Unstable States: Community dynamics in a chaotic world

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Eigenvalues indicate system dynamics.

A system is stable if it eventually returns to a fixed point after a disturbance.

A system is unstable if it is not stable.

For imaginary (complex) eigenvalues:

- Negative Real Part Stable. Damping is a requirement of stability.
- $\bullet\,$ Positive Real Part Unstable. Increasing amplitude is unstable.
- Zero Real Part Unstable. Oscillators are unstable, as they will not go back to a steady state post-disturbance
- The complex part will not affect the stability.

Real eigenvalues:

• Zero eigenvalues: unstable

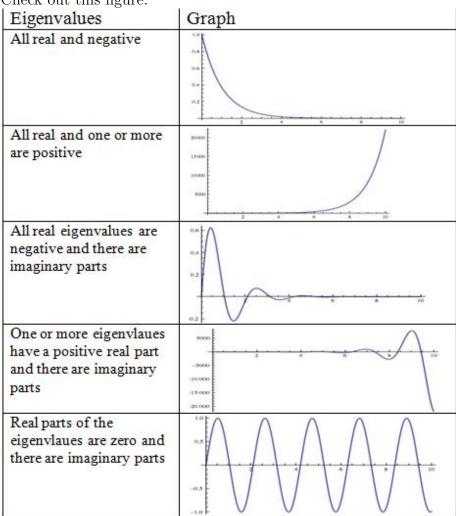
• All Positive: unstable

• All Negative: unstable

• Both positive and negative: unstable, saddle.

• Repeated: needs further analysis. For the special case of two eigenvalues, both positive = unstable, both negative = stable.

Check out this figure:



```
> source('../src/sens2_ews.R')
> library(ecodist)
> fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second
> x <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second pa
> x <- x[,-1]
> x <- x[,-1]
> x <- x[1:12500,]
> x[x<0] <- 0
> x <- t(x)
>
> ##Obtain equations
>
```

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>

Foundations of Resilience Thinking

- Resilience = size and steepness of attractor basins
- Distinguishes between: near equilibrium behavior and long term persistence
- Panarchy = Processes operating at different scales permit organization

How does NMDS treat mutual information?

```
> source('../src/sens2_ews.R')
> library(ecodist)
```

> fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second

```
> x <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second ps
> x <- x[,-1]
> x <- x[1:12500,]
> x[x<0] <- 0
> x <- t(x)
> ###kernel density information
> d1 <- density(x[1,])
> d2 <- density(x[2,])
> plot(density(x[1,])$x,density(x[3,])$x)
> points(-2:12,-2:12)
> ###
>
```

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Mutual Information for Time Series

This is an attempt to implement methods from Moniz et al. (2007) Application of Information Theory Methods to Food Web Reconstruction

```
> source('../src/sens2_ews.R')
> library(entropy)
> library(vegan)
> library(ecodist)
```

```
> fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second
> x <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second page 1.00 page
> x < -x[,-1]
> x <- x[1:12500,]
> x[x<0] <- 0
> x \leftarrow t(x)
> ###
> ## mid.m <- array(0,dim=c(nrow(x),nrow(x)))
> ## for (i in 1:nrow(x)){
> ## for (j in 1:nrow(x)){
                           mid.m[i,j] \leftarrow mi.Dirichlet(x[c(i,j),],0.5)
> ##
> ##
                        }
> ## }
> mid.m <- dget(file='../data/midm.rda')</pre>
> mid.m <- round(mid.m,7)</pre>
> pca <- princomp(as.dist(mid.m))</pre>
> nms <- nmds.min(nmds(as.dist(mid.m),2,2))</pre>
> plot(nms)
> plot(envfit(env=fps,ord=nms))
> envfit(env=fps,ord=nms)
>
              library(entropy)
> y2d <- rbind( sample(1:100,100), sample(1:100,100) )</pre>
> y2d. = rbind( sample(1:100,100), sample(1:100,100) )
> test <- test. <- 0
```

```
> a <- seq(0,1,by=0.01)
> for (i in 1:length(a)){test[i] <- mi.Dirichlet(y2d,a=a[i])}
> for (i in 1:length(a)){test.[i] <- mi.Dirichlet(y2d.,a=a[i])}
> plot(test.~test);abline(lm(range(test.)~range(test)))
>
```

