# Ejemplos

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Paper: Modelos ocultos de Markov: una aplicación en series de tiempo

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https://github.com/lizbethna/HMMaplica

Este archivo muestra las instrucciones para correr los códigos de R y Stan.

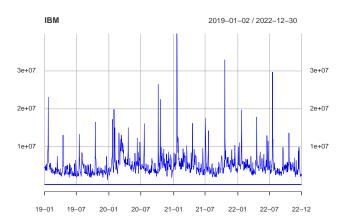
# Markov switching GARCH

```
library(ggplot2)
library(rstan) # RStan
library(quantmod) # Quantitative Financial Modelling Framework
```

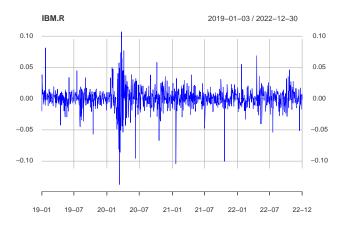
```
plot_statepath <- function(zstar) {</pre>
  K <- length(unique(as.vector(zstar)))</pre>
  x <- index(zstar)
  t <- 1:dim(zstar)[1]
  opar <- par(no.readonly = TRUE)</pre>
  zcol <- (1:K)[zstar]</pre>
  layout(matrix(c(1, 2), nrow = 2, ncol = 1), heights = c(0.95, 0.05))
  plot(x = x, y = zstar,
    xlab = bquote(t), ylab = bquote(hat(z)[t]),
    main = bquote("Secuencia mas probable de estados ocultos"),
    ylim = c(1, K), type = 'l', col = 'gray')
  points(x=x, y=zstar,
           pch = 21, bg = zcol, col = zcol, cex = 0.7)
  par(mai = c(0, 0, 0, 0))
  plot.new()
  legend(x = "center",
         legend = c('Trayectoria mas probable', paste('Estado', 1:K)),
         pch = c(NA, rep(21, K)),
         lwd = c(2, rep(NA, K)),
         col = c('lightgray', 1:K),
         pt.bg = c('lightgray', 1:K),
         bty = 'n', cex = 0.7,
```

```
horiz = TRUE)
par(opar)
}
```

## **Datos**







# Código Stan

```
# Markov-switching GARCH
msgarch_fit <- function(y) {
  rstan_options(auto_write = TRUE)
  options(mc.cores = parallel::detectCores())

stan.model = 'hmm_garch.stan'</pre>
```

```
y <- as.vector(coredata(y));</pre>
  stan.data = list(
   T = length(y),
   y = y
  stan(file = stan.model,
      data = stan.data, verbose = T,
       iter = 1000, warmup = 500,
       thin = 1, chains = 1,
       cores = 1, seed = 900)
}
# Fit GARCH
fit <- msgarch_fit(IBM.R)</pre>
TRANSLATING MODEL 'hmm_garch' FROM Stan CODE TO C++ CODE NOW.
successful in parsing the Stan model 'hmm_garch'.
CHECKING DATA AND PREPROCESSING FOR MODEL 'hmm_garch' NOW.
COMPILING MODEL 'hmm_garch' NOW.
STARTING SAMPLER FOR MODEL 'hmm_garch' NOW.
SAMPLING FOR MODEL 'hmm_garch' NOW (CHAIN 1).
Chain 1: Gradient evaluation took 0.001646 seconds
Chain 1: 1000 transitions using 10 leapfrog steps per transition would take 16.46 seconds.
Chain 1: Adjust your expectations accordingly!
Chain 1:
Chain 1:
Chain 1: Iteration:
                     1 / 1000 [ 0%]
                                        (Warmup)
Chain 1: Iteration: 100 / 1000 [ 10%]
                                        (Warmup)
Chain 1: Iteration: 200 / 1000 [ 20%]
                                        (Warmup)
Chain 1: Iteration: 300 / 1000 [ 30%]
                                        (Warmup)
Chain 1: Iteration: 400 / 1000 [ 40%]
                                        (Warmup)
Chain 1: Iteration: 500 / 1000 [ 50%]
                                        (Warmup)
Chain 1: Iteration: 501 / 1000 [ 50%]
                                        (Sampling)
Chain 1: Iteration: 600 / 1000 [ 60%]
                                        (Sampling)
Chain 1: Iteration: 700 / 1000 [ 70%]
                                        (Sampling)
Chain 1: Iteration: 800 / 1000 [ 80%]
                                        (Sampling)
Chain 1: Iteration: 900 / 1000 [ 90%]
                                        (Sampling)
Chain 1: Iteration: 1000 / 1000 [100%]
                                         (Sampling)
Chain 1:
Chain 1: Elapsed Time: 13.3548 seconds (Warm-up)
Chain 1:
                        12.1476 seconds (Sampling)
Chain 1:
                        25.5024 seconds (Total)
Chain 1:
```

