



Time Series in R: Forecasting and Visualisation

Forecast evaluation

29 May 2017

Outline

- 1 Forecasting residuals
 - 2 Evaluating forecast accuracy
 - 3 Forecasting benchmark methods
 - 4 Lab session 7
 - 5 Time series cross-validation
 - 6 Lab session 8

Fitted values

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

Forecasting residuals

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

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Assumptions

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

Forecasting residuals

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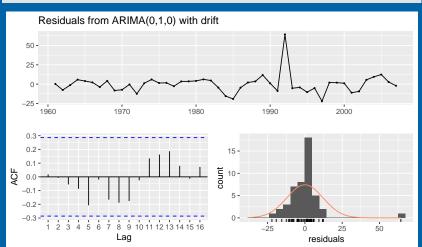
Assumptions

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Useful properties (for prediction intervals)

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

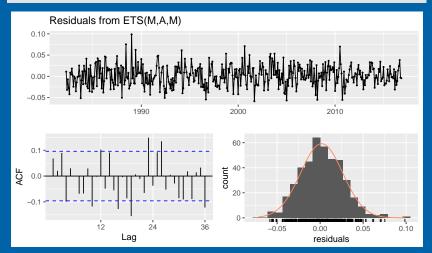
livestock %>% auto.arima %>% checkresiduals(test=FALSE)



livestock %>% auto.arima %>% checkresiduals(plot=FALSE)

```
##
## Ljung-Box test
##
## data: Residuals from ARIMA(0,1,0) with drift
## Q* = 8.6, df = 9, p-value = 0.5
##
## Model df: 1. Total lags used: 10
```





```
auscafe %>% ets %>% checkresiduals(plot=FALSE)
```

```
##
## Ljung-Box test
##
## data: Residuals from ETS(M,A,M)
## Q* = 64, df = 8, p-value = 7e-11
##
## Model df: 16. Total lags used: 24
```

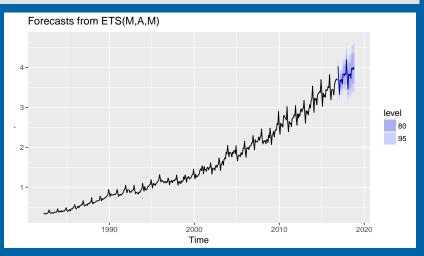
Residuals and forecasting

- Autocorrelations left in residuals suggest the forecast method can be improved (in theory).
- Small autocorrelations have little effect, even if significant.
- Non-Gaussian residuals can be handled using bootstrapped forecast intervals:

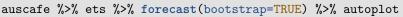
```
forecast(..., bootstrap=TRUE)
```

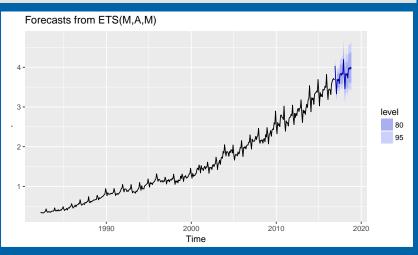
Bootstrapped forecast intervals

auscafe %>% ets %>% forecast %>% autoplot



Bootstrapped forecast intervals





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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 errors

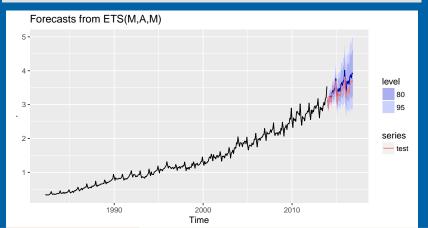
Forecast "error": the difference between an observed value and its forecast.

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

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 $\hat{y}_{T+h|T}$.

```
training <- window(auscafe, end=c(2013,12))
test <- window(auscafe, start=c(2014,1))
training %>% ets %>% forecast(h=length(test)) -> fc
autoplot(fc) + autolayer(test)
```



Let $\hat{y}_{t+h|t}$ denote the forecast of y_{t+h} using data up to time t.