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Reading Beltrami:

```
> ##For N. = rN
> ##N(t) = N(0)e^rt
>
> NO <- seq(0,100,by=0.1)
> t <- 0:100
> Nt <- sapply(NO, function(NO,t,r) NO*exp(r*t),t=t,r=2)
> plot(Nt[,1]~t,xlim=range(t),ylim=range(Nt),type='1',)
> for (i in 2:ncol(Nt)){lines(Nt[,i]~t)}
> plot(apply(Nt,2,max)~NO,type='1')
> plot(Nt[,ncol(Nt)]~t,xlim=range(t),ylim=range(Nt),type='1',)
> plot(Nt[,ncol(Nt)]~t,xlim=range(t),ylim=range(Nt),type='1',)
```

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```
> library('nimble')
>
```

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Alternative state detection? Should we be looking for break points?

```
> source('../src/sens2_ews.R')
              fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second
> x <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second page 1.00 page
> x < -x[,-1]
> x <- x[1:12500,]
> x[x<0] <- 0
> ###
> my.col <- rainbow(nlevels(factor(fps$Prey)))[as.numeric(factor(fps$Prey))]</pre>
> ###
> cx <- x[,5]
> cx[cx <= 0.03] <- 0
> plot(cx)
> abline(v=find.days(cx)$start)
> abline(v=find.days(cx)$end)
> days <- find.days(cx)</pre>
> x.days <- apply(x,2,function(x,d) x[d],d=sort(unlist(days)))</pre>
> x.cor <- list()
> for (k in 1:ncol(x)){
               out <- 0
               for (i in 1:(length(days)-1)){
                       d1 <- days[[i]]</pre>
                       d2 <- days[[(i+1)]]</pre>
+
                       xt \leftarrow x[d1[1:min(c(length(d1), length(d2)))],k]
```

```
xt1 \leftarrow x[d2[1:min(c(length(d1), length(d2)))],k]
      out[i] <- cor(xt,xt1,method='k')</pre>
+
   }
+ x.cor[[k]] <- out
+ }
> plot(x.cor[[1]],type='l',ylim=c(-0.5,1),xlab='Day',col=my.col[[1]])
> for (i in 1:length(x.cor)){lines(x.cor[[i]],col=my.col[[i]])}
> ###
> par(mfcol=c(2,1))
> i <- 1
> plot(density(x[,i]),main=",xlab=expression(0[2]),
       xlim=range(x)+c(-1,1), ylim=c(0,1), col=my.col[i])
> for (i in 2:ncol(x)){
    x.dens <- density(x[,i])</pre>
    x.dens\$y <- x.dens\$y/max(x.dens\$y)
    x.dens$y <- x.dens$y+runif(1,-0.025,0.025)
    lines(x.dens,col=my.col[i])
+ }
> legend('topright',legend=levels(factor(fps$Prey)),lty=1,col=rainbow(nlevels(factor
> ##
> my.col <- rainbow(nlevels(factor(fps$Prey)))[as.numeric(factor(fps$Prey))]</pre>
> i <- 1
> plot(1:nrow(x)~x[,i],col=my.col[i],xlab=expression(0[2]),ylab='time',type='1')
> for (i in 2:ncol(x)){
    lines(1:nrow(x)~x[,i],col=my.col[i],xlab=expression(0[2]),ylab='time',type='1')
+ }
```

