

34 FOURIER ANALYSIS & APPLICATIONS - Noise & Variability IV
 Nonstationary Processes.

Stationarity allowed us two things:

- ① We could obtain new samples from \mathcal{R} simply by waiting a long time & sampling again
- ② We could simplify the characterization of moments, since $\langle x(t)x(t+\tau) \rangle$ depended only on τ .

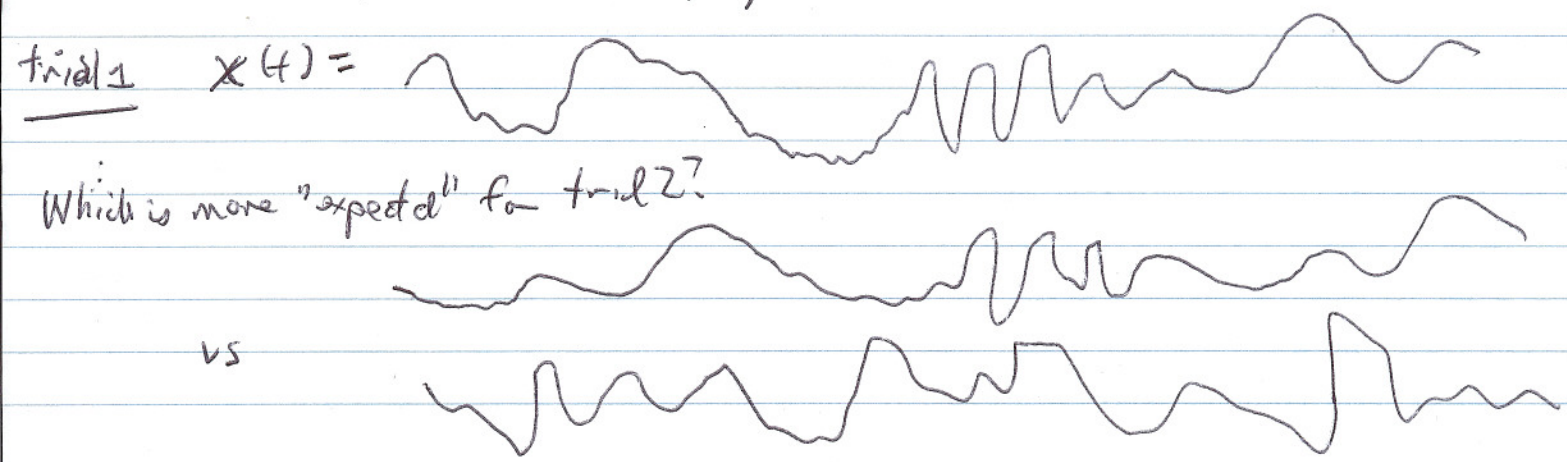
Now we relax both of these assumptions.

Still need a way to obtain multiple samples of \mathcal{R}
 e.g., "trials", each trial has a start time [or, parallel preps!]

Still need to deal with the inter-dependence of direct time-domain estimates

$$\left. \begin{aligned} &\langle x(t)x(t+\tau) \rangle \\ &\langle x(t)x(t+\tau') \rangle \\ &\langle x(t+\tau)x(t+\tau') \rangle \end{aligned} \right\} \text{are mutually dependent.}$$

Still want to try to use intuitions about general properties of biological systems, e.g., say you observe

