

Stability properties for MAR(1) models

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Background

There is growing interest in the use of first-order vector autoregressive, or VAR(1), models in ecology where they are often referred to as multivariate autoregressive, or MAR(1), models (*e.g.*, Ives *et al.* 2003 *Ecological Monographs* 73:301–330).

Assume a MAR(1) model of the general form

$$\mathbf{x}_t = \mathbf{a} + \mathbf{B}(\mathbf{x}_{t-1} - \mathbf{a}) + \mathbf{w}_t$$

where \mathbf{x}_t is an $n \times 1$ vector of state variates at time t , \mathbf{a} is an $n \times 1$ vector of underlying levels (means) for each of the states, \mathbf{B} is an $n \times n$ interaction matrix, and \mathbf{w}_t is an $n \times 1$ vector of multivariate normal process errors; $\mathbf{w}_t \sim \text{MVN}(\mathbf{0}, \mathbf{Q})$.

Of particular interest here is that the variance-covariance matrix of the stationary distribution for \mathbf{x}_t as $t \rightarrow \infty$ gives an indication of the relative stability of the system.

State-space form

This MAR(1) model can be used within a state-space framework, wherein \mathbf{a} is instead included as part of a second model for the observed data \mathbf{y} , such that

$$\begin{aligned}\mathbf{x}_t &= \mathbf{B}\mathbf{x}_{t-1} + \mathbf{w}_t, \\ \mathbf{y}_t &= \mathbf{x}_t + \mathbf{a} + \mathbf{v}_t,\end{aligned}$$

and \mathbf{v}_t is an $n \times 1$ vector of observation errors, the statistical distribution of which does not affect the variance of the stationary distribution.

Variance of the stationary distribution

We will restrict this treatment to stationary models wherein all of the eigenvalues of \mathbf{B} lie within the unit circle. Because $t = (t - 1)$ as $t \rightarrow \infty$, under assumptions of stationarity we can write the process equation from the above state-space model as

$$\mathbf{x}_t = \mathbf{B}\mathbf{x}_t + \mathbf{w}_t.$$

From this, it follows that

$$\text{Var}(\mathbf{x}_t) = \mathbf{B}\text{Var}(\mathbf{x}_t)\mathbf{B}^\top + \text{Var}(\mathbf{w}_t).$$

If we define $\mathbf{\Sigma} = \text{Var}(\mathbf{x}_t)$, then

$$\mathbf{\Sigma} = \mathbf{B}\mathbf{\Sigma}\mathbf{B}^\top + \mathbf{Q}.$$

Unfortunately, however, there is no closed-form solution for $\mathbf{\Sigma}$ when written in this form.

The *vec* operator

It turns out that we can use the *vec* operator to derive an explicit solution for $\mathbf{\Sigma}$. The *vec* operator converts an $i \times j$ matrix into an $(ij) \times 1$ column vector. For example, if

$$\mathbf{M} = \begin{bmatrix} 1 & 3 \\ 2 & 4 \end{bmatrix},$$

then

$$\text{vec}(\mathbf{M}) = \begin{bmatrix} 1 \\ 2 \\ 3 \\ 4 \end{bmatrix}.$$

Solution

Thus, if \mathbf{I} is an $n \times n$ identity matrix, and we define $\mathcal{I} = (\mathbf{I} \otimes \mathbf{I})$ and $\mathcal{B} = (\mathbf{B} \otimes \mathbf{B})$, then

$$\text{vec}(\mathbf{\Sigma}) = (\mathcal{I} - \mathcal{B})^{-1}\text{vec}(\mathbf{Q}).$$

Importance of species interactions to stability

Among the many ways of classifying stability, I am interested in the extent to which community interactions, relative to environmental forcing, contribute to the overall variance of the stationary distribution. In a stable system, any perturbation affecting one or more of the community members does not amplify as it moves throughout the community as a whole, such that the variances in log-densities over time would be driven almost entirely by random environmental variation. Given that same magnitude of random environmental variation, less stable systems are characterized by greater variances in the temporal dynamics of their constituents.

Here we will use determinants to measure the “volume” of a matrix. Looking back to the matrix form of the equation for the variance of the stationary distribution, we see that

$$\mathbf{\Sigma} = \mathbf{B}\mathbf{\Sigma}\mathbf{B}^\top + \mathbf{Q},$$

and hence

$$\Sigma - \mathbf{Q} = \mathbf{B}\Sigma\mathbf{B}^\top.$$

Therefore, the volume of the difference $\Sigma - \mathbf{Q}$ provides a measure of how much species interactions contribute to the variance of the stationary distribution. Taking determinants of both sides, we get

$$\begin{aligned}\det(\Sigma - \mathbf{Q}) &= \det(\mathbf{B}\Sigma\mathbf{B}^\top) \\ &= \det(\mathbf{B})\det(\Sigma)\det(\mathbf{B}^\top) \\ &= \det(\mathbf{B})\det(\Sigma)\det(\mathbf{B}).\end{aligned}$$

