### Time Series Analysis

Week 4 - Introduction to Stochastic Processes, Operators and Linear Systems

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## Week 4: Outline of the lecture

- Operators; the backward shift operator; sec. 4.5.
- Stochastic processes in general: Sec 5.1, 5.2, 5.3 [except 5.3.2].

## Operators; The backwards shift operator B

- An operator A is (here) a function of a time series  $\{x_t\}$  (or a stochastic process  $\{X_t\}$ ).
- ▶ Application of an operator on a time series {x<sub>t</sub>} yields a new time series {Ax<sub>t</sub>}. Likewise of a stochastic process {AX<sub>t</sub>}.

# Operators; The backwards shift operator B

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- ▶ Application of an operator on a time series {x<sub>t</sub>} yields a new time series {Ax<sub>t</sub>}. Likewise of a stochastic process {AX<sub>t</sub>}.
- ► Most important operator for us: The backwards shift operator  $B : Bx_t = x_{t-1}$ . Notice  $B^j x_t = x_{t-j}$ .
- ▶ All other operators we shall consider in this lecture may be expressed in terms of *B*.

### Question

					t+1		
$x_t$	0.2	0.1	1.3	0.8	0.4	0.1	0.5
$B^j x_t$			0.2	0.1	1.3	0.8	0.4

What is j?

### Question

	t-3	t-2	t-1	t	t+1	t+2	t+3
$x_t$	0.2	0.1	1.3	0.8	0.4	0.1	0.5
$B^j x_t$	0.1	1.3	0.8	0.4	0.1	0.5	

What is *j*?

# The forward shift ${\cal F}$

The forward shift operator

► 
$$Fx_t = x_{t+1}$$
;  $F^j x_t = x_{t+j}$ ;

# The forward shift F

#### The forward shift operator

- F $x_t = x_{t+1}$ ;  $F^j x_t = x_{t+j}$ ; How can F be expressed in terms of B? ( $Bx_t = x_{t-1}$ )
- Combining a forward and backward shift yields the identity operator;  $BFx_t = Bx_{t+1} = x_t$ , ie. F and B are each others inverse:  $B^{-1} = F$  and  $F^{-1} = B$ .

Question					1	I .			
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# The difference operator $\nabla$

#### The difference operator

 $\triangleright \nabla x_t = x_t - x_{t-1}$ 

• How can  $\nabla$  be expressed with B? A:  $\nabla = 1 - B$  or B:  $\nabla = 1 + B^{-1}$ 

# The difference operator $\nabla$

#### The difference operator

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• How can  $\nabla$  be expressed with B? A:  $\nabla = 1 - B$  or B:  $\nabla = 1 + B^{-1}$ 

 $= (1-B)x_t.$ 

▶ Thus:  $\nabla = 1 - B$ .

# The summation S

$$Sx_t = x_t + x_{t-1} + x_{t-2} + \dots$$
  
=  $x_t + Bx_t + B^2 x_t \dots$   
=  $(1 + B + B^2 + \dots) x_t$ 

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Summation, then difference (using 
$$Sx_t = x_t + Sx_{t-1}$$
)  
 $\nabla Sx_t = Sx_t - Sx_{t-1} = x_t + Sx_{t-1} - Sx_{t-1} = x_t$ 

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Difference, then summation

$$S\nabla x_t = (1 + B + B^2 \dots) x_t - (1 + B + B^2 \dots) x_{t-1}$$
  
= (1 + B + B^2 \dots) x\_t - (B + B^2 \dots) x\_t = x\_t

• So  $\nabla$  and S are each others inverse:

$$\nabla^{-1} = \frac{1}{1-B} = 1 + B + B^2 + \ldots = S$$

#### In R:

x <- 1:10
x
diff(x)
cumsum(x)
# Not so nice in the ends!
cumsum(diff(x))
diff(cumsum(x))
# Put a zero in beginning to fix it
cumsum(diff(c(0,x)))
diff(cumsum(c(0,x)))</pre>

Shift operator is simply the most important:

```
lagvec <- function(x, lag){
    if (lag > 0) {
        ## Lag x, i.e. delay x lag steps
        return(c(rep(NA, lag), x[1:(length(x) - lag)]))
    }else if(lag < 0) {
        ## Lag x, i.e. delay x lag steps
        return(c(x[(abs(lag) + 1):length(x)], rep(NA, abs(lag))))
    }else{
        ## lag = 0, return x
        return(x)
    }
}</pre>
```