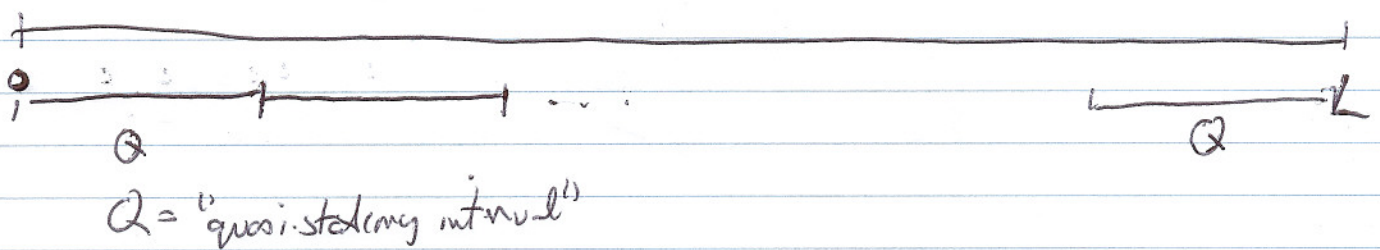


The intuition is that there are gradual changes in system properties over time, & interactions tend to be local.

This has a natural embodiment in terms of spectral estimation:

Get the best local spectral estimate, & see how it changes over time.

Full "twiddle"

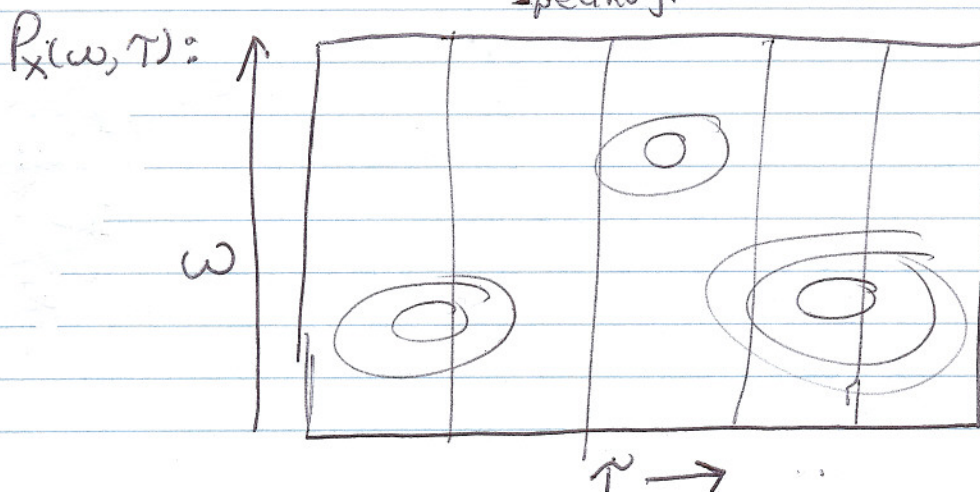


So everything we said before about getting the best estimate in length L still applies, but now over a period of time Q :

Multitaper estimate beginning of time τ , for signal from $t = \tau$ to $\tau + Q$:

$$P_x(\omega, \tau) = \frac{1}{K} \sum_{j=1}^K \left| \int_0^Q x(t + \tau) W_j(t, Q) e^{-i\omega t} dt \right|^2$$

where the $W_j(t, Q)$ are multitaper functions on $[0, Q]$
spectrogram



Frequency resolution is $\frac{2\pi K}{Q}$.