The proportion $\pi_{\mathbf{B}}$ of the volume of Σ attributable to species interactions is then

$$\pi_{\mathbf{B}} = \frac{\det(\Sigma - \mathbf{Q})}{\det(\Sigma)} = \det(\mathbf{B})^2.$$

The effects of covariates

Often we would like to examine the potential effects of covariates on community dynamics. In that case, the MARSS(1) model becomes

$$\begin{aligned}\mathbf{x}_t &= \mathbf{B}\mathbf{x}_{t-1} + \mathbf{C}\mathbf{c}_t + \mathbf{w}_t, \\ \mathbf{y}_t &= \mathbf{x}_t + \mathbf{a} + \mathbf{v}_t,\end{aligned}$$

where \mathbf{C} is an $n \times p$ matrix of covariate effects, and \mathbf{c}_t is a $p \times 1$ vector of covariates at time t . Although the variance of the stationary distribution will be affected by any covariate effects, the method described above for estimating the relative effects of species interactions is unaffected. That is, if Ξ is the variance-covariance matrix of the covariates in \mathbf{c} , then analogous to above the variance of the stationary distribution will be

$$\begin{aligned}\Sigma &= \mathbf{B}\Sigma\mathbf{B}^\top + \mathbf{C}\Xi\mathbf{C}^\top + \mathbf{Q}, \\ \Sigma - \mathbf{C}\Xi\mathbf{C}^\top - \mathbf{Q} &= \mathbf{B}\Sigma\mathbf{B}^\top,\end{aligned}$$

and hence the proportion $\pi_{\mathbf{B}}$ of the volume of Σ attributable to species interactions is, as before, given by

$$\pi_{\mathbf{B}} = \frac{\det(\Sigma - \mathbf{C}\Xi\mathbf{C}^\top - \mathbf{Q})}{\det(\Sigma)} = \det(\mathbf{B})^2.$$

Effect of density-dependence

In a MAR(1) model of community dynamics, the matrix \mathbf{B} maps the vector of log-densities from one time step to another. In particular, the diagonal elements of \mathbf{B} control the degree of density-dependence via the degree of so-called mean reversion. Because we are interested in stationary dynamics, the diagonals of \mathbf{B} will all be less than 1 in absolute value. As such, here are a few generalizations:

1. As $|b_{ii}| \rightarrow 0$ the strength of density dependence increases (*i.e.*, the degree of mean reversion increases);
2. As $b_{ii} \rightarrow 1$ the strength of density dependence decreases to the point where there is no density dependence when $b_{ii} = 1$ (*i.e.*, the temporal dynamics become a non-stationary random walk); and
3. Negative values of \mathbf{B} ($-1 < b_{ii} < 0$) result in a time series with a 2-point oscillation.

Furthermore, the determinant (and the trace) of a matrix are functions of the eigenvalues λ_i , such that for an $N \times N$ matrix \mathbf{B} :

$$\det(\mathbf{B}) = \prod_{i=1}^N \lambda_i.$$

The eigenvalues themselves are a function of all of the elements in \mathbf{B} , but they are particularly sensitive to the elements along the diagonal. We can formally examine the sensitivity of $\det(\mathbf{B})$ to each of the elements in \mathbf{B} via the following relationship:

$$\frac{\partial \det(\mathbf{B})}{\partial \mathbf{B}_{ij}} = \det(\mathbf{B})(\mathbf{B}^{-1})_{ji}.$$

Thus, for any matrix \mathbf{B} we can ask whether its determinant is most sensitive to a diagonal versus an off-diagonal element.