#### Lag to a data frame:

```
lagdf <- function(x, lag) {
    # Lag x, i.e. delay x lag steps
    if (lag > 0) {
        x[(lag+1):nrow(x), ] <- x[1:(nrow(x)-lag), ]
        x[1:lag, ] <- NA
    }else if(lag < 0) {
        # Lag x "ahead in time"
        x[1:(nrow(x)-abs(lag)), ] <- x[(abs(lag)+1):nrow(x), ]
        x[(nrow(x)-abs(lag)+1):nrow(x), ] <- NA
    }
    return(x)
}</pre>
```

## Properties of B, F, $\nabla$ and S

► The operators are all linear, ie.

$$H[ax_t + by_t] = aH[x_t] + bH[y_t]$$

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$$H[ax_t + by_t] = aH[x_t] + bH[y_t]$$

The operators may be combined into new operators: The power series

$$a(z) = \sum_{i=0}^{\infty} a_i z^i$$

defines a new operator from an operator H by linear combinations:

$$a(H) = \sum_{i=0}^{\infty} a_i H^i$$

# Examples of combined operators

 $\blacktriangleright$   $\nabla^{-1}$ :

$$\frac{1}{1-z} = \sum_{i=0}^{\infty} z^i$$
 so  $\nabla^{-1} = \frac{1}{1-B} = \sum_{i=0}^{\infty} B^i = S$ 

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Operator polynomial of order q:

$$heta(z) = \sum_{i=0}^q heta_i z^i$$

ie.  $\theta_i = 0$  for i > q.

$$\theta(B) = (1 + \theta_1 B + \dots + \theta_q B^q)$$

where  $\theta_0$  is chosen to be 1

# Stochastic Processes - in general

- Function:  $X(t, \omega)$
- ▶ Time:  $t \in T$
- Realization:  $\omega \in \Omega$
- lndex set: T
- Sample Space: Ω
- $X(\cdot, \cdot)$  is a stochastic process
- $X(t, \cdot)$  is a random variable

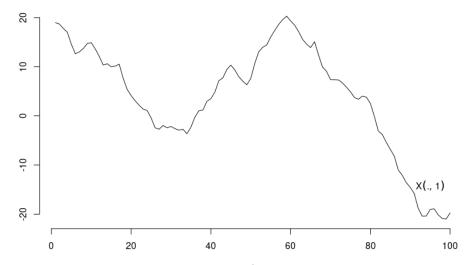
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- $X(t, \omega)$  is an observation. This is what we often denote  $x_t$ .

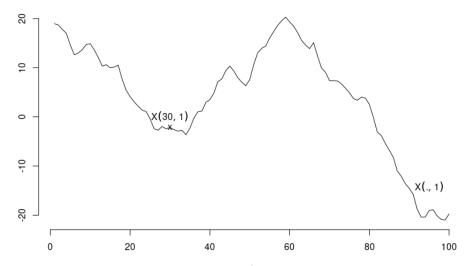
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- $X(t, \omega)$  is an observation. This is what we often denote  $x_t$ .
- In this course we restrict ourselves to the case where time is discrete, and the realizations take values on the real numbers (continuous range).

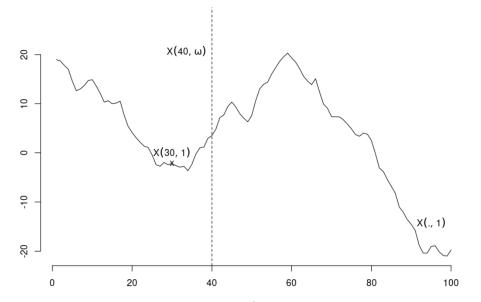
## Stochastic Processes - illustration



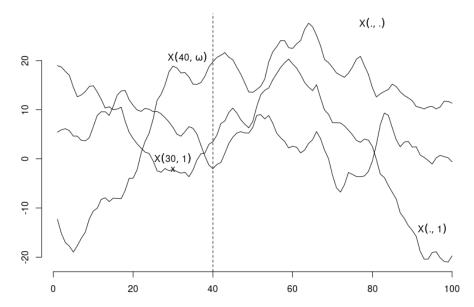
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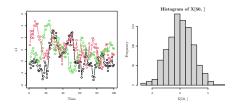
## Stochastic Processes - illustration