```
> 
###Calculate the correlation between days
>
```

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```
- Sub-sample - Standardize, detrend and decycle - Euclidean distance - Plot of eigen
values - PC plot - overlay vectors
> library(vegan)
> library(ecodist)
> source('../src/sens2_ews.R')
    fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second
> x <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second page 1.5.
> x < -x[,-1]
> x < -x[1:12500,]
> x[x<0] <- 0
> ###standardize, detrend and decycle
> sdd.x <- list()
> index <- apply(fps,1,function(x) paste(x[3:6],collapse=' '))</pre>
> cn <- (1:nrow(fps))[fps$Prey==0&index==paste(fps[1,3:6],collapse=' ')]</pre>
> for (i in 1:ncol(x)){
    if (all(x[,i]==0)){sdd.x[[i]] \leftarrow NA}else{
      cn <- (1:nrow(fps))[fps$Prey==0&index==paste(fps[i,3:6],collapse=' ')]</pre>
      sdd.x[[i]] \leftarrow sddSens(x[,i],x[,cn],eval=6250)
```

```
+ }
+ }
> sdd.x <- do.call(cbind,sdd.x)</pre>
> sdd.x[is.na(sdd.x)] <- 0
> ###ordinate
> sdd.d <- dist(t(sdd.x))</pre>
> hist(sdd.d)
> plot(hclust(sdd.d))
> ##
> par(mfrow=c(1,3))
> plot(sdd.x[,1],xlab='time',ylab=expression(0[2]),ylim=c(-1,1),pch=19,cex=0.05)
> for (i in 1:ncol(sdd.x)){points(sdd.x[,i],pch=19,cex=0.05)}
> sdd.ord <- princomp(sdd.d)</pre>
> names(sdd.ord)
> plot(sdd.ord$sdev/sum(sdd.ord$sdev))
> abline(h=0.05,lty=2)
> plot(sdd.ord$scores[,1:2])
> plot(envfit(ord=sdd.ord$scores[,1:2],env=fps))
> ###
> envfit(ord=sdd.ord$scores[,1:2],env=fps)
> envfit(ord=sdd.ord$scores[,2:3],env=fps)
> envfit(ord=sdd.ord$scores[,1:3],env=fps)
> pairs(fps)
> pairs(sdd.ord$scores[,1:3],pch=19,cex=1)
> par(mfrow=c(3,3))
> for (i in 1:3){
```

```
for (j in 1:3){
      if (i==j){
        plot(c(-1,1),c(-1,1),pch="',xlab="',ylab="',xaxt='n',yaxt='n')
        text(c(0,0),c(0,0),labels=i,cex=5)
      }else{
        plot(sdd.ord$scores[,j:i],pch=19,cex=0.75)
        plot(envfit(ord=sdd.ord$scores[,j:i],env=fps))
   }
   }
+ }
> library(rgl)
> plot3d(sdd.ord$scores[,1:3],type='s',col=rainbow(nlevels(factor((fps$Prey))))[as.n
> plot3d(sdd.ord$scores[,1:3],type='s',col=rainbow(nlevels(factor((fps$b))))[as.nume
> plot3d(sdd.ord$scores[,1:3],type='s',col=rainbow(nlevels(factor((fps$a))))[as.nume
> plot3d(sdd.ord$scores[,1:3],type='s',col=rainbow(nlevels(factor((fps$K))))[as.nume
> plot3d(sdd.ord$scores[,1:3],type='s',col=rainbow(nlevels(factor((fps$d))))[as.nume
>
```

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Plot tests (ordination and smoothed-detrended-decycled) are up on github.

EWS for simulations:

- Write a script that will record the early warning signals for all 180 simulations.

```
> source('../src/sens2_ews.R')
```

> fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and secon

```
> x <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second page 1
> x < -x[,-1]
> #x <- x[1000:5400,]
> x[x<0] <- 0
> ###Isolate controls
> my.col <- rainbow(nlevels(factor(fps$Prey)))[as.numeric(factor(fps$Prey))]</pre>
> sdd.x <- list()
> index <- apply(fps,1,function(x) paste(x[3:6],collapse=' '))</pre>
> cn <- (1:nrow(fps))[fps$Prey==0&index==paste(fps[1,3:6],collapse=' ')]</pre>
> plot(sdd.x[[1]] \leftarrow sddSens(x[,1],x[,cn]),ylim=c(-1,1),pch=19,
       cex=0.25,ylab=expression(0[2]),xlab='t',col=my.col[1])
> for (i in 1:ncol(x)){
    if (all(x[,i]==0)){sdd.x[[i]] \leftarrow NA}else{
      cn <- (1:nrow(fps))[fps$Prey==0&index==paste(fps[i,3:6],collapse=' ')]</pre>
      points(sdd.x[[i]] \leftarrow sddSens(x[,i],x[,cn]),pch=19,cex=0.25,col=my.col[i])
+
    }
+ }
> legend('topright',legend=unique(fps$Prey),col=rainbow(nlevels(factor(fps$Prey))),p
> ###
> par(mfrow=c(1,2))
> plot(x[,1])
> plot(sdd.x[[1]])
> ###
> my.col <- rainbow(nlevels(factor(fps$b)))[as.numeric(factor(fps$b))]
> sdd.x <- list()
> index <- apply(fps,1,function(x) paste(x[3:6],collapse=' '))</pre>
```

```
> cn <- (1:nrow(fps))[fps$Prey==0&index==paste(fps[1,3:6],collapse=' ')]</pre>
> plot(sdd.x[[1]] \leftarrow sddSens(x[,1],x[,cn]),ylim=c(-1,1),pch=19,
       cex=0.25, ylab=expression(0[2]), xlab='t', col=my.col[1])
> for (i in 1:ncol(x)){
    if (all(x[,i]==0)){sdd.x[[i]] \leftarrow NA}else{
      cn <- (1:nrow(fps))[fps$Prey==0&index==paste(fps[i,3:6],collapse=' ')]</pre>
      points(sdd.x[[i]] \leftarrow sddSens(x[,i],x[,cn]),pch=19,cex=0.25,col=my.col[i])
   }
+
+ }
> legend('topright',legend=levels(factor(fps$b)),col=rainbow(nlevels(factor(fps$b)))
> library(strucchange)
>
>
  Look at the temporal correlation structure among simulations:
> ##Using the 100 min averages
> ##Runs are in columns
    fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second
> x <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second page 1
> x < -x[,-1]
> plot(x[,1],xlab='t',ylab=expression(0[2]),ylim=c(0,max(x)),type='1',lwd=0.5,col=rai
> for (i in 2:ncol(x)){
    lines(1:nrow(x),x[,i],lwd=0.5,col=rainbow(ncol(x))[i])
+ }
> abline(v=c(1000,5400))
> x <- x[1000:5400,]
```

```
> x[x<0] <- 0
> d.x <- dist(t(x))
> library(vegan)
> library(ecodist)
> eig.x <- eigen(d.x)
> biplot(princomp(d.x))
> vd.x <- vegdist(t(rbind(x,rep(1,ncol(x)))))
> nmds.x <- nmds(vd.x,2,2)
> envf.x <- envfit(env=fps,ord=nmds.min(nmds.x))
> plot(apply(nmds.min(nmds.x),2,jitter,factor=50),col=rainbow(ncol(x)))
> plot(envf.x)
> envf.x
```

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Generate plots for the tipping points 2

- Model description is in SarrModel-20140709-AME.lyx
- Note that the a and b terms for calculating w(t) are coupled in order to keep the integral (i.e. the area under the decomposition curve) equal across simulations
- $\bullet \ \ \text{Simulation output is in } \ \ \text{Raw}_runs.csv, ten_min_average.csv, hundred_min_average.csvPAR to O2.docx descriptions and the property of the prope$

• Model was coded in mathematica (Pitcher $Plant_Threshold_Model.txt$) Free parameter space for sim

```
##free parameter space
                   fps <- read.csv('~/Dropbox/Tipping Point MS/Model sensitivity analysis and second
> ##Using the 100 min averages
> ##Runs are in columns
> x <- read.csv('/Users/Aeolus/Dropbox/Tipping Point MS/Model sensitivity analysis a
> x < -x[,-1]
> plot(x[,1],pch=19,cex=0.1)
> for (i in 2:ncol(x)){
            points(1:length(x[,i]),x[,i],pch=19,cex=0.1)
+ }
> par(mfrow=c(1,2))
> i <- 2
> plot(x[,i],pch=19,cex=0.1,col=rainbow(length(x[,i])))
> plot(x[2:length(x[,i]),i]^x[1:(length(x[,i])-1),i],pch=19,cex=0.1,col=rainbow(length(x[,i]),i]^x[1:(length(x[,i])-1),i],pch=19,cex=0.1,col=rainbow(length(x[,i]),i]^x[1:(length(x[,i])-1),i],pch=19,cex=0.1,col=rainbow(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(l
> par(mfrow=c(1,2))
> for (i in 1:ncol(x)){
                    plot(x[,i],pch=19,cex=0.1,col=rainbow(length(x[,i])),
                                               main=paste(paste(colnames(fps)[2:ncol(fps)],fps[i,2:ncol(fps)],sep='='),col
+
                    plot(x[2:length(x[,i]),i]^x[1:(length(x[,i])-1),i],pch=19,cex=0.1,col=rainbow(length(x[,i]),i]^x[1:(length(x[,i])-1),i],pch=19,cex=0.1,col=rainbow(length(x[,i]),i]^x[1:(length(x[,i])-1),i],pch=19,cex=0.1,col=rainbow(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(length(x[,i])-1),i]^x[1:(len
                                                main=paste(paste(colnames(fps)[2:ncol(fps)],fps[i,2:ncol(fps)],sep='='),col
+
                    play(x[,i]^3/max(x^3))
                    Sys.sleep(0.15)
+ }
```