A simulation study

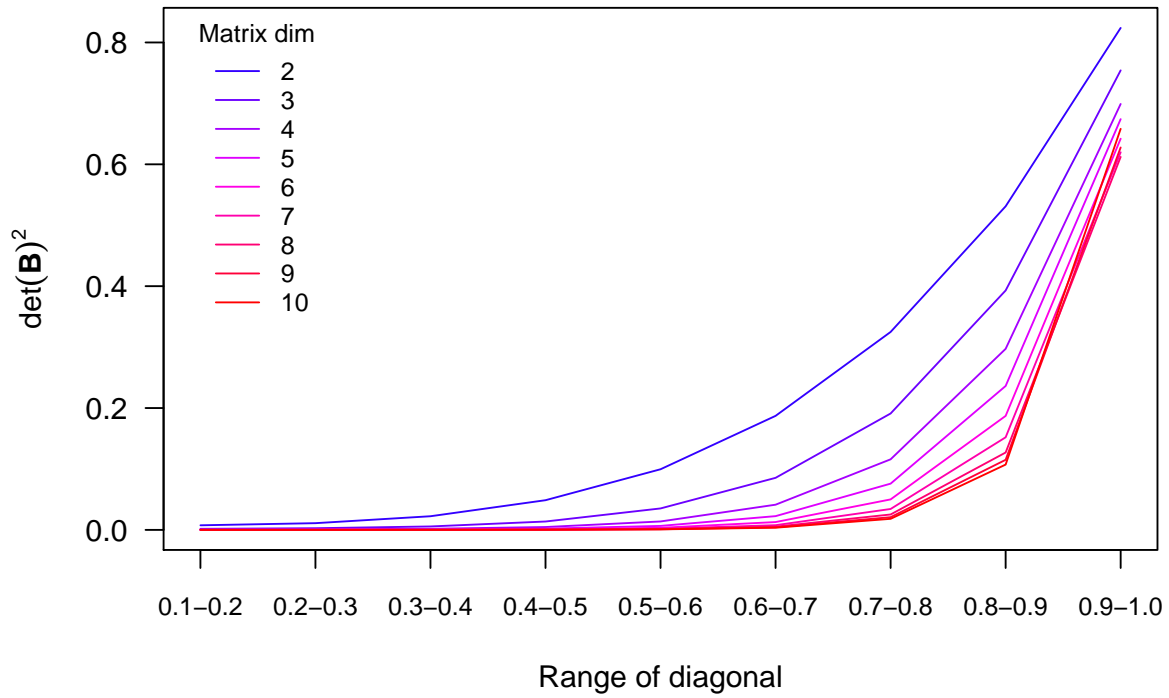
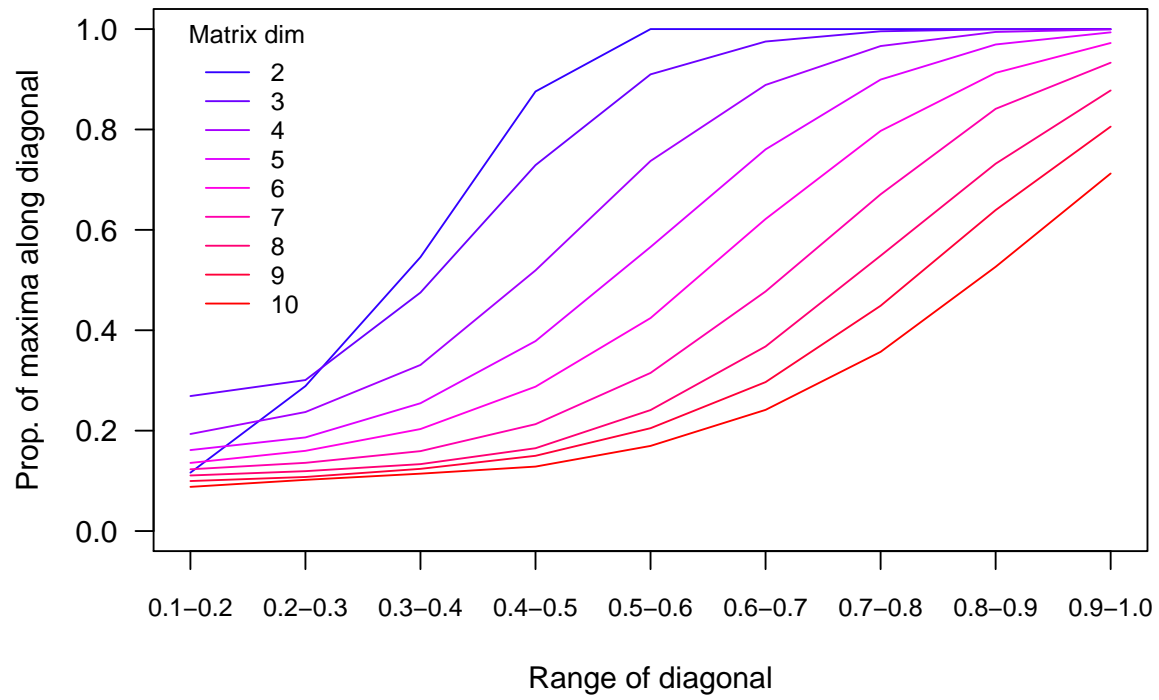
Let's examine the determinant and the location of its maximum partial derivative for whole bunch of matrices. We'll do so over a range of matrix sizes and entries. Here's the pseudo code:

- Choose a matrix size
- Do the following many times
 - Populate the entire matrix with random values over a specified range
 - Replace the diagonal with random values over a specified range
 - Calculate partial derivatives and note location of max in abs value
 - For the same matrix, increase the range of values in the diagonal
 - Calculate partial derivatives and note location of max in abs value
 - Calculate $\det(\mathbf{B})^2$

```

## maximum matrix dim
pp <- 10
## number of samples
mc <- 10000
## maximum diagonal value (x10)
dd <- 10
## min off-diagonal
dn <- -0.5
## min off-diagonal
dx <- 0.5
## sample over matrix size and diagonal
sam <- dtm <- matrix(NA,mc,dd-1)
res1 <- res2 <- matrix(NA,dd-1,pp-1)
for(ii in 2:pp) {
  for(jj in 1:mc) {
    mat <- matrix(runif(ii^2,dn,dx),ii,ii)
    for(kk in 2:dd) {
      tmp <- mat
      diag(tmp) <- runif(ii,kk-1,kk)/10
      rc <- which(abs(t(solve(tmp)))==max(abs(t(solve(tmp)))))
      sam[jj,kk-1] <- rc %in% seq(1,ii^2,ii+1)
      dtm[jj,kk-1] <- det(tmp)^2
    }
  }
  res1[,ii-1] <- apply(sam, 2, sum)/mc
  res2[,ii-1] <- apply(dtm, 2, mean)
}
## plot the results
par(mfrow=c(2,1), mai=c(0.9,0.9,0.1,0.1), omi=rep(0,4))
matplot(res1, type="l", lty="solid", col=rainbow(pp-1, start=0.7, end=1),
        ylim=c(0,1), las=1,
        xaxt="n", ylab="Prop. of maxima along diagonal", xlab="Range of diagonal")
axis(1, at=seq(pp-1), cex.axis=0.8, labels=c("0.1-0.2", "0.2-0.3", "0.3-0.4",
        "0.4-0.5", "0.5-0.6", "0.6-0.7",
        "0.7-0.8", "0.8-0.9", "0.9-1.0"))
legend(x="topleft", legend = seq(2,pp), col=rainbow(pp-1, start=0.7, end=1),
       lty="solid", bty="n", title="Matrix dim", cex=0.8, inset=c(0.03,0.01))
matplot(res2, type="l", lty="solid", col=rainbow(pp-1, start=0.7, end=1),
        las=1,
        xaxt="n", ylab=expression(det(bold(B))^2), xlab="Range of diagonal")
axis(1, at=seq(pp-1), cex.axis=0.8, labels=c("0.1-0.2", "0.2-0.3", "0.3-0.4",
        "0.4-0.5", "0.5-0.6", "0.6-0.7",
        "0.7-0.8", "0.8-0.9", "0.9-1.0"))
legend(x="topleft", legend = seq(2,pp), col=rainbow(pp-1, start=0.7, end=1),
       lty="solid", bty="n", title="Matrix dim", cex=0.8, inset=c(0.03,0.01))

```



An example

As an example, I will fit a MARSS(1) model to some plankton data from Lake Washington (USA). These data were made publicly available by Dr. Daniel Schindler of the University of Washington,

and they have been included as part of the MARSS package for **R**. To speed up model fitting, I will focus on a representative decade from 1980 through 1989, and only use some of the plankton groups.

```
library(MARSS)
## load the data (2 datasets here)
data(lakeWAp plankton)
raw <- as.data.frame(lakeWAp planktonRaw)
## keep only 1977-
dat <- subset(raw, Year>=1980 & Year<=1989)
## contents
colnames(dat)

## [1] "Year" "Month"
## [3] "Temp" "TP"
## [5] "pH" "Cryptomonas"
## [7] "Diatoms" "Greens"
## [9] "Bluegreens" "Unicells"
## [11] "Other.algae" "Conochilus"
## [13] "Cyclops" "Daphnia"
## [15] "Diaptomus" "Epischura"
## [17] "Leptodora" "Neomysis"
## [19] "Non.daphnid.cladocerans" "Non.colonial.rotifers"

## phytoplankton names (omitting bluegreens)
phyto_names <- c("Cryptomonas", "Diatoms", "Greens")
## zooplankton names (omitting Leptodora & Neomysis)
zoops_names <- c("Cyclops", "Diaptomus", # copepods
                 "Daphnia", "Non.daphnid.cladocerans") # cladocerans
all_names <- c(phyto_names, zoops_names)
## number of taxa
nP <- length(phyto_names)
nZ <- length(zoops_names)
nn <- nP + nZ
## length of ts
TT <- dim(dat)[1]
```

In addition to the biotic interactions, I will include the effects of water temperature on the plankton.

Model definition

Now we need to set up the various vectors and matrices that define our MARSS(1) model. Here I will make the following assumptions:

1. independent and identically distributed observation errors for all of the taxa within each of the two plankton groups;
2. a block-diagonal form for **Q** wherein all of the taxa within a plankton group have the same variance, and no covariance within and among blocks; and
3. fixed monthly effects to eliminate seasonal components.