#### Demo i R:

```
par(mfrow=c(1,2))
# Generate realizations
n <- 100
x1 <- filter(rnorm(n), 0.9, "recursive")
plot(x1, type="b", ylim=c(-6,6))
x2 <- filter(rnorm(n), 0.9, "recursive")
lines(x2, type="b", col=2)
x3 <- filter(rnorm(n), 0.9, "recursive")
lines(x3, type="b", col=3)</pre>
```

```
# Generate 1000 realizations
X <- matrix(filter(rnorm(1000*n), 0.9, "recursive"), nrow=n)
# One realization (i.e. a time series)
X[ ,1]
# Realization of the stochastic variable at one time point
hist(X[50, ])</pre>
```



# Complete Characterization

In the end a stochastic process is just a multivariate variable.

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In the end a stochastic process is just a multivariate variable. Thus, from lecture 1, it is characterised by its n-dimensional probability density:

 $f_{X(t_1),\ldots,X(t_n)}(x_1,\ldots,x_n)$ 

# 2nd order moment representation

Mean function:

$$\mu(t) = E[X(t)] = \int_{-\infty}^{\infty} x f_{X(t)}(x) dx,$$

Autocovariance function:

$$\begin{aligned} \gamma_{XX}(t_1, t_2) &= \gamma(t_1, t_2) = \operatorname{Cov} \left[ X(t_1), X(t_2) \right] \\ &= E \left[ (X(t_1) - \mu(t_1)) (X(t_2) - \mu(t_2)) \right] \end{aligned}$$

The variance function is obtained from  $\gamma(t_1, t_2)$  when  $t_1 = t_2 = t$ :

$$\sigma^{2}(t) = V[X(t)] = E\left[(X(t) - \mu(t))^{2}\right]$$

# Stationarity

▶ A process {*X*(*t*)} is said to be *strongly stationary* if all finite-dimensional distributions are invariant for changes in time, i.e. for every *n*, and for any set (*t*<sub>1</sub>, *t*<sub>2</sub>, ..., *t*<sub>n</sub>) and for any *h* it holds

$$f_{X(t_1),\dots,X(t_n)}(x_1,\dots,x_n) = f_{X(t_1+h),\dots,X(t_n+h)}(x_1,\dots,x_n)$$

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$$f_{X(t_1),\dots,X(t_n)}(x_1,\dots,x_n) = f_{X(t_1+h),\dots,X(t_n+h)}(x_1,\dots,x_n)$$

- ► A process {X(t)} is said to be *weakly stationary of order* k if all the first k moments are invariant to changes in time
- A weakly stationary process of order 2 is simply called *weakly stationary* or just *stationary*:

$$\mu(t) = \mu$$
  $\sigma^2(t) = \sigma^2$   $\gamma(t_1, t_2) = \gamma(t_1 - t_2)$ 

# Ergodicity

- ▶ In time series analysis we normally assume that we have access to one realization only
- We therefore need to be able to determine characteristics of the process  $X_t$  from one realization  $x_t$

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- ▶ In time series analysis we normally assume that we have access to one realization only
- We therefore need to be able to determine characteristics of the process  $X_t$  from one realization  $x_t$
- It is often enough to require the process to be mean-ergodic:

$$E[X(t)] = \int_{\Omega} x(t,\omega) f(\omega) \, d\omega = \lim_{T \to \infty} \frac{1}{2T} \int_{-T}^{T} x(t,\omega) \, dt$$

i.e. if the mean of the ensemble equals the mean over time

Some intuitive examples, not directly related to time series: http://news.softpedia.com/news/What-is-ergodicity-15686.shtml

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- Markov processes: The conditional distribution depends only on the latest state of the process:

 $P\{X(t_n) \le x | X(t_{n-1}), \cdots, X(t_1)\} = P\{X(t_n) \le x | X(t_{n-1})\}$ 

