ARMA(1,1)-GARCH(1,1)  
Estimation and forecast using rugarch 1.2-2  

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1 Introduction

First we specify a model ARMA(1,1)-GARCH(1,1) that we want to estimate. Secondly we touch upon the matter of fixing certain parameters of the model. Thirdly we get some data to estimate the model on and estimate the model on the data. After having estimated the model we inspect the created R-object from the fitting of the model.

Then we use the model for making a forecast: 1) A simple forecast and 2) a rolling forecast and 3) a rolling forecast with reestimation of model.

Throughout I make some comments on calculation of VaR.

Reference manual, source code and a very helpful vignette is available at:
http://cran.r-project.org/web/packages/rugarch/index.html

I refer to the reference manual and vignette when relevant so if not for any other reason - although other reasons are plenty - it is a good idea to download and read these.

The author of the rugarch package Alexios Ghalanos has a blog:
http://www.unstarched.net/blog/

I'm using rugarch: Univariate GARCH models R-package version 1.2-2 by Alexios Ghalanos.

2 Model specification - »uGARCHspec«

To fit a GARCH-model the first step is to create an instance of the S4-class uGARCHspec. The created object serves the purpose of specifying the model
to be estimated. The chosen model can be estimated with all parameters free or alternatively with some parameters fixed.

2.1 Modelspecification (free parameters only)

The uGARCHspec class is documented on page 86 of the rugarch reference manual and the usual help page is found typing ?ugarchspec in R-console. On the help page we find the following peace of code:

```r
model=ugarchspec(
  variance.model = list(model = "sGARCH", garchOrder = c(1, 1)),
  mean.model = list(armaOrder = c(1, 1), include.mean = TRUE),
  distribution.model = "norm"
)
```

For simplicity I have left out some options we are not going to use which are explicited on the helppage. Also I have added the `model=`... so that our instance of the uGARCHspec class i called `model`.

So when creating the `model` we specify the `variance.model`, the `mean.model` and the `distribution.model`. Notice that the first two arguments `mean.model` and `variance.model` takes arguments that are objects of the type `list`.

We have set the argument `model` in the `variance.model` to have the value "sGARCH" with `garchOrder=c(1,1)` so that our choice of variancemodel is:

\[ \sigma_t^2 = \omega + \alpha \epsilon_{t-1}^2 + \beta \sigma_{t-1}^2 \]  

as explained in the vignette on page 6 with \( p=1, q=1 \) and \( m=0 \) since we are ignoring external regressors.

The meanmodel is chosen to have `armaOrder=c(1,1)` and we include a mean - constant - by `mean=TRUE` so that our chosen model is:

\[ r_t = \mu + \theta_1 (r_{t-1} - \mu) + \theta_2 \epsilon_{t-1} + \epsilon_t \]  

When specifying the distribution we have chosen a normaldistribution using the argument "norm" so:

\[ \epsilon_t = \sigma_t z_t \quad z_t \sim N(0,1) \]  

often one would choose other distributions due to the stylised fact of financial series having fat tails. One option would be to use the Student-t distribution `distribution.model="std"` or a skewed version of it "sstd". The conditional distributions are discussed on page 13 of the vignette.
2.2 Model specification (fixed parameters)

We might want to fix some or all of the parameters in the model. We could fix some of the parameters and estimate the rest for example with the purpose of making likelihood ratio comparison. We could also fix all parameter values with the purpose of using the model for forecasting or simulation.

To fix the parameters we need to use the argument `fixed.pars = list()` in the `ugarchspec()` see helppage `?ugarchspec`. We also need to find the name of the parameter we want to fix. If you have chosen a model type for example the sGARCH in our case you can find the name of the relevant parameters by typing `spec@model.pars`. You need to substitute `spec` with the name you gave to your instance of the uGARCHspec class, which in our case was `model` so we simply type `model@model.pars`.