The following code also makes use of various shorthand character description of vectors and matrices in MARSS.

```

mod_list <- list()
## process model
mod_list$B <- "unconstrained"
mod_list$U <- "zero"
## include seasonal effect
mod_list$C <- "unconstrained"
mod_list$c <- matrix(NA,2,TT)
mod_list$c[1,] <- cos(2*pi*seq(TT)/12)
mod_list$c[2,] <- sin(2*pi*seq(TT)/12)
## var-cov of process (environmental errors)
mod_list$Q <- "diagonal and unequal"
## observation model
mod_list$Z <- "identity"
## set levels to zero & subtract the mean from the data below
mod_list$A <- "zero"
mod_list$D <- "zero"
mod_list$d <- "zero"
## observation var-cov
mod_list$R <- matrix(list(0), nn, nn)
diag(mod_list$R) <- c(rep("r_phyto", nP), rep("r_zoops", nZ))

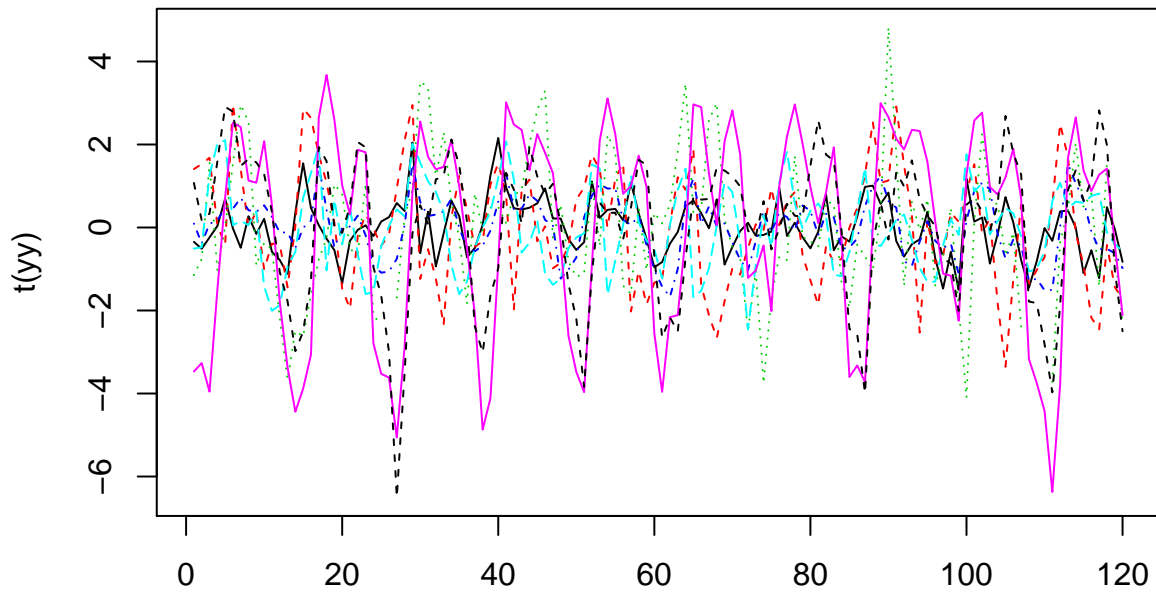
```

The model we are fitting is based upon log-densities, so we'll need to transform the data before proceeding. In this case, zeroes in the data set are indicative of the sampling program and do not indicate true absences; therefore I will convert them to NA.

```

yy <- dat[,all_names]
yy[yy==0] <- NA
## log-transform variates
## data must also be n x T, so transpose
yy <- t(log(yy))
## subtract the mean from each ts
yy <- yy - apply(yy, 1, mean, na.rm=TRUE)
matplot(t(yy), type="l")

```

Now we can fit the model. Note that the output from MARSS is rather voluminous, so I will suppress it here.

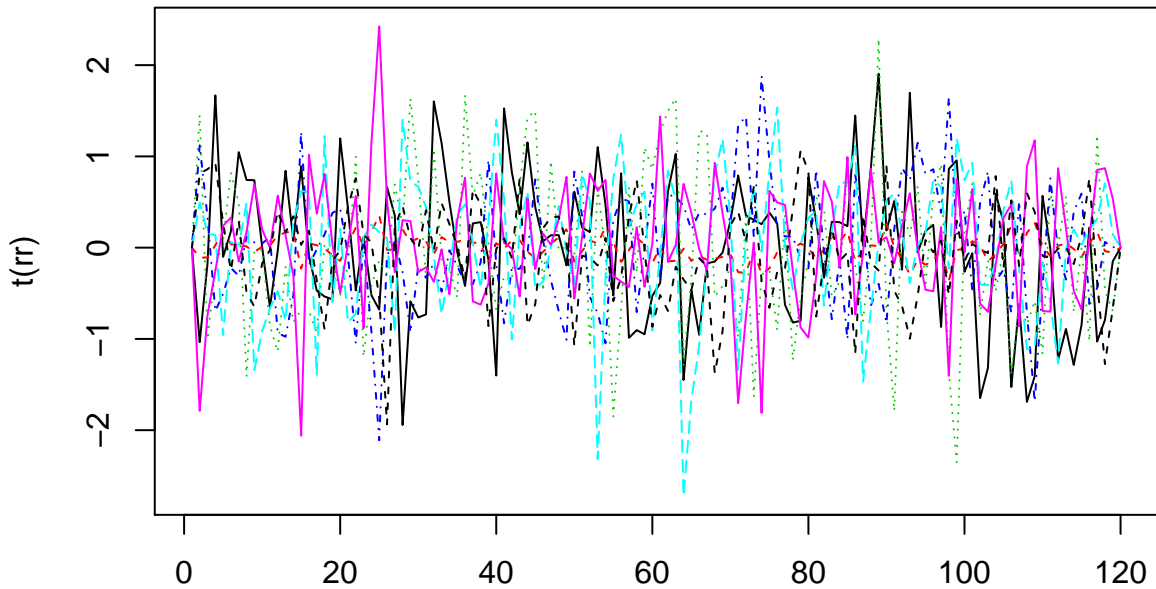
```
## fit model
mod <- MARSS(yy, model=mod_list, control=list(maxit=5000, safe=TRUE))

## Warning! Abstol convergence only. Maxit (=5000) reached before log-log convergence.
##
## MARSS fit is
## Estimation method: kem
## Convergence test: conv.test.slope.tol = 0.5, abstol = 0.001
## WARNING: Abstol convergence only no log-log convergence.
## maxit (=5000) reached before log-log convergence.
## The likelihood and params might not be at the ML values.
## Try setting control$maxit higher.
## Log-likelihood: -962.6041
## AIC: 2083.208 AICc: 2099.928
##
##
##                                     Estimate
## R.r_phyto                          0.369889
## R.r_zoops                          0.160168
## B.(1,1)                            1.036310
## B.(2,1)                            3.563123
## B.(3,1)                            0.770864
## B.(4,1)                            1.365007
## B.(5,1)                           -0.695990
## B.(6,1)                            2.538670
## B.(7,1)                            1.328608
## B.(1,2)                           -0.172489
## B.(2,2)                            0.474807
## B.(3,2)                           -0.011794
## B.(4,2)                           -0.382594
## B.(5,2)                           -0.269544
```

