteau tests for autocorrelation

• If data are WN, Q^* has χ^2 distribution with (h - K) degrees of freedom where K = no. parameters in model. That is

$$Q^* \sim \chi^2_{(h-K)}$$

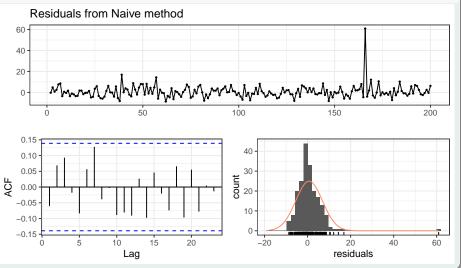
• When applied to raw data, set K = 0.

For the Google example:

```
# lag=h and fitdf=K
Box.test(res, lag=10, fitdf=0, type="Lj")
##
## Box-Ljung test
##
## data: res
## X-squared = 11.031, df = 10, p-value = 0.3551
```

Muting the Q^* statitics and plotting graphs only.

checkresiduals(naive(goog200),test = FALSE)



checkresiduals function

Muting the plot and getting only the Q^* statitics

```
checkresiduals(naive(goog200),plot = FALSE)

##

## Ljung-Box test

##

## data: Residuals from Naive method

## Q* = 11.031, df = 10, p-value = 0.3551

##

## Model df: 0. Total lags used: 10
```

Your turn

Compute seasonal naïve forecasts for quarterly Australian beer production from 1992.

```
beer <- window(ausbeer, start=1992)
fbeer <- snaive(beer)
autoplot(fbeer)</pre>
```

Test if the residuals are white noise.

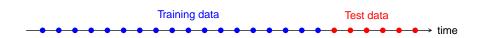
```
checkresiduals(fbeer)
```

What do you conclude?

Section 4

Evaluating forecast accuracy

Training and test sets



- A model which fits the training data well will not necessarily forecast well.
- A perfect fit can always be obtained by using a model with enough parameters.
- Over-fitting a model to data is just as bad as failing to identify a systematic pattern in the data.
 - The test set must not be used for *any* aspect of model development or calculation of forecasts.
 - Forecast accuracy is based only on the test set.

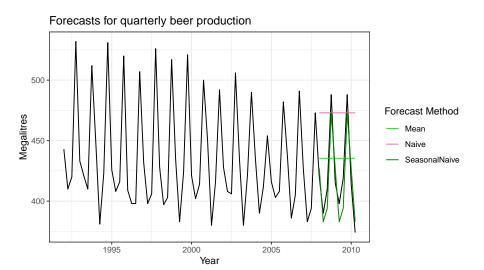
Forecast errors

Forecast "error" the difference between an observed value and its forecast.

$$e_{T+h} = y_{T+h} - \hat{y}_{T+h|T},$$

where the training data is given by $\{y_1, \ldots, y_T\}$

- Unlike residuals, forecast errors on the test set involve multi-step forecasts.
- These are *true* forecast errors as the test data is not used in computing $\hat{y}_{T+h|T}$.



```
beer2 <- window(ausbeer, start=1992, end=c(2007,4))</pre>
beerfit1 <- meanf(beer2,h=10)
beerfit2 <- rwf(beer2,h=10)
beerfit3 <- snaive(beer2,h=10)
tmp <- cbind(Data=window(ausbeer, start=1992),</pre>
             Mean=beerfit1[["mean"]],
             Naive=beerfit2[["mean"]],
             SeasonalNaive=beerfit3[["mean"]])
autoplot(tmp) +
  labs(title = "Forecasts for quarterly beer production",
    x = "Year", y = "Megalitres") +
scale color manual(values=c('black', 'green', 'hotpink',
  'green4'),
  breaks=c("Mean", "Naive", "SeasonalNaive"),
    name="Forecast Method")
```

$$y_{T+h} = (T+h)$$
th observation, $h = 1, ..., H$
 $\hat{y}_{T+h|T} =$ its forecast based on data up to time T .
 $e_{T+h} = y_{T+h} - \hat{y}_{T+h|T}$

$$\begin{aligned} \text{MAE} &= \text{mean}(|e_{T+h}|)\\ \text{MSE} &= \text{mean}(e_{T+h}^2)\\ \text{RMSE} &= \sqrt{\text{mean}(e_{T+h}^2)}\\ \text{MAPE} &= 100\text{mean}(|e_{T+h}|/|y_{T+h}|) \end{aligned}$$

- MAE, MSE, RMSE are all scale dependent.
- MAPE is scale independent but is only sensible if $y_t \gg 0$ for all t, and y has a natural zero.

Mean Absolute Scaled Error

$$MASE = mean(|e_{T+h}|/Q)$$

where Q is a stable measure of the scale of the time series $\{y_t\}$.

Proposed by Hyndman and Koehler (IJF, 2006).

For non-seasonal time series,

$$Q = (T-1)^{-1} \sum_{t=2}^{T} |y_t - y_{t-1}|$$

works well. Then MASE is equivalent to MAE relative to a naïve method.

Mean Absolute Scaled Error

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where Q is a stable measure of the scale of the time series $\{y_t\}$.

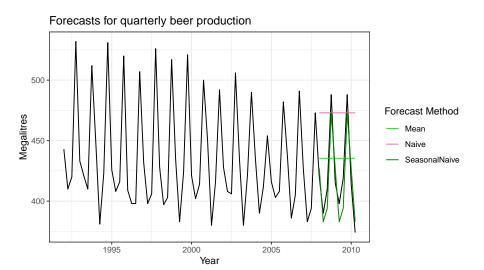
Proposed by Hyndman and Koehler (IJF, 2006).