The pitcher plant respiration model:

Terms:

- t = time
- x = oxygen
- A = environmental factor promoting oxygenation (i.e. PAR = light)
- f(w,x) = loss/decay of oxygen
- \bullet g(x) = recovery of oxygen (augmented by mineralized nutrients)
- w = prey mass

$$\frac{dx}{dt} = A - f(w, x) + g(x) \tag{1}$$

$$x_{t+1} = a_t * \sin(2\pi f t) - (m + a_t \frac{w_{t-1}}{K_w + w_{t-1}}) + D_t(x_t, a_t')$$
(2)

>

$$w(t+1) = ae^{-bw_0t} (3)$$

Note that this is a correction from the printed version in PNAS

```
> y <- NULL
> a <- 20
> ##b
> for (i in 1:17) { y[i] <- 20*exp(-.1395*(i-1)) }
> plot(c(0:16), y, xlim=c(0,16), ylim=c(0,20))
> par(list(pin=c(3,3),las=1))
> integrand \leftarrow function(x, a=4) {20*exp(-a*x)}
> plot(integrand, xlim=c(0,16), xlab="Days", ylab="Mass of prey remaining (mg)",
                        lwd=2, col="red", font=2, font.lab=2)
> integrate(integrand, lower=0, upper=16)
> ###
> a <- 20
> b <- 4
> e <- exp(1)
> w <- 75/1000
> for (t in 2:20){
+ w[t] <- exp(-w[t-1])
+ }
> plot(w,type='1')
```

Here is a link to the pitcher-plant tipping point dropbox. To start, go to ../Model sensitivity analysis and second paper/data/ and read the metadata.csv file. The various raw and average files are the ones to think about plotting. We can talk about it later today.

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Question: Are transitions between chaotic states detectable?

Question: Are there classes of models that possess detectable chaotic transition warning signals?

Modeling Issues:

- 1. a in S2.2 can be confused with a in S2.1
- 2. What is the exponent in S2.2?
- 3. Why is b (S2.2) in days rather than minutes?
- 4. Why is w(t-1) used in S2.3?

```
> ##Initial stab at the spo2 model
> a0 <- 10
> f <- 1/(60*24) #total minutes in a day
> t <- rep((1:(60*24)),11)
> At <- a0 * sin(2*pi*f*t)
> At[t>720] <- 0
> plot(At~I(1:length(t)),type='l')
> a <- 20
> b <- 4
> wt <- 20
> wt <- a*exp(1)^(-b*wt)
> spo2 <- function(x,a,f,t,m,w,Kw,s,d){
+ At <- a0 * sin(2*pi*f*t)</pre>
```

+

+ }

>

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Chatted with Aaron

The data don't support a general pattern of alternative states.

If systems are inherently chaotic or stochastic, then what?

The data:

- Climate isn't stable
- Fossil records don't show stability
- Tropical systems don't show stability
- Even temperate systems break the rule of stability

Two goals:

- 1. Can we tell when we transition between states?
- 2. Is there an alternative philosphy? Mathematical framework?