From the top: First we choose the sGARCH-model but lets try it with higher order ARMA and GARCH and a skewed version og the Student-t:

```r
model2=ugarchspec(
  variance.model = list(model = "sGARCH", garchOrder = c(2, 2)),
  mean.model = list(armaOrder = c(2, 2), include.mean = TRUE),
  distribution.model = "sstd"
)
```

Then to inspect the parameter names:

```r
model2@model$pars
```

And you should get a printout similar to the one in tabel 1. From the table the names of the parameters are available in the first column and the fourth column takes the value 1 where the relevant parameter is included 0 otherwise. The fifth column - called »Estimate« - indicates whether the parameter is estimated or fixed. So $\beta$ in our GARCH-model when the `garchOrder=c(1,1)` is probably called `beta1` if there is any justice in the world. So to fix the parameter we do:

```r
model=ugarchspec(
  variance.model = list(model = "sGARCH", garchOrder = c(1, 1)),
  mean.model = list(armaOrder = c(1, 1), include.mean = TRUE),
  distribution.model = "norm"
  fixed.pars = list(beta1=0.86) )
```

You can try `model@model$pars` to see if the relevant parameter has been set to the chosen value.
### Table 1: R-print from `model2@model$pars`

<table>
<thead>
<tr>
<th></th>
<th>Level</th>
<th>Fixed</th>
<th>Include</th>
<th>Estimate</th>
<th>LB</th>
<th>UB</th>
</tr>
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<td>0</td>
<td>0</td>
<td>0</td>
<td>NA</td>
<td>NA</td>
</tr>
</tbody>
</table>

#### 3 Getting data

Before we can estimate the model we have to get some data. You can use the SP500 index which is included in the rugarch package simply type:

```r
data(sp500ret)
```

and then the sp500-index is loaded and is called `sp500ret`. Alternatively you can load data from Yahoo using the `get.hist.quote()` function from the package `tseries`:

```r
library(tseries)
sp500.prices = get.hist.quote(
  instrument = "^GSPC",
  quote = "Adj",
  provider = c("yahoo"), method = NULL,
  origin = "1899-12-30", compression = "d",
  retclass = c("zoo"), quiet = FALSE, drop = FALSE
)
sp500 = as.data.frame(sp500.prices)
N = length(sp500[,1])
sp500.returns = 100*(log(sp500[2:N,])-log(sp500[1:(N-1),]))
```

This downloads the adjusted closing prices - adjusted for dividends - and calculates continuously compounded returns.
4 Estimating the model

To estimate the model we use the model-object created and the data sp500ret and provide them as arguments to the uGARCHfit() function:

```
modelfit=ugarchfit(spec=model,data=sp500ret)
```

This creates an object we have chosen to call modelfit which is an instance of the S4 class uGARCHfit specific to the rugarch package as documented on page 65 of the rugarch reference manual.

This object is far more interesting in than the model in the sense that it contains information we need to be able to manipulate and as a minimum print. First of all simply typing the objectname prints the parameter estimates of the model along with standard errors and a lot of interesting statistics (as is custom in R):

```
modelfit
```

I'm not going to delve on these statistics they are explained better elsewhere. References for the relevant litterature can be found in the rugarch vignette. However I will make a comment on S4-objects. Just to repeat: Estimating the model creates the object we have chosen to call modelfit as an instance of the class uGARCHfit to get the class type:

```
class(modelfit)
#[1] "uGARCHfit"
#attr(,"package")
#[1] "rugarch"
```

Another useful feature is that you can get the structure of the object by typing `str(modelfit)`. This results in a somewhat »messy« print but if you look for the @ you should be able to spot the following structure in the printout:

```
Formal class 'uGARCHfit' [package "rugarch"] with 2 slots
..@ fit
..@ model
```

Since we have an S4 object the object has what is called slots and in this case there are 2 slots the first is called fit the second model. This is also apparent typing `slotNames(modelfit)`. Lets think of the slots as places to store data and in each of the slots there are »containers« where every container is indicated by $(the containers are of course R-objects like vectors,
matrices or lists). Reading the print from `str(modelfit)` you will notice that the first $ is followed by `hessian`:

Formal class 'uGARCHfit' [package "rugarch"] with 2 slots
..@ fit :List of 25
   ..$ hessian