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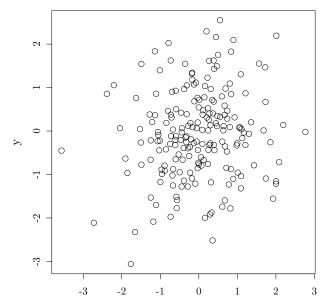
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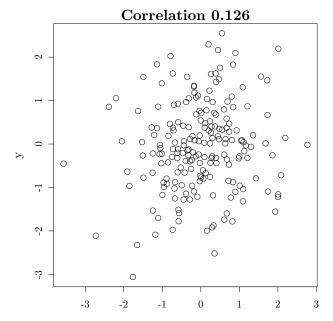
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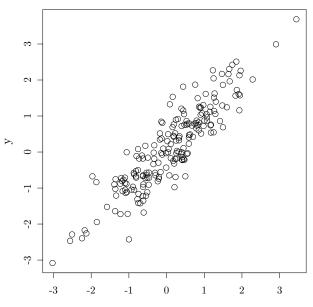
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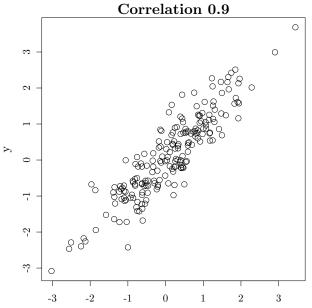
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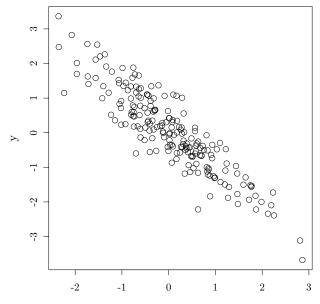
- Deterministic processes: Can be predicted without uncertainty from past observations
- Pure stochastic processes: Can be written as a linear combination of uncorrelated random variables
- Decomposition:  $X_t = S_t + D_t$ , where  $S_t$  is a pure stochastic process and  $D_t$  is a deterministic process.

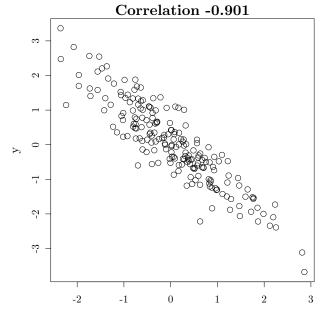


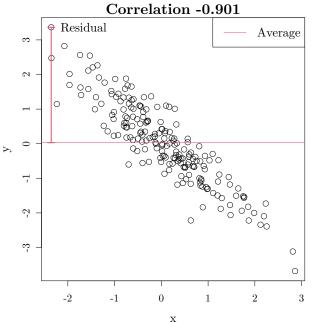






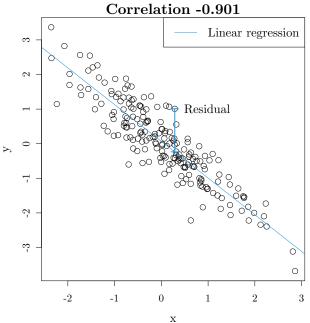






Summed Squared Residuals

$$SS_{average} = \sum_{i}^{n} Residual_{i}^{2}$$



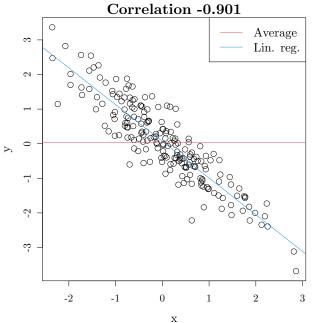
#### Summed Squared Residuals

$$SS_{\mathsf{lin}} = \sum_{i}^{n} Residual_{i}^{2}$$

#### How to calculate correlation?

Two vectors with corresponding elements  $x_i$  and  $y_i$ .

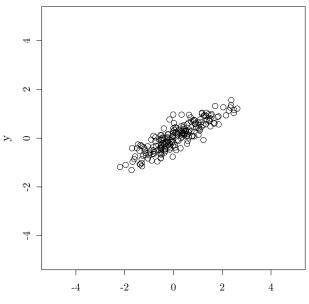
$$R^2 = \frac{SS_{\text{average}} - SS_{\text{lin}}}{SS_{\text{average}}}$$

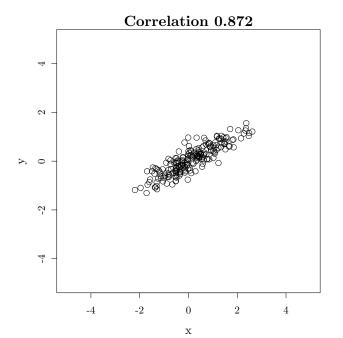


$$R^{2} = \frac{SS_{\text{average}} - SS_{\text{lin}}}{SS_{\text{average}}}$$
$$= \frac{270.548 - 50.81256}{270.548}$$
$$= 0.8121865$$