Training set measures:

$$\begin{aligned} \text{MAE} &= \tfrac{1}{T} \sum_{t=1}^{T} |y_t - \hat{y}_{t|t-1}| \\ \text{MSE} &= \tfrac{1}{T} \sum_{t=1}^{T} (y_t - \hat{y}_{t|t-1})^2 \\ \text{MAPE} &= \tfrac{100}{T} \sum_{t=1}^{T} |y_t - \hat{y}_{t|t-1}| / |y_t| \end{aligned}$$

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- 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.

Let $\hat{y}_{t+h|t}$ denote the forecast of y_{t+h} using data up to time t.

Test set measures:

$$\begin{aligned} \text{MAE} &= \frac{1}{H} \sum_{h=1}^{H} |y_{T+h} - \hat{y}_{T+h|T}| \\ \text{MSE} &= \frac{1}{H} \sum_{h=1}^{H} (y_{T+h} - \hat{y}_{T+h|T})^2 \qquad \text{RMSE} = \sqrt{\frac{1}{H} \sum_{h=1}^{H} (y_{T+h} - \hat{y}_{T+h|T})^2} \\ \text{MAPE} &= \frac{100}{H} \sum_{h=1}^{H} |y_{T+h} - \hat{y}_{T+h|T}| / |y_t| \end{aligned}$$

- MAE, MSE, RMSE are all scale dependent.
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Mean Absolute Scaled Error

MASE =
$$\frac{1}{H} \sum_{h=1}^{H} |y_{T+h} - \hat{y}_{T+h|T}|/Q$$

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,

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

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

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Proposed by Hyndman and Koehler (IJF, 2006).

For seasonal time series,

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

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

```
training <- window(auscafe, end=c(2013,12))
test <- window(auscafe, start=c(2014,1))
training %>% ets %>% forecast(h=length(test)) -> fc
accuracy(fc, test)
```

```
## Training set 0.001482 0.03816 0.02761 0.09342
## Test set -0.121597 0.15318 0.13016 -3.51895
## MAPE MASE ACF1 Theil's U
## Training set 2.056 0.2881 0.2006 NA
## Test set 3.780 1.3583 0.6323 0.7647
```

Poll: true or false?

- Good forecast methods should have normally distributed residuals.
- A model with small residuals will give good forecasts.
- The best measure of forecast accuracy is MAPE.
- If your model doesn't forecast well, you should make it more complicated.
- Always choose the model with the best forecast accuracy as measured on the test set.

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Average method

Forecasts equal to mean of historical data.

Naïve method

- Forecasts equal to last observed value.
- Consequence of efficient market hypothesis.

Seasonal naïve method

Forecasts equal to last value from same season.

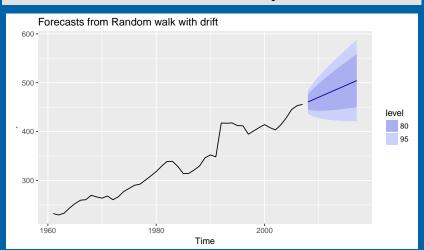
Drift method

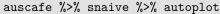
- Forecasts equal to last value plus average change.
- Equivalent to extrapolating a line drawn between first and last observations.

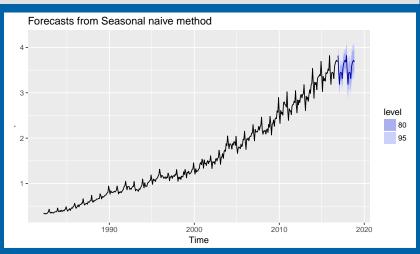
- Mean: meanf(y, h=20)
- Naïve: naive(y, h=20)
- Seasonal naïve: snaive(y, h=20)
- Drift: rwf(y, drift=TRUE, h=20)

Check that your method does better than these standard benchmark methods.

livestock %>% rwf(drift=TRUE) %>% autoplot







```
training %>% ets %>% forecast(h=length(test)) -> fc_ets
training %>% snaive(h=length(test)) -> fc_snaive
accuracy(fc_ets, test)
```

```
## ME RMSE MAE MPE
## Training set 0.001482 0.03816 0.02761 0.09342
## Test set -0.121597 0.15318 0.13016 -3.51895
## MAPE MASE ACF1 Theil's U
## Training set 2.056 0.2881 0.2006 NA
## Test set 3.780 1.3583 0.6323 0.7647
```

accuracy(fc_snaive,test)

```
## ME RMSE MAE MPE MAPE
## Training set 0.08569 0.1226 0.09583 6.529 7.286
## Test set 0.41363 0.4344 0.41363 12.183 12.183
## MASE ACF1 Theil's U
## Training set 1.000 0.8425 NA
## Test set 4.317 0.6438 2.165
```