## B.(6,2)	-2.793273
## B.(7,2)	-2.271962
## B.(1,3)	0.003898
## B.(2,3)	0.078826
## B.(3,3)	0.396992
## B.(4,3)	0.033682
## B.(5,3)	0.013971
## B.(6,3)	-0.039329
## B.(7,3)	-0.016341
## B.(1,4)	0.003095
## B.(2,4)	-2.855293
## B.(3,4)	-0.428257
## B.(4,4)	0.868483
## B.(5,4)	1.181447
## B.(6,4)	8.890468
## B.(7,4)	7.359708
## B.(1,5)	0.018770
## B.(2,5)	-0.204729
## B.(3,5)	0.122953
## B.(4,5)	-0.025568
## B.(5,5)	0.262689
## B.(6,5)	0.332045
## B.(7,5)	-0.067550
## B.(1,6)	-0.089099
## B.(2,6)	-0.000669
## B.(3,6)	0.123341
## B.(4,6)	-0.166642
## B.(5,6)	-0.428653
## B.(6,6)	-0.570045
## B.(7,6)	-1.169429
## B.(1,7)	-0.014456
## B.(2,7)	0.152225
## B.(3,7)	0.031708
## B.(4,7)	-0.019874
## B.(5,7)	0.065571
## B.(6,7)	-0.432458
## B.(7,7)	0.404846
## Q.(X.Cryptomonas,X.Cryptomonas)	0.030365
## Q.(X.Diatoms,X.Diatoms)	0.001534
## Q.(X.Greens,X.Greens)	1.274952
## Q.(X.Cyclops,X.Cyclops)	0.004020
## Q.(X.Diaptomus,X.Diaptomus)	0.236876
## Q.(X.Daphnia,X.Daphnia)	0.161245
## Q.(X.Non.daphnid.cladocerans,X.Non.daphnid.cladocerans)	0.099368
## x0.X.Cryptomonas	-0.182630
## x0.X.Diatoms	-1.447394
## x0.X.Greens	-0.171765
## x0.X.Cyclops	-0.883487
## x0.X.Diaptomus	-2.832596
## x0.X.Daphnia	-2.355732
## x0.X.Non.daphnid.cladocerans	2.421227

```
## C.(X.Cryptomonas,1) -0.127349
## C.(X.Diatoms,1) -0.538701
## C.(X.Greens,1) -0.749682
## C.(X.Cyclops,1) -0.180890
## C.(X.Diaptomus,1) -0.631246
## C.(X.Daphnia,1) 1.180092
## C.(X.Non.daphnid.cladocerans,1) 0.727710
## C.(X.Cryptomonas,2) 0.012359
## C.(X.Diatoms,2) -0.177790
## C.(X.Greens,2) -0.068852
## C.(X.Cyclops,2) 0.017268
## C.(X.Diaptomus,2) -0.060490
## C.(X.Daphnia,2) 0.854117
## C.(X.Non.daphnid.cladocerans,2) 0.207171
##
## Standard errors have not been calculated.
## Use MARSSparamCIs to compute CIs and bias estimates.
##
## Convergence warnings
## Warning: the Q.(X.Diatoms,X.Diatoms) parameter value has not converged.
## Type MARSSinfo("convergence") for more info on this warning.

rr <- residuals(mod)$state.residuals
matplot(t(rr), type="l")
```