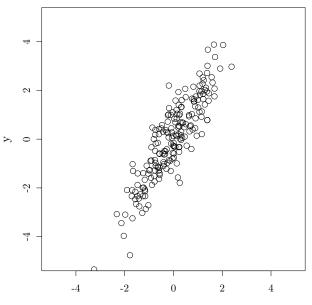
and

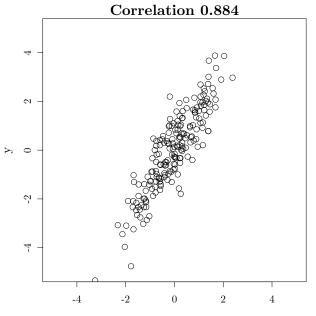
 $Cor = sign(slope) \cdot R$ = -0.9012139

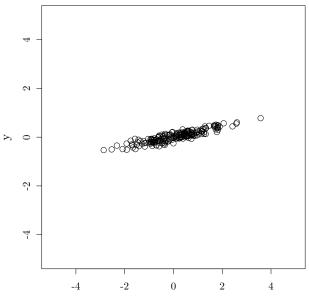


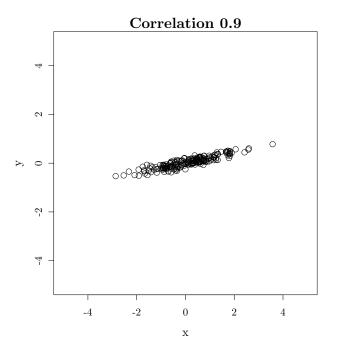


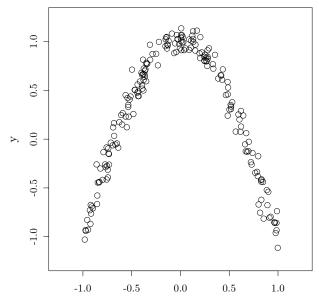
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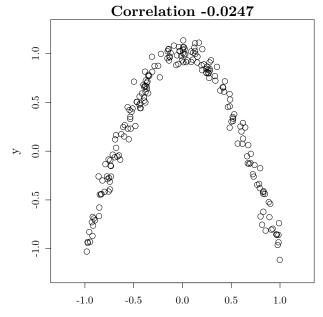












5.2.2 Covariance and correlation functions

### Autocovariance and autocorrelation

▶ For stationary processes: Only dependent on the time difference  $au = t_2 - t_1$ 

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Autocovariance:

$$\gamma(\tau) = \gamma_{XX}(\tau) = \operatorname{Cov}[X(t), X(t+\tau)] = E[X(t)X(t+\tau)]$$

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$$ho( au) = 
ho_{XX}( au) = \gamma_{XX}( au)/\gamma_{XX}(0) = \gamma_{XX}( au)/\sigma_X^2$$

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$$\rho(\tau) = \rho_{XX}(\tau) = \gamma_{XX}(\tau)/\gamma_{XX}(0) = \gamma_{XX}(\tau)/\sigma_X^2$$

Some properties of the autocovariance function:

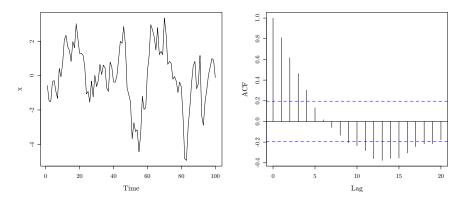
$$\blacktriangleright \ \gamma(\tau) = \gamma(-\tau)$$

 $\blacktriangleright |\boldsymbol{\gamma}(\tau)| \leq \boldsymbol{\gamma}(0)$ 

# ACF

```
# Simulate a process
n <- 100
x <- filter(rnorm(n), 0.9, "recursive")
plot(x)</pre>
```

```
# The acf and "height" of "lag 1 bar"
val <- acf(x)
val[1]
# The correlation with lag 1
cor(x, lagvec(x,1), use="complete.obs")</pre>
```



#### White noise

• Def. 5.9:  $\{\varepsilon_t\}$  is a sequence of mutually uncorrelated identically distributed random variables, where:

$$\blacktriangleright \ \mu_t = \mathsf{E}[\varepsilon_t] = 0$$

• 
$$\sigma_t = \operatorname{Var}[\varepsilon_t] = \sigma_{\varepsilon}^2$$

• 
$$\gamma_{\varepsilon}(k) = \operatorname{Cov}[\varepsilon_t, \varepsilon_{t+k}] = 0$$
, for  $k \neq 0$ 

How many of the ACF bars stick of the CI if white noise?