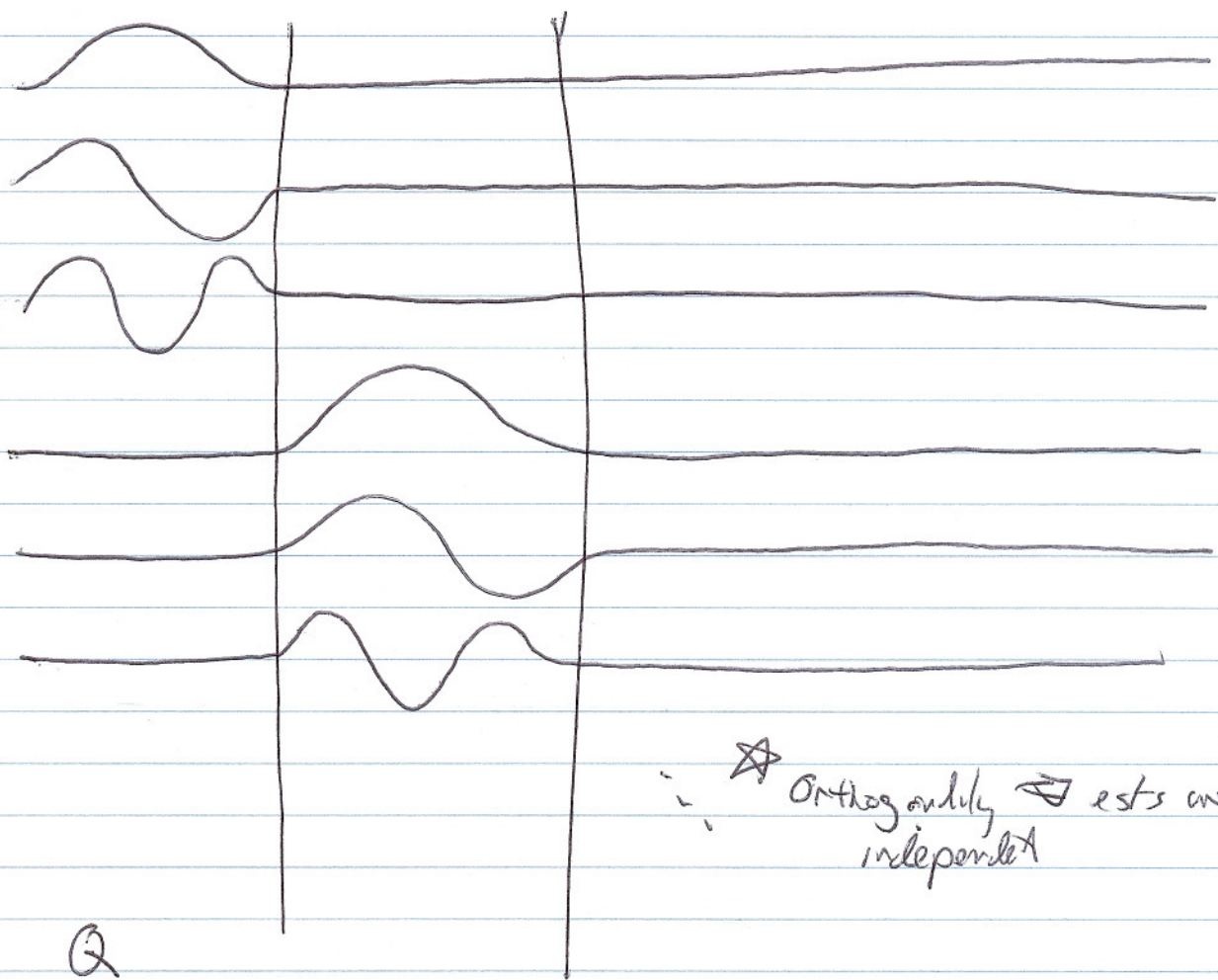
Time resolution is Q .

Larger $K \rightarrow$ better freqs but worse resolution

36

Essentially, we've made multiple local estimates of the power spectrum by multiplying $X(t)$ by tapers & their shifts.

($k=3$)



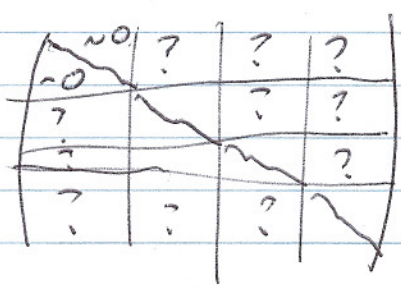
★ Orthogonality \Rightarrow ests are independent

and calculated Fourier estimates of frequencies $\gg \frac{2\pi}{Q}$.