Stability analysis

Now we can calculate the proportion π_B of the variance of the stationary distribution Σ owing to species interactions.

```
## get B matrix
BB <- coef(mod, type="matrix")$B
```

```

round(BB,2)

##      [,1] [,2] [,3] [,4] [,5] [,6] [,7]
## [1,] 1.04 -0.17 0.00 0.00 0.02 -0.09 -0.01
## [2,] 3.56 0.47 0.08 -2.86 -0.20 0.00 0.15
## [3,] 0.77 -0.01 0.40 -0.43 0.12 0.12 0.03
## [4,] 1.37 -0.38 0.03 0.87 -0.03 -0.17 -0.02
## [5,] -0.70 -0.27 0.01 1.18 0.26 -0.43 0.07
## [6,] 2.54 -2.79 -0.04 8.89 0.33 -0.57 -0.43
## [7,] 1.33 -2.27 -0.02 7.36 -0.07 -1.17 0.40

## calc pi_B
det(BB)^2

## [1] 0.0003752867

## Ives adjustment for other community comparisons
det(BB)^(2/nn)

## [1] 0.3240584

```

It looks like π_B is very small, which suggests that most of the temporal dynamics are driven by environmental forcing.

From this we can also calculate the ratio ρ of the importance of density-dependent interactions relative to other interactions.

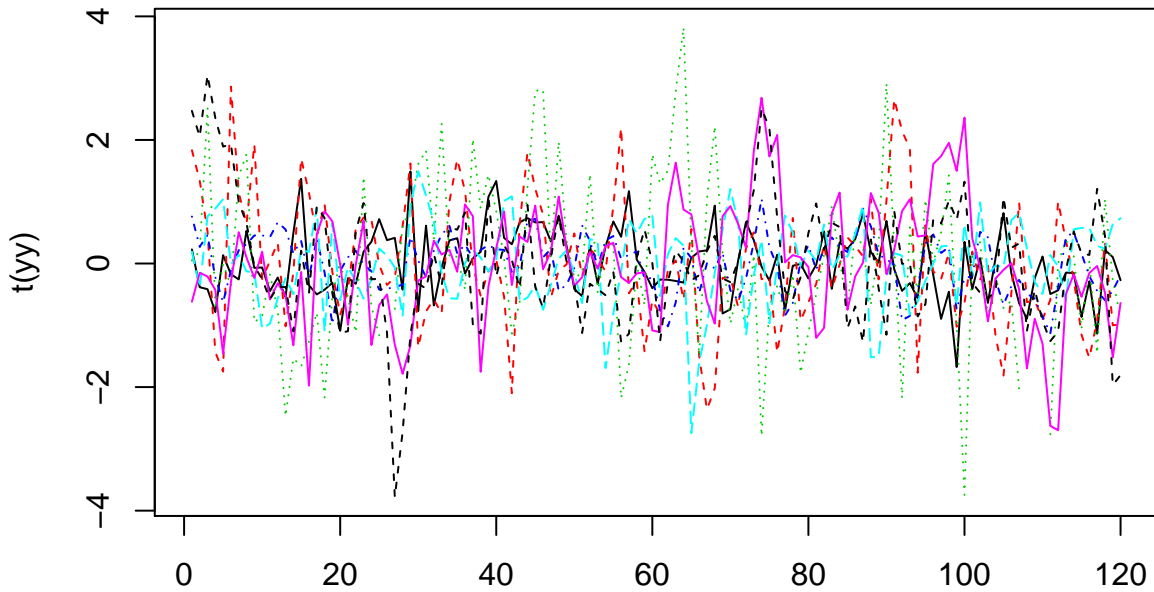
Example via decomposition

Now let's fit the MARSS model to the L Washington plankton data after we have removed any seasonal effect via decomposition (*i.e.*, removed the monthly means), rather than fitting some dummy covariates for season.

```

yy <- dat[,all_names]
yy[yy==0] <- NA
## log-transform variates
## data must also be n x T, so transpose
yy <- t(log(yy))
## subtract monthly means
for(i in 1:12) {
  idx <- seq(from=i, by=12, length.out=TT/12)
  mm <- apply(yy[,idx], 1, mean, na.rm=TRUE)
  yy[,idx] <- yy[,idx] - mm
}
## plot the decomposed data
matplot(t(yy), type="l")

```



Now we can fit the model. Note that I set **C** and **c** to "zero" because we have eliminated the seasonal effect.

```
mod_list$C <- "zero"
mod_list$c <- "zero"
## fit model
mod <- MARSS(yy, model=mod_list, control=list(maxit=5000, safe=TRUE))
```

```
## Warning! Abstol convergence only. Maxit (=5000) reached before log-log convergence.
##
## MARSS fit is
## Estimation method: kem
## Convergence test: conv.test.slope.tol = 0.5, abstol = 0.001
## WARNING: Abstol convergence only no log-log convergence.
## maxit (=5000) reached before log-log convergence.
## The likelihood and params might not be at the ML values.
## Try setting control$maxit higher.
## Log-likelihood: -859.1017
## AIC: 1848.203   AICc: 1859.346
##
##
##                                     Estimate
## R.r_phyto                        2.92e-01
## R.r_zoops                        5.31e-04
## B.(1,1)                          1.29e-01
## B.(2,1)                        -1.64e+00
## B.(3,1)                        -6.16e-01
## B.(4,1)                        -1.25e+00
## B.(5,1)                         1.01e-01
## B.(6,1)                        -7.98e+00
## B.(7,1)                        -5.84e+00
## B.(1,2)                         1.04e-02
## B.(2,2)                         4.36e-01
## B.(3,2)                         6.02e-02
```