This is interesting because it tells you how to access the information stored in the object by typing:

```r
modelfit@fit$hessian
```

You could get the parameter estimates by typing:

```r
modelfit@fit$coef
```

The fitted values:

```r
modelfit@fit$fitted.values
```

The point is that how to access the data is indicated from the `str(modelfit)` printout. However you could also use the designed `methods` which are documented on page 65 of the reference manual:

```r
coeff(modelfit)
infocriteria(modelfit)
sigma(modelfit)
fitted(modelfit)
residuals(modelfit)
```

To name a few. Be sure to update R and rugarch package since the available methods possibly change from version to version (I had problems with the method quantile used when calculating VaR). To calculate VaR use the quantile method:

```r
VaR=quantile(modelfit,0.01)
```

or in our case with standardized normal distribution we could do:

```r
VaR=modelfit@fit$sigma*qnorm(0.01)+modelfit@fit$fitted.values
```

Since the model is estimated on the observations from the periods for which we calculate VaR this is called in-sample VaR. As such it may seem uninteresting since what would be the purpose of calculating a risk measure of an already realized contingency? Hence later we do forecasting.
5 Forecasting

Having estimated a model we might want to use it for forecasting. Reading the helpfile ?ugarchforecast it says:

\[
\text{ugarchforecast}(\text{fitORspec, data = NULL, n.ahead = 10, n.roll = 0, out.sample = 0})
\]

So the first argument is an estimated model - \textit{fit} - or a specified model - \textit{spec}. Using a specified model all parameter values need to be fixed and the \textit{data} argument cannot be \textit{NULL}. Using a fitted object the data is contained in the fitted object and hence do not need to be supplied, so the \textit{data=\text{NULL}} can be used to forecast from the same series as used in estimation.

In our case the fitted model \textit{modelfit} is used for forecasting. First lets do a simple forecast:

\[
\text{modelfor=ugarchforecast(modelfit, data = NULL, n.ahead = 10, n.roll = 0, out.sample = 0)}
\]

The last observation of the used dataset has the date 2009-01-30 and this is \textit{T_0} in the forecast:

\begin{verbatim}
0-roll forecast [T0=2009-01-30]:
 Series Sigma
T+1 0.0016724 0.02480
T+2 0.0015259 0.02474
T+3 0.0013981 0.02467
T+4 0.0012865 0.02461
T+5 0.0011891 0.02455
T+6 0.0011041 0.02448
T+7 0.0010299 0.02442
T+8 0.0009651 0.02436
T+9 0.0009086 0.02430
T+10 0.0008592 0.02424
\end{verbatim}

The object \textit{modelfor} is an instance of the class \textit{uGARCHforecast} page 71 of the reference manual. The object has two slots with names:

\[
\text{slotNames(modelfor)}
\]

# [1] "forecast" "model"

You can access the forecast of the series with method \texttt{fitted(modelfor)} and the sigma (the conditional standard deviation) as \texttt{sigma(modelfor)}. Also available are some plots \texttt{plot(modelfor)} and then select an appropri-
ate option.

We could forecast with a larger n.ahead=50

modelfor=ugarchforecast(modelfit, data = NULL, n.ahead = 50, n.roll = 0, out.sample = 0)

An using str(modelfor) we can see that the adress for the seriesforecast is modelfor@forecast$seriesFor and we can do a simple plot:

plot(modelfor@forecast$seriesFor)

exhibiting exponential decay to unconditional mean modelfit@fit$coef["mu"] or by method uncmean(modelfit).