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Lab Session 7

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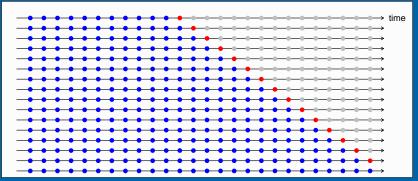
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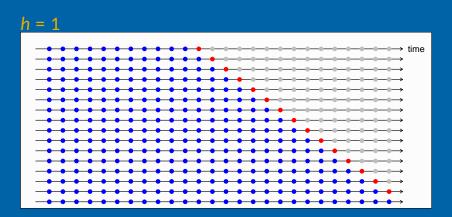
Traditional evaluation



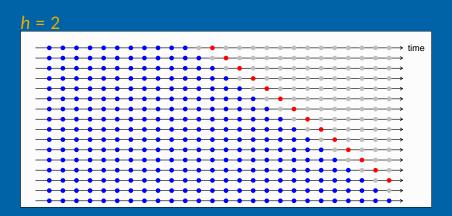
Traditional evaluation



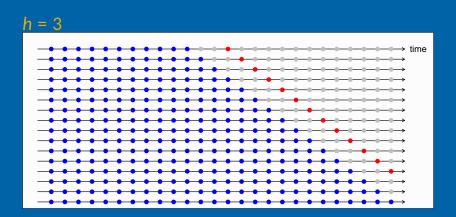




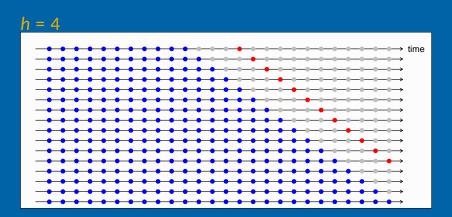
- Forecast accuracy averaged over test sets.
- Also known as "evaluation on a rolling forecasting origin"



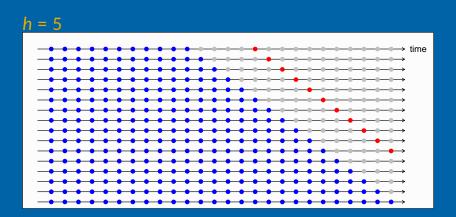
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tsCV function:

```
e <- tsCV(ts, forecastfunction, h=1, ...)
e1 <- tsCV(auscafe, stlf,
  etsmodel="AAN", damped=FALSE, lambda=0)
autoplot(e1)
   0.1 -
o 0.0 -
  -0.1 -
                                    2000
                                                    2010
                   1990
                                   Time
```

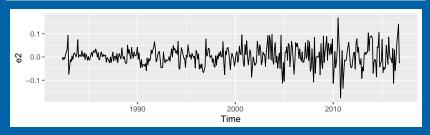
[1] 0.0259

sqrt(mean((e1/auscafe)^2, na.rm=TRUE))

tsCV function:

For ets and auto.arima, you need to write a single forecasting function:

```
fets <- function(x, h, model="ZZZ", damped=NULL, ...) {
  forecast(ets(x, model=model, damped=damped), h=h)
}
e2 <- tsCV(auscafe, fets, model="MAM", damped=FALSE)
autoplot(e2)</pre>
```



```
sqrt(mean((e2/auscafe)^2, na.rm=TRUE))
```

tsCV function:

Comparison should be over the same observations:

```
pe1 <- window(100*e1/auscafe, start=1985)
pe2 <- window(100*e2/auscafe, start=1985)
sqrt(mean(pe1^2, na.rm=TRUE))</pre>
```

```
## [1] 2.571
```

```
sqrt(mean(pe2^2, na.rm=TRUE))
```

```
## [1] 2.733
```

- A good way to choose the best forecasting model is to find the model with the smallest RMSE computed using time series cross-validation.
- Minimizing AICc is asymptotically equivalent to minimizing tscv with h = 1.

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Lab Session 8