For **seasonal** time series,

$$Q = (T - m)^{-1} \sum_{t=m+1}^{T} |y_t - y_{t-m}|$$

works well.

Then MASE is equivalent to MAE relative to a seasonal naïve method.



```
beer2 <- window(ausbeer, start=1992, end=c(2007,4))
beer3 <- window(ausbeer, start=2008)</pre>
beerfit1 <- meanf(beer2, h=10)
beerfit2 <- rwf(beer2, h=10)
beerfit3 <- snaive(beer2, h=10)
accuracy(beerfit1, beer3)
accuracy(beerfit2, beer3)
accuracy(beerfit3, beer3)
# Create a table to store the accuracy results.
tab <- matrix(NA,ncol=4,nrow=3)
tab[1,] \leftarrow accuracy(beerfit1, beer3)[2,c(2,3,5,6)]
tab[2,] \leftarrow accuracy(beerfit2, beer3)[2,c(2,3,5,6)]
tab[3,] \leftarrow accuracy(beerfit3, beer3)[2,c(2,3,5,6)]
colnames(tab) <- c("RMSE", "MAE", "MAPE", "MASE")</pre>
rownames(tab) <- c("Mean method", "Naïve method",</pre>
                   "Seasonal naïve method")
knitr::kable(tab, digits=2)
```

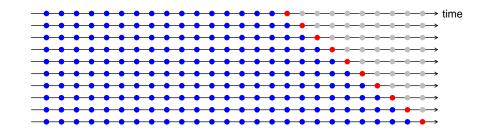
	RMSE	MAE	MAPE	MASE
Mean method	38.45	34.83	8.28	2.44
Naïve method	62.69	57.40	14.18	4.01
Seasonal naïve method	14.31	13.40	3.17	0.94

Time series cross-validation

Traditional evaluation



Time series cross-validation



tsCV function:

[1] 6.168928

```
e <- tsCV(goog200, rwf, drift=TRUE, h=1)
sqrt(mean(e^2, na.rm=TRUE))

## [1] 6.233245
sqrt(mean(residuals(rwf(goog200, drift=TRUE))^2,na.rm=TRUE))</pre>
```

A good way to choose the best forecasting model is to find the model with the smallest RMSE computed using time series cross-validation.

Section 5

Digression

Pipe function

Ugly code:

```
sqrt(mean(e^2, na.rm=TRUE))
## [1] 6.233245
sqrt(mean(residuals(rwf(goog200, drift=TRUE))^2,na.rm=TRUE))
```

Better with a pipe:

[1] 6.168928

```
goog200 %>% tsCV(forecastfunction=rwf, drift=TRUE, h=1) -> e
e^2 %>% mean(na.rm=TRUE) %>% sqrt
## [1] 6.233245
```

```
goog200 %>% rwf(drift=TRUE) %>% residuals -> res
res^2 %>% mean(na.rm=TRUE) %>% sqrt
```

e <- tsCV(goog200, rwf, drift=TRUE, h=1)

[1] 6.168928

Section 6

Prediction intervals

- A forecast $\hat{y}_{T+h|T}$ is (usually) the mean of the conditional distribution $y_{T+h} \mid y_1, \dots, y_T$.
- A prediction interval gives a region within which we expect y_{T+h} to lie with a specified probability.
- \bullet Assuming forecast errors are normally distributed, then a 95% PI is

$$\hat{y}_{T+h|T} \pm 1.96\hat{\sigma}_h$$

where $\hat{\sigma}_h$ is the st dev of the h-step distribution.

• When h = 1, $\hat{\sigma}_h$ can be estimated from the residuals.

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Naive forecast with prediction interval:

```
res_sd <- sqrt(mean(res^2, na.rm=TRUE))</pre>
c(tail(goog200,1)) + 1.96 * res sd * c(-1,1)
## [1] 519.3872 543.5694
naive(goog200, level=95)
##
      Point Forecast Lo 95 Hi 95
## 201
             531,4783 519,3105 543,6460
## 202
             531,4783 514,2705 548,6861
## 203
             531,4783 510,4031 552,5534
## 204
             531,4783,507,1428,555,8138
## 205
             531,4783 504,2704 558,6862
## 206
             531,4783 501,6735 561,2830
## 207
             531,4783,499,2854,563,6711
## 208
             531,4783,497,0627,565,8939
## 209
             531,4783,494,9750,567,9815
```

531.4783 493.0005 569.9561

- Point forecasts are often useless without prediction intervals.
- Prediction intervals require a stochastic model (with random errors, etc).
- Multi-step forecasts for time series require a more sophisticated approach (with PI getting wider as the forecast horizon increases).

Assume residuals are normal, uncorrelated, sd = $\hat{\sigma}$:

Mean forecasts:
$$\hat{\sigma}_h = \hat{\sigma}\sqrt{1+1/T}$$

Naïve forecasts:
$$\hat{\sigma}_h = \hat{\sigma}\sqrt{h}$$

Seasonal naïve forecasts
$$\hat{\sigma}_h = \hat{\sigma}\sqrt{k+1}$$

Drift forecasts:
$$\hat{\sigma}_h = \hat{\sigma} \sqrt{h(1+h/T)}$$
.

where k is the integer part of (h-1)/m.

Note that when h=1 and T is large, these all give the same approximate value $\hat{\sigma}$.

- Computed automatically using: naive(), snaive(), rwf(), meanf(), etc.
- Use level argument to control coverage.
- Check residual assumptions before believing them.
- Usually too narrow due to unaccounted uncertainty.