5.1 Rolling forecast

Using a model for forecast it might be interesting to see how well the model forecasts so we need some observations to compare with the forecasted values. For this reason it is possible to fit the model with the argument out.sample different than 0. In this case it becomes possible to do rolling 1-step-head forecasts. To use this option we have to estimate the model not on all observations hence we use the out.sample argument. For simplicity I choose the out.sample=2 and I choose a GARCH-model with a simpler meanequation ARMA(0,0):

model=ugarchspec (  
  variance.model = list(model = "sGARCH", garchOrder = c(1, 1)),  
  mean.model = list(armaOrder = c(0, 0)),  
  distribution.model = "norm"  
)

modelfit=ugarchfit(model,data=sp500ret,out.sample=2)

So fx. if we have 500 observations and choose out.sample=100 we would be estimating the model on the first 400 observations. With the dataset sp500ret we have $N = 5523$ observations so the index of the last observation on which the model is estimated on is $T = 5521$ since we have chosen out.sample=2:

modelfit@model$modeldata$T  # 5521

Then we do the forecasts:

modelfor=ugarchforecast(modelfit, data = NULL, n.ahead = 1, n.roll = 2, out.sample = 2)
and we look at the forecasted values:

```
sigma(modelfor)
# 2009-01-28 2009-01-29 2009-01-30
# T+1 0.02413003 0.02512952 0.02492642
fitted(modelfor)
# 2009-01-28 2009-01-29 2009-01-30
# T+1 0.0005205228 0.0005205228 0.0005205228
```

The values are forecasted values based on the information of the listed date, the values are not forecasts for the expected value of the specific date. Observation \( T = 5521 \) is the \( T_0 \) observation which in \texttt{sp500ret} has the date 2009-01-28 and being part of the sample on which the model is estimated there can be no forecast for the date (following conventional terminology and only calling something a forecast if it is out-of-sample).

Anyway since we have a simple model it should be easy to control for this by manual calculation. The mean is constant in our model so the forecast of the returns series for any period is simply the estimated constant mean 

\[
\text{coef(modelfit)}["mu"]
\]

For the period \( T = 5521 \) we have an estimated residual

\[
\text{residuals(modelfit)}[5521]
\]

but for the two periods for which we are forecasting we lack these but can calculate them comparing forecast of the series - the constant mean - to the observed return. With these residuals we can forecast the sigmas.

We could redo the forecasting manually along above lines with the following script:

```
#Getting the expected return from the estimated model
mu=coef(modelfit)["mu"]
#Getting the relevant observed return from last two periods of sp500ret
return=sp500ret[5522:5523,]
#Getting residual of period T=5521 from modelfit
e5521=as.vector(residuals(modelfit)[5521])
#Calculating residuals of period T=5522 and 5523
e5522=return[1]-mu
e5523=return[2]-mu
#Taking estimated parameters from modelfit
theta=coef(modelfit)
#Making function for forecast
.fgarch=function(e,sigma_0,theta) {
  omega=theta["omega"]
  alpha=theta["alpha1"]
```
beta=theta["beta1"]
sigma_1 = sqrt(omega + alpha*e^2 + beta*sigma_0^2)
names(sigma_1)=NULL
return(sigma_1)
}

# Getting estimated sigma for period T=5521 from modelfit
sigma5521=as.vector(sigma(modelfit)[5521])

# Forecast sigma_5522 and comparing with rugarch forecast
sigma5522=.fgarch(e5521,sigma_0=sigma5521,theta)
sigma5522
sigma(modelfor)[1]

# Forecast sigma_5523 and comparing with rugarch forecast
sigma5523=.fgarch(e5522,sigma_0=sigma5522,theta)
sigma5523
sigma(modelfor)[2]

# Forecast sigma_5524 and comparing with rugarch forecast
sigma5524=.fgarch(e5523,sigma_0=sigma5523,theta)
sigma5524
sigma(modelfor)[3]

And we are able to get the same values manually as using the rugarch function \texttt{ugarchforecast}. Following the script hopefully makes it clear what is being reported calling \texttt{sigma(modelfor)} and \texttt{fitted(modelfor)} if that was somehow unclear to begin with.