Looked at more abstractly, we are hoping to pre full covariance matrix of X by a change of basis - "local Fourier components" -

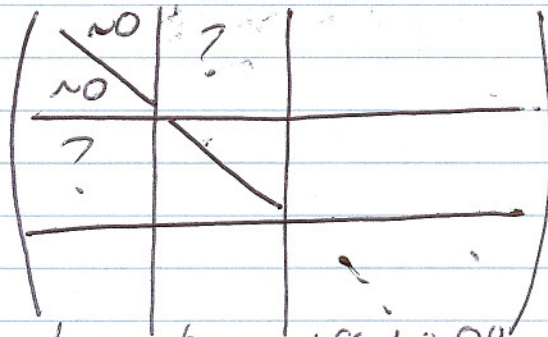
"local" time-invariance means that within each block, the covariance matrix is diagonal.

Across blocks structure is not known -



(37)

What about across the blocks?



Could (i) reanalyze with a different "Q" to look for long-term structure

(ii) "look at the results" - ROI's

(iii) spectrum of spectra dimensional reduction methods

33

Rel. of spectrogram to spectrum -

With proper normalization (e.g., Chronux spectrogram c.m., etc.)

$$\int P_x(\omega, \tau) d\tau = P_x(\omega)$$

$$\int P_x(\omega, \tau) d\omega = \langle x(\tau)^2 \rangle$$

The nice properties of spectra generalize: if X_1 & X_2 are independent

$$P_{X_1+X_2} = P_{X_1} + P_{X_2}$$

For linear filters (and are slow on a timescale Q), $x \rightarrow [L] \rightarrow y$

$$P_y(\omega, \tau) = |L(\omega)|^2 P_x(\omega, \tau)$$

even $P_y(\omega, \tau) = |L(\omega, \tau)|^2 P_x(\omega, \tau)$ if L changes slowly

Things generalize to cross-spectrum & coherence

Cross spectrum:
$$P_{xy}(\omega) \cong \frac{1}{K} \sum_{j=1}^K \left(\int_0^T x(t+\tau) w_j(t, \tau) e^{-i\omega t} dt \right) \left(\int_0^T y(t) w_j(t, \tau) e^{-i\omega t} dt \right)$$

becomes

$$P_{xy}(\omega, \tau) \cong \frac{1}{K} \sum_{j=1}^K \left(\int_0^Q x(t+\tau) w_j(t, \tau) e^{-i\omega t} dt \right) \left(\int_0^T y(t) w_j(t, \tau) e^{-i\omega t} dt \right)$$

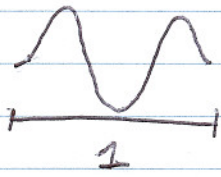
$$\frac{P_{xy}(\omega, \tau)}{\sqrt{P_x(\omega, \tau)} \sqrt{P_y(\omega, \tau)}}$$

(coherency)