## Resultados

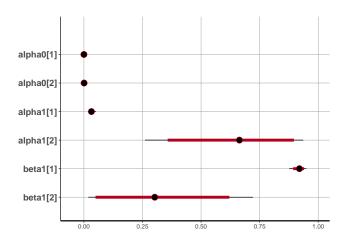
```
round(summary(fit, pars=c("alpha0","alpha1","beta1","A"))$summary,3)
          mean se_mean
                          sd 2.5%
                                      25%
                                            50%
                                                 75% 97.5%
                                                             n eff Rhat
alpha0[1] 0.000
                 0.000 0.000 0.000 0.000 0.000 0.000 0.000 329.528 1.001
alpha0[2] 0.001
                 0.000 0.000 0.000 0.001 0.001 0.001 0.002 293.842 0.998
                 0.000 0.009 0.017 0.027 0.032 0.038 0.051 309.071 0.999
alpha1[1] 0.033
alpha1[2] 0.646
                 0.012 0.200 0.261 0.482 0.663 0.817 0.936 296.904 0.998
beta1[1] 0.918
                 0.001 0.019 0.876 0.908 0.920 0.930 0.950 290.352 1.002
                 0.012 0.210 0.018 0.139 0.303 0.496 0.722 289.117 0.998
beta1[2] 0.321
A[1,1]
         0.937
                 0.001 0.016 0.902 0.927 0.938 0.950 0.963 447.746 1.000
A[1,2]
         0.063
                 0.001 0.016 0.037 0.050 0.062 0.073 0.098 447.746 1.000
A[2,1]
         0.542
                 0.005 0.099 0.383 0.463 0.543 0.610 0.752 375.599 1.004
A[2,2]
         0.458
                 0.005 0.099 0.248 0.390 0.457 0.537 0.617 375.599 1.004
round(summary(fit, pars=c("alpha0", "alpha1", "beta1", "A"))$c_summary,3)
```

#### , , chains = chain:1

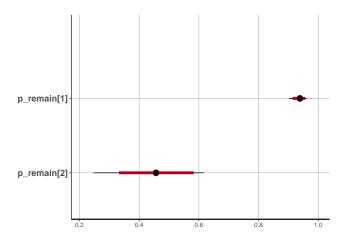
#### stats

```
parameter
                                25%
                                       50%
                                             75% 97.5%
             mean
                     sd 2.5%
  alpha0[1] 0.000 0.000 0.000 0.000 0.000 0.000 0.000
  alpha0[2] 0.001 0.000 0.000 0.001 0.001 0.001 0.002
  alpha1[1] 0.033 0.009 0.017 0.027 0.032 0.038 0.051
  alpha1[2] 0.646 0.200 0.261 0.482 0.663 0.817 0.936
  beta1[1]
           0.918 0.019 0.876 0.908 0.920 0.930 0.950
  beta1[2] 0.321 0.210 0.018 0.139 0.303 0.496 0.722
  A[1,1]
            0.937 \ 0.016 \ 0.902 \ 0.927 \ 0.938 \ 0.950 \ 0.963
            0.063 0.016 0.037 0.050 0.062 0.073 0.098
  A[1,2]
  A[2,1]
            0.542 0.099 0.383 0.463 0.543 0.610 0.752
  A[2,2]
            0.458 0.099 0.248 0.390 0.457 0.537 0.617
```

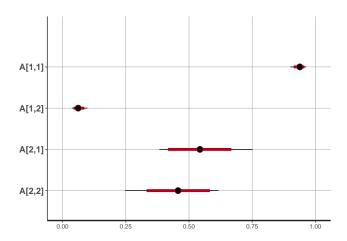
#### plot(fit,pars=c("alpha0","alpha1","beta1"))

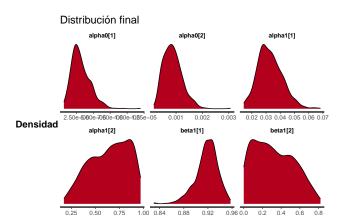


# plot(fit,pars="p\_remain")

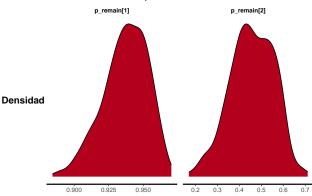


# plot(fit,pars="A")

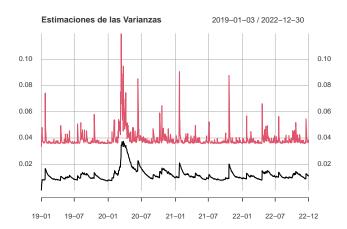


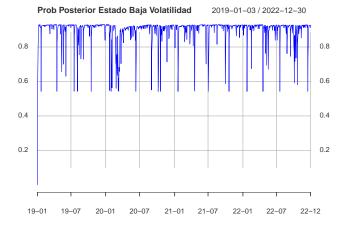


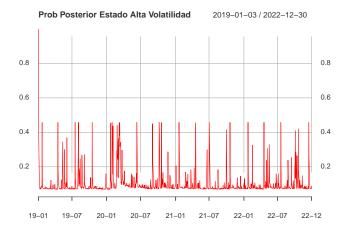
#### Distribución final de p\_remain



# Densidad | Densidad |







## Secuencia mas probable de estados ocultos