If the mean equation was ARMA(1,1) it would be a little more complicated due to the expected return of the next period not simply being a constant. First we calculate the variance for $T_1$:

$$\sigma_{T_1}^2 = \hat{\omega} + \hat{\alpha} \epsilon_{T_0}^2 + \hat{\beta} \sigma_{T_0}^2$$

and we forecast the return series using the mean model:

$$r_{T_1} = \hat{\mu} + \hat{\theta}_1 (r_{T_0} - \hat{\mu}) + \hat{\theta}_2 \epsilon_{T_0}$$

Comparing the predicted return $r_{T_1}$ to the observed $r_{+1}$ we can calculate a new residual $\epsilon_{T_1} = r_{+1} - r_{T_1}$. And given this residual we can calculate a new 1-step-head forecast for period $T_2$ for both the sigma and the return series. With the forecast $r_{T_2}$ calculated we can calculate $\epsilon_{T_2} = r_{+2} - r_{T_2}$ and so on by iteration. Since we constantly allow ourselves to use last periods return in forecasting the return for the next period we say $\texttt{n.ahead}=1$, hence conditioning on the information $\Omega_{t-1}$. 

10
Since we have calculated sigmas out-of-sample we could use these to calculate VaR out of sample. Since the density of our model is normal the 1-step-head density is normal with the forecasted standard deviation given that the model is valid. To calculate VaR - using the simple GARCH model with the constant meanequation - we would do:

\[ \text{qnorm}(0.05) \times \text{sigma(modelfor)} + \text{coef(modelfit)}["mu"] \]

or using the built-in function of the rugarch package:

\[ \text{quantile(modelfor,0.05)} \]

If not using the model with the constant mean but instead using the ARMA(1,1) for meanequation (or another meanequation for that matter) do:

\[ \text{qnorm}(0.05) \times \text{sigma(modelfor)} + \text{fitted(modelfor)} \]

or using the built-in function of the rugarch package:

\[ \text{quantile(modelfor,0.05)} \]

Having used the forecast for a period where actual observations are available we gain two new plot options using: \text{plot(modelfor)} where we now can do the rolling options.

### 6 Rolling forecast with reestimation

If we want to know how well our model does VaR calculation we would probably want to calculate VaR out of sample since this mimics the actual practice closer. Calculating VaR on a daily basis it probably would not be a problem to reestimate the model every day hence reestimating for each 1-step-head VaR calculation. However if we want to leave 1000 observations out of sample to calculate 1000 out-of-sample VaR measures we would have to reestimate the model 1000 times and this would be computationally heavy and timeconsuming. On the other hand simply calculating 1000 VaR measures without reestimation would probably not mimic actual practice and would due to changing parameters give bad VaR estimates.

A compromise using the \text{sp500ret} would be to \text{out.sample=1000} hence first estimating the model on 5523−1000 = 4523 observations. Then calculate 50 1-step-ahead forecasts and out-of-sample VaR estimates and the reestimate the model on the observations 1, 2, ..., 4535, ..., 4583. Based on the new parameter estimates we could then do yet another 50 1-step-ahead forecasts continuing like this until we have forecasts and VaR calculated for all 1000
6.1 Doing a rolling forecast

The rugarch package allows this to be done quite easily using the `ugarchroll`. Let's try doing this with the ARMA(1,1)-GARCH(1,1)-model leaving only 100 observations out-of-sample and refitting every 50 observations so as not to be overwhelmed by the complexity of the returned object:

```r
model = ugarchspec (  
  variance.model = list(model = "sGARCH", garchOrder = c(1, 1)),  
  mean.model = list(armaOrder = c(1, 1)),  
  distribution.model = "norm"  
)  
modelroll = ugarchroll (  
  spec = model, data = sp500ret, n.ahead = 1, forecast.length = 100,  
  n.start = NULL, refit.every = 50, refit.window = c("recursive"),  
  window.size = NULL, solver = "hybrid", fit.control = list(),  
  solver.control = list(), calculate.VaR = TRUE, VaR.alpha = c(0.01, 0.05),  
  cluster = NULL, keep.coef = TRUE  
)
```

So we choose the model using `ugarchspec` and give this as an argument value to `ugarchroll` by `spec=model`. Since the model has not been estimated we have to decide what data to use `data=sp500ret`. We use the `n.ahead=1` and as said on the `?ugarchroll` help page » Only rolling 1-ahead forecasts are supported spanning the dataset, which should be useful for backtesting models. Anything more complicated should be wrapped by the user by making use of the underlying functions in the package.«