$$\sqrt{P_x(\omega, \tau)} \sqrt{P_y(\omega, \tau)}$$

39

Alternative approaches

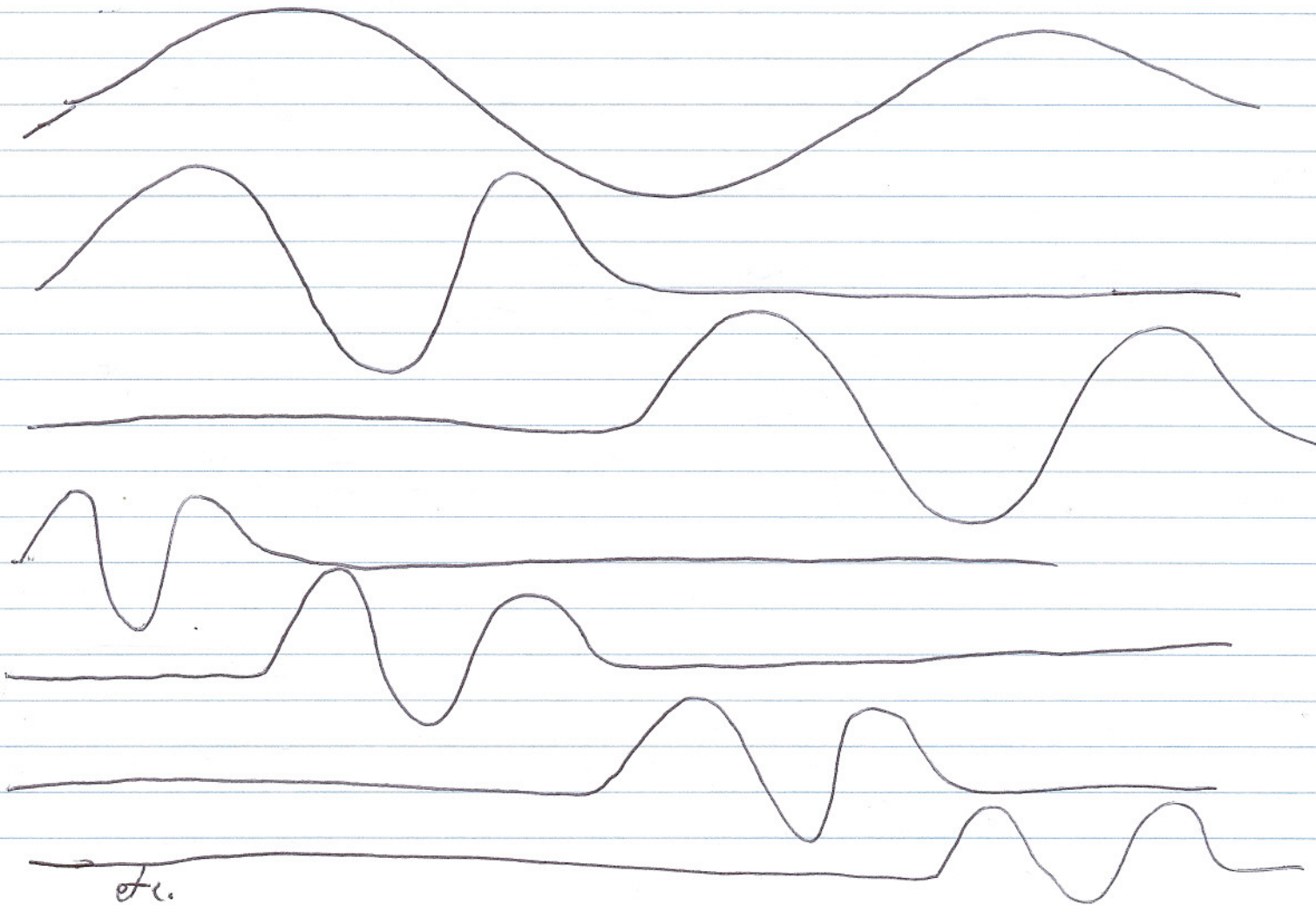
Wavelets Choose a "mother wavelet" $M(t)$, eg. 

† calculate

$$M_x(\tau, s) = \int x(t + \tau) M(t/s) dt = \int x(t) M\left(\frac{t - \tau}{s}\right) dt$$

for a range of values of s , typically in powers of 2, and for values of τ spaced by k/s . Highest $s = T$.

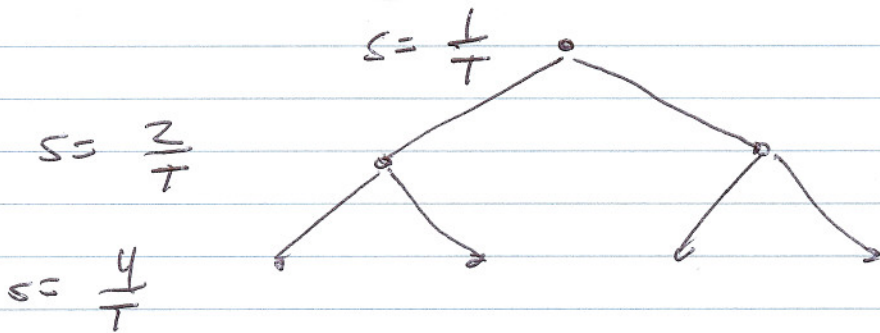
Essentially, projecting X onto



(40)

The wavelets form a basis, so

$M_X(\uparrow, S)$ has transform X into orthonormal basis set



so $\langle |M_X(\uparrow, S)|^2 \rangle$ is analogous to $P_X(\uparrow, \frac{1}{S})$

• one can also calculate $\langle M_X(\uparrow, S) \overline{M_Y(\uparrow, S)} \rangle$, etc.