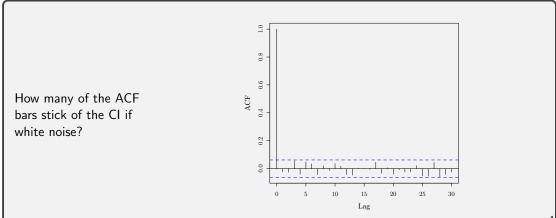
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, for  $k \neq 0$ 



A small game! Go try it now: Can you make white noise? Can you make an alternating pattern? Can you make an alternating pattern with "longer period"? Can you make a slow decay to zero?

```
x <- readline("Write a sequence of numbers from 0 to 9\n")
x <- as.numeric(strsplit(x,"")[[1]])
acf(x)</pre>
```

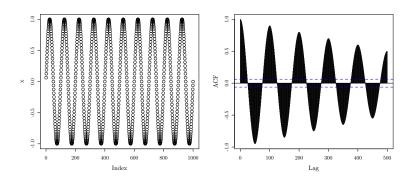
### More ACF...

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# Plot it
par(mfrow=c(1,2))
plot(x)
# The ACF is actually also a sine!
acf(x, lag.max=500)

# Can you "screw it" and make it look like white noise up by altering a single value in x?

#x[100] <- ?? #acf(x, lag.max=500)



#### Linear processes

• A linear process  $\{Y_t\}$  is a process that can be written on the form

$$Y_t - \mu = \sum_{i=0}^\infty \psi_i arepsilon_{t-i}$$

where  $\mu$  is the mean value of the process and

- $\triangleright$  { $\varepsilon_t$ } is white noise, i.e. a sequence of uncorrelated, identically distributed random variables.
- $\{\varepsilon_t\}$  can be scaled so that  $\psi_0 = 1$
- Without loss of generality we assume  $\mu = 0$

### $\psi$ - and $\pi$ -weights

Transfer function and linear process:

$$\psi(B) = 1 + \sum_{i=1}^{\infty} \psi_i B^i$$
  $Y_t = \psi(B)\varepsilon_t$ 

Inverse operator (if it exists) and the linear process:

$$\pi(B) = 1 + \sum_{i=1}^{\infty} \pi_i B^i \qquad \pi(B) Y_t = \varepsilon_t,$$

• Autocovariance using  $\psi$ -weights:

$$\gamma(k) = \operatorname{Cov}\left[Y_t, Y_{t+k}\right] = \operatorname{Cov}\left[\sum_{i=0}^{\infty} \psi_i \varepsilon_{t-i}, \sum_{i=0}^{\infty} \psi_i \varepsilon_{t+k-i}\right] = \sigma_{\varepsilon}^2 \sum_{i=0}^{\infty} \psi_i \psi_{i+k}$$

#### Stationarity and invertibility

• The linear process  $Y_t = \psi(B)\varepsilon_t$  is stationary if

$$\psi(z) = \sum_{i=0}^{\infty} \psi_i z^{-i}$$

converges for  $|z| \ge 1$  (i.e. old values of  $\varepsilon_t$  are down-weighted)

• The linear process  $\pi(B) Y_t = \varepsilon_t$  is said to be *invertible* if

$$\pi(z) = \sum_{i=0}^{\infty} \pi_i z^{-i}$$

converges for  $|z| \ge 1$  (i.e.  $\varepsilon_t$  can be calculated from recent values of  $Y_t$ )

#### Linear process as a statistical model?

$$Y_t = -\pi_1 Y_{t-1} - \pi_2 Y_{t-2} \dots + \varepsilon_t + \psi_1 \varepsilon_{t-1} + \psi_2 \varepsilon_{t-2} + \psi_3 \varepsilon_{t-3} + \dots$$

- Observations:  $Y_1$ ,  $Y_2$ ,  $Y_3$ , ...,  $Y_N$
- ▶ Task: Find an infinite number of parameters from *N* observations!
- Solution: Restrict the sequence 1,  $\psi_1$ ,  $\psi_2$ ,  $\psi_3$ , ... Which gives us the famous ARMA model.