Let's have a look at the returned object:

```r
slotNames(modelroll)  
# [1] "model" "forecast"
```

Object returned has two slots and we take a closer look on the second:

```r
str(modelroll@forecast)  
#List of 2  
#$ VaR :'data.frame': 100 obs. of 3 variables:  
# ..$ alpha(1%): num [1:100] -0.032 -0.0378 -0.0364 -0.036 -0.0345  
# ..$ alpha(5%): num [1:100] -0.0224 -0.0262 -0.0253 -0.0252 -0.0241  
# ..$ realized : num [1:100] -0.03473 0.00613 0.01371 0.00212 -0.04828
```
# $ density:'data.frame': 100 obs. of 6 variables:
#   ...$ Mu : num [1:100] 0.000792 0.001651 0.001418 0.001015 0.000935
#   ...$ Sigma: num [1:100] 0.0141 0.0169 0.0162 0.0159 0.0152
#   ...$ Skew : num [1:100] 0 0 0 0 0
#   ...$ Shape: num [1:100] 0 0 0 0 0
#   ...$ Shape(GIG): num [1:100] 0 0 0 0 0
#   ...$ Realized: num [1:100] -0.03473 0.00613 0.01371 0.00212 -0.04828

We see that we get some objects with 100 recorded values since we left 100 observations out of sample. We get the Value at Risk measure for the long position $\alpha = 0.01$ and $\alpha = 0.05$ because we chose calculate.VaR = TRUE, VaR.alpha = c(0.01, 0.05) in the uGarchroll() call. Also included in the VaR-data-frame is the realized values which is the actual returns from the period associated with the calculated VaR which is a nice feature when we have to calculate whether or not a VaR-violation has occurred - a VaR-violation being defined as an instance where the realized return is smaller than the VaR-measure (long position). Lets just calculate the rate of VaR-violations for:

```r
VaR = modelroll@forecast$VaR[, "alpha(1%)"]
return = modelroll@forecast$VaR[, "realized"]
Hit = return < VaR
sum(Hit)
# 5
```

So there are 5 VaR-violations out of 100 observations which is exactly the 5% expected given a $\alpha = 0.05$ (so the model would pass the Kupiec test at least for $\alpha = 0.05$).

We also get a density data-frame where the values of the forecasted expected return is stored under `modelroll@forecast$density[, "Mu"]` and the forecasted volatility is stored as `modelroll@forecast$density[, "Sigma"]`. Have a look at the reference manual on page 80 for the »uGARCHroll« in order to see methods for this class.

The Skew, Shape and Shape(GIG) are all 0 because they relate to parameters of the chosen distribution of the model and since we are using standardizad normal distribution these are all 0.

So one question could be: How do I calculate the VaR of another $\alpha$-level? One option would be to reestimate - hence doing the uGarchroll again but as should be apparent from earlier calculations this is unnecessary. Simply do:

```r
q_st = qnorm(0.025)
sigma = modelroll@forecast$density[, "Sigma"]
```
If you are using another distribution you should use the `qdist()` function as explained on page 13 of the rugarch vignette. This function should be given the relevant parameter values as explained on page 13-18 of the rugarch vignette. Here is an example using the Skewed Student t distribution in an ARMA(1,1)-GARCH(1,1) model on the sp500ret dataset:

```r
model = ugarchspec(
  variance.model = list(model = "sGARCH", garchOrder = c(1, 1)),
  mean.model = list(armaOrder = c(1, 1)),
  distribution.model = "sstd"
)
modelroll = ugarchroll(
  model, data = sp500ret, n.ahead = 1, forecast.length = 100,
  n.start = NULL, refit.every = 50, refit.window = c("recursive"),
  window.size = NULL, solver = "hybrid", fit.control = list(),
  solver.control = list(), calculate.VaR = TRUE, VaR.alpha = c(0.01, 0.05),
  cluster = NULL, keep.coef = TRUE
)

# getting the estimated skew from modelroll
skew.estimate = modelroll@forecast$density[, "Skew"]
# getting the estimated shape from modelroll
shape.estimate = modelroll@forecast$density[, "Shape"]
# using quantile function from rugarch package
# setting skew and shape according to estimates
# other values kept as specified in vignette
q_st = qdist(distribution = "sstd", p = 0.05, mu = 0,
             sigma = 1, lambda = -0.5, skew = skew.estimate, shape = shape.estimate)
mu = modelroll@forecast$density[, "Mu"]
sigma = modelroll@forecast$density[, "Sigma"]
VaR = sigma * q_st + mu
```