## B.(4,2)	5.11e-02
## B.(5,2)	-4.18e-02
## B.(6,2)	1.79e-01
## B.(7,2)	-2.81e-02
## B.(1,3)	-9.66e-03
## B.(2,3)	1.70e-02
## B.(3,3)	4.56e-01
## B.(4,3)	3.24e-02
## B.(5,3)	1.41e-02
## B.(6,3)	7.74e-02
## B.(7,3)	3.88e-02
## B.(1,4)	5.93e-03
## B.(2,4)	4.54e-02
## B.(3,4)	-1.44e-01
## B.(4,4)	5.62e-01
## B.(5,4)	-1.57e-01
## B.(6,4)	1.60e-01
## B.(7,4)	3.84e-01
## B.(1,5)	-1.83e-02
## B.(2,5)	-4.29e-02
## B.(3,5)	-9.26e-02
## B.(4,5)	-4.89e-02
## B.(5,5)	2.68e-01
## B.(6,5)	-7.26e-02
## B.(7,5)	-4.02e-02
## B.(1,6)	-1.26e-02
## B.(2,6)	-1.44e-01
## B.(3,6)	-2.30e-02
## B.(4,6)	-5.06e-02
## B.(5,6)	-1.12e-01
## B.(6,6)	4.06e-01
## B.(7,6)	-2.13e-01
## B.(1,7)	7.03e-03
## B.(2,7)	-4.08e-03
## B.(3,7)	1.05e-02
## B.(4,7)	-5.74e-02
## B.(5,7)	9.06e-02
## B.(6,7)	6.46e-02
## B.(7,7)	6.42e-01
## Q.(X.Cryptomonas,X.Cryptomonas)	6.07e-03
## Q.(X.Diatoms,X.Diatoms)	5.43e-01
## Q.(X.Greens,X.Greens)	1.18e+00
## Q.(X.Cyclops,X.Cyclops)	1.13e-01
## Q.(X.Diaptomus,X.Diaptomus)	4.00e-01
## Q.(X.Daphnia,X.Daphnia)	1.66e-01
## Q.(X.Non.daphnid.cladocerans,X.Non.daphnid.cladocerans)	3.62e-01
## x0.X.Cryptomonas	-1.55e+00
## x0.X.Diatoms	-9.58e+00
## x0.X.Greens	-4.71e+00
## x0.X.Cyclops	-5.32e+00
## x0.X.Diaptomus	-7.71e+00

```

## x0.X.Daphnia -2.36e+01
## x0.X.Non.daphnid.cladocerans -1.55e+01
##
## Standard errors have not been calculated.
## Use MARSSparamCIs to compute CIs and bias estimates.
##
## Convergence warnings
## 12 warnings. First 10 shown. Type cat(object$errors) to see the full list.
## Warning: the R.r_zoops parameter value has not converged.
## Warning: the Q.(X.Cryptomonas,X.Cryptomonas) parameter value has not converged.
## Warning: the Q.(X.Daphnia,X.Daphnia) parameter value has not converged.
## Warning: the x0.X.Cryptomonas parameter value has not converged.
## Warning: the x0.X.Diatoms parameter value has not converged.
## Warning: the x0.X.Greens parameter value has not converged.
## Warning: the x0.X.Cyclops parameter value has not converged.
## Warning: the x0.X.Diaptomus parameter value has not converged.
## Warning: the x0.X.Daphnia parameter value has not converged.
## Warning: the x0.X.Non.daphnid.cladocerans parameter value has not converged.
## get B matrix
BB <- coef(mod, type="matrix")$B
round(BB,2)

##      [,1] [,2] [,3] [,4] [,5] [,6] [,7]
## [1,] 0.13 0.01 -0.01 0.01 -0.02 -0.01 0.01
## [2,] -1.64 0.44 0.02 0.05 -0.04 -0.14 0.00
## [3,] -0.62 0.06 0.46 -0.14 -0.09 -0.02 0.01
## [4,] -1.25 0.05 0.03 0.56 -0.05 -0.05 -0.06
## [5,] 0.10 -0.04 0.01 -0.16 0.27 -0.11 0.09
## [6,] -7.98 0.18 0.08 0.16 -0.07 0.41 0.06
## [7,] -5.84 -0.03 0.04 0.38 -0.04 -0.21 0.64
## calc pi_B
det(BB)^2

## [1] 1.444155e-07
## Ives adjustment for other community comparisons
det(BB)^(2/nn)

## [1] 0.1053906

```