Run simulation increasing r with variance over the chaos threshold

- > ##hold ni constant
- > ##hold sd constant at 0.1
- > ##record r
- > ##record n

```
##record ews
    source('../src/unstable_states.R')
> library(earlywarnings)
> cb8.16 <- 2.57 #choatic boundary, 8-16 cycles
> ri <- (cb8.16)-(0.02/2)
> rf <- (cb8.16)+(0.02/2)
> dmc <- list()</pre>
> ews <- list()
> mcn <- 100
> for (i in 1:mcn){
   print(i)
   dmc[[i]] <- disrupt.mc(N=10,sd=0.01,ri=ri,rf=rf,dump=TRUE)</pre>
    ews[[i]] <- generic_ews(dmc[[i]]$N)</pre>
    dev.off()
+ }
> ## dput(dmc,'../results/dmc.out')
> ## dput(ews,'../results/ews.out')
> ###
> ###
> ###
> dmc <- dget('../results/dmc.out')</pre>
> ews <- dget('../results/ews.out')</pre>
> ##
> yl <- apply(do.call(rbind,ews),2,min)</pre>
> yu <- apply(do.call(rbind,ews),2,max)</pre>
> ##
```

```
> par(mfrow=c(2,(ncol(ews[[1]])-1)/2))
> for (i in 2:ncol(ews[[1]])){
   for (j in 1:length(ews)){
      if (j==1){
       plot(ews[[j]][,i]~ews[[j]][,1],ylim=c(yl[i],yu[i]),
             ylab=colnames(ews[[1]])[i],xlab='t',
             type='1',1wd=0.25)
      }else{
        lines(ews[[j]][,i]~ews[[j]][,1],lwd=0.25)
      }
+
+
   }
+ }
> ###Average plots
> ews. <- do.call(rbind,ews)
> t <- do.call(rbind,dmc)$t
> r <- do.call(rbind,dmc)$r
> r. <- r
> r[r.>=cb8.16] <- 2
> r[r.<cb8.16] <- 1
> r \leftarrow r[t>=ews[[1]][,1][1]]
> #pairs(ews.,cex=0.05,pch=19,col=r)
> par(mfrow=c(1,1))
> plot(ews.[,c(3,4)],pch=19,col=r,cex=(0.01+(0.5*(ews.[,1]/max(ews.[,1])))))
> unique(ews.[ews.[,3]>30.15&ews.[,4]>0.185,1])
> ews. <- apply(ews.,2,function(x,t) tapply(x,t,mean),t=ews.[,1])
> pairs(ews.,cex=0.10,pch=19)
```

>

Reading Hastings Reading Sheffer Reading Dakos

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- the distribution of the average of r is uniform
- ensemble distribution is normal
- EWS stats not correlated between 8to16 and 16to8
- Phase (Ni to Nf) spaces for EWS correlations
- EWS stats intercorrelations show correlations and break points

```
> ###Run repeated simulations for ews time series
```

```
> source('../src/unstable_states.R')
```