### 6.2 Different roll-procedures

Doing a forecast we could choose to specify `forecast.length`, `refit.every` and `refit.window`. Above we used the argument `forecast.length=100` to leave 100 observations out-of-sample leaving the rest 5423 observations in-sample. As a consequence the first model is estimated on the 5423 in-
sample observations. Also we chose to refit the model every 50 observations refit.every=50 meaning that the first model would be used for 50 forecasts, that is fore period $5423 + 1, 5423 + 2, ..., 5423 + 50$. Then the model is reestimated and another 50 forecasts are made for the periods $5473 + 1, 5473 + 2, ..., 5473 + 50$ after which forecasts have been made for all periods left as out of sample. Leaving out 100 observations but reestimating for every 51 observations results in the first model being used for 51 forecasts and the other for 49 and in general the last model estimated will be used for less forecasts. But on what observations are the models estimated? The window could be recursive in which case the model is estimated on all observations with lower index than the index of the first period for which the model is used to forecast. The second model in our example would be estimated on observations $1, 2, 3, ..., 5473$ however if instead using a moving window refit.window="moving" the second model would only be estimated on observations $51, ..., 5473$ moving forward the first observation by refit.every=50 every time a new model is estimated. Using the first procedure each model is estimated on an increasing sample size while using the second procedure keeps the sample size constant.

Instead of specifying forecast.length, refit.every and refit.window one could leave out forecast.length and instead choose n.start indication where in dataset to start rolling forecast. As an example lets say number of observations are $N = 1000$ and we choose n.start=800 then the forecast length is 200. If you at the same time specify the forecast.length=100 it will automatically be overwritten to have the value forecast.length=200. Otherwise the procedure is the same so with the first model i estimated on $1, ..., 800$ and with refit.every=50 used to forecast for $800 + 1, ..., 800 + 50$ and then the second model is estimated on $1, ..., 850$ if refit.window=c("recursive") and on $51, ..., 850$ if refit.window=c("moving").

Another way the determine the procedure is to specify forecast.length, refit.every and window.size and refit.window=c("window"). Assuming we have $N = 1000$ observations and forecast.length=100 so that 100 observations are left out-of-sample (as we now know only for the initial model so it would be more correct to say that choosing forecast.length determines for how many periods forecasts are going to be calculated thus explaining the name of the argument). Assume we choose refit.every=50 as a consequence the estimation of two models would be sufficient. Assume further that we choose a window.size=500 this result is that each model is estimated in 500 observations the first observation for the first model being $N$-forecast.length-window.size=1000-100-500=400. The first model is estimated on observations 400, 401, ..., 900 the second on 450, ..., 950 because refit.every=50. If refit.window=c("recursive") the window.size is
ignored so that the first model is estimated on 1, ..., 900 and the second on 1, ..., 950. But what if the total sample $N$ is too small for the first model to be estimated on a window with the specified window size? Let $N = 1000$ and \texttt{forecast.length=300} and \texttt{refit.every=50} then 700 observations are available to estimate the first model on and 750 for the second and 800 for the third etc. As a consequence choosing \texttt{window.size=770} will result in the first model being estimated on 1, ..., 700 (hence less than chosen because only 700 are available) and the second model will be estimated on observations 1, ..., 750 hence also less than chosen and the third model will be estimated on 30, 31, ..., 800.

Again remember that the created object can be inspected by \texttt{str()} so to investigate the consequences of different choices have a look at \texttt{str(modelroll @model)} to see how \texttt{$rollind} and \texttt{$out.sample} among others change depending on how you choose to specify the forecasting procedure.