- > library(earlywarnings)
- > cb8.16 <- 2.57 #choatic boundary, 8-16 cycles
- > ri <- (cb8.16)-(0.02/2)
- > rf <- (cb8.16)+(0.02/2)
- > ##Visualizing the error in r
- > r8.16 <- list()
- > r16.8 <- list()
- > for (i in 1:138){
- + print(i)
- + r8.16[[i]] <- disrupt.mc(N=i,sd=0.01,ri=ri,rf=rf,dump=TRUE)\$r
- + r16.8[[i]] <- disrupt.mc(N=i,sd=0.01,ri=rf,rf=ri,dump=TRUE)\$r

```
+ }
> rmu8.16 <- apply(do.call(rbind,r8.16),2,mean)</pre>
> rmu16.8 <- apply(do.call(rbind,r16.8),2,mean)</pre>
> par(mfrow=c(1,2))
> plot(density(rmu8.16),main='',xlab='r')
> for (i in 1:length(r8.16)){
    lines(density(r8.16[[i]]),col='grey',lwd=0.5)
+ }
> abline(v=cb8.16,lty=2);abline(v=mean(rmu8.16),lty=2,col='darkgrey')
> plot(density(rmu16.8),xlab='r',main='')
> for (i in 1:length(r16.8)){
    lines(density(r16.8[[i]]),col='grey',lwd=0.5)
+ }
> abline(v=cb8.16,lty=2);abline(v=mean(rmu16.8),lty=2,col='darkgrey')
> ###Determine average threshold
> ##What is the point at which rbar has crossed the threshold?
> ##Directionality depends on direction of r
> rt8.16 <- (1:length(rmu8.16))[rmu8.16>=cb8.16][1]
> rt16.8 <- (1:length(rmu16.8))[rmu16.8<=cb8.16][1]</pre>
> rt8.16
> rt16.8
> ###EWS stats
> stats8.16 <- dget(file='../results/stats816.rdata')</pre>
> stats16.8 <- dget(file='../results/stats168.rdata')</pre>
> ###
> stats8.16 <- na.omit(do.call(rbind,stats8.16))</pre>
```

```
> stats16.8 <- na.omit(do.call(rbind,stats16.8))</pre>
> ###
> stats8.16 <- stats8.16[1:min(c(nrow(stats8.16),nrow(stats16.8))),]
> stats16.8 <- stats16.8[1:min(c(nrow(stats8.16),nrow(stats16.8))),]</pre>
> ###
> par(mfrow=c(2,ncol(stats8.16)/2),
      mai=c(0.25,0.01,0.25,0.01))
> for (i in 1:ncol(stats8.16)){
    plot(density(stats8.16[,i]),
         main=colnames(stats8.16)[i],
         xlim=c(min(c(stats8.16[,i],stats16.8[,i])),
+
           max(c(stats8.16[,i],stats16.8[,i]))),
         xaxt='n', yaxt='n', bty='n')
    lines(density(stats16.8[,i]),lty=2)
+ }
> ###
> par(mfrow=c(2,ncol(stats16.8)/2))
> for (i in 1:ncol(stats8.16)){
    plot(stats16.8[,i]~stats8.16[,i],
         xlab=paste(colnames(stats8.16)[i],'8.16'),
         ylab=paste(colnames(stats16.8)[i],'16.8'))
    abline(lm(stats16.8[,i]~stats8.16[,i]))
+ }
> ###
> par(mfrow=c(4,ncol(stats8.16)/2),
      mai=c(0,0,0,0)
```

```
> for (i in 1:ncol(stats8.16)){
   plot(stats8.16[1:(nrow(stats8.16)-1),i]~stats8.16[2:(nrow(stats8.16)),i],
         type='1', xaxt='n', yaxt='n', bty='n')
+ }
> for (i in 1:ncol(stats16.8)){
   plot(stats16.8[1:(nrow(stats16.8)-1),i]~stats16.8[2:(nrow(stats16.8)),i],
         type='1',col='darkgrey',xaxt='n',yaxt='n',bty='n')
+ }
> ###Ensemble N~stats
> pairs(data.frame(Ni=(1:nrow(stats8.16)),stats8.16),pch=19,cex=0.10,col='black')
> pairs(data.frame(Ni=(1:nrow(stats16.8)),stats16.8),cex=0.10,col='black')
>
> library(earlywarnings)
> set.seed(1)
> drmc8.16 <- disrupt.mc(sd=0.01,ri=ri,rf=rf,dump=TRUE)</pre>
> set.seed(1)
> drmc16.8 <- disrupt.mc(sd=0.01,ri=rf,rf=ri,dump=TRUE)</pre>
> ews8.16 <- generic_ews(drmc8.16$N)</pre>
> ews16.8 <- generic_ews(drmc16.8$N)</pre>
> stats8.16 <- cor(ews8.16,method='ken')
> stats16.8 <- cor(ews8.16,method='ken')</pre>
>
    ##Re-doing noise in r shifting up and down across 8-16
> source('../src/unstable_states.R')
```

```
> cb8.16 <- 2.57 #choatic boundary, 8-16 cycles
> ri <- (cb8.16)-(0.02/2)
> rf <- (cb8.16)+(0.02/2)
> set.seed(1)
> drmc8.16 <- disrupt.mc(sd=0.01,ri=ri,rf=rf,dump=TRUE)</pre>
> set.seed(1)
> drmc16.8 <- disrupt.mc(sd=0.01,ri=rf,rf=ri,dump=TRUE)</pre>
> par(mfrow=c(3,2))
> plot(drmc8.16$r,xlab='time',ylab='r')
> abline(h=cb8.16,col=2)
> plot(drmc16.8$r,xlab='time',ylab='r')
> abline(h=cb8.16,col=2)
> plot(drmc16.8$N,xlab='time',ylab='N')
> plot(drmc8.16$N,xlab='time',ylab='N')
> plot(drmc16.8$N~drmc16.8$r,xlab='r',ylab='N',type='l')
> plot(drmc8.16\$N^drmc8.16\$r,xlab='r',ylab='N',type='l')
> ###Phase space
> par(mfrow=c(1,2))
>
                                             #plus n time steps
> for (n in 25:50){
   plot(drmc16.8$N[(1:(length(drmc16.8$N)-n))],drmc16.8$N[((n+1):(length(drmc16.8$
         xlab='N', ylab='N+1', type='l'
    plot(drmc8.16\$N[(1:(length(drmc8.16\$N)-n))],drmc8.16\$N[((n+1):(length(drmc8.16\$N)-n))]
         xlab='N', ylab='N+1', type='l')
+ }
> #EWS
```

```
> library(earlywarnings)
> ews8.16 <- generic_ews(drmc8.16$N)
> ews16.8 <- generic_ews(drmc16.8$N)
>
```

Summary to date (going back in time):

- At the 8-16 cycle threshold, error in r leads to early, sudden shifts
- Slow ramping can also be seen visually
- Sudden jumps across cycle boundaries can be seen visually