Compare & contrast:

The wavelets are not orthogonal. They are not quite as confined in frequency space.

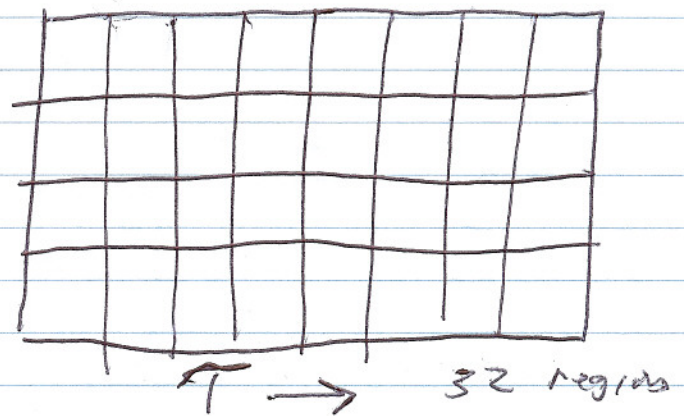
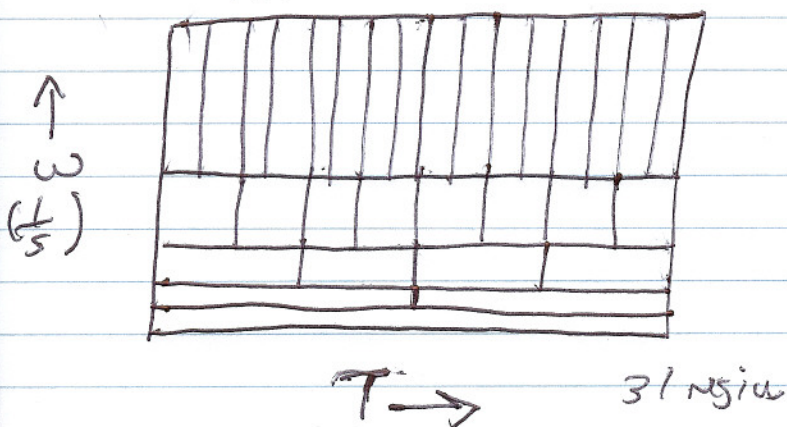
↓
so sets are not independent

↓
so you could get fooled by oscillations.

But they are an alternative way to "tile" time-freq space

WAVELETS

SPECTROGRAMS



(91)

With wavelets, increase, freq. res. \rightarrow decrease time resolution,
shape is controlled by "mother" wavelet - typical Gaussian \times sine
 \times cos
 \times polynomial

No position to exploit local time-invariance with wavelets - i.e., in MT
method, values within each "column" in the time-freq. decomposition
are \sim independent

However - if it is \sim true that shorter events have higher frequency,
then wavelets will be more effective representations.

E.g., natural scenes.

But $\begin{cases} ? \\ ? \end{cases}$ natural sounds
brain activity

\downarrow
* Group theoretic interpretation:
 $s(t) \rightarrow s(t+\tau)$
 $s(t) \rightarrow s(\alpha t)$

[two transformations]

* Look for a representation in which
these two functions are "simple"

* $t \rightarrow t+\tau$ $t \rightarrow \tau$ wavelets

$t \rightarrow \alpha t$ $t \rightarrow$ move up the stack
of wavelets

MT + wavelet ideas generalize to spatial signals

local resolution invariance

\rightarrow
+ approx.
independence
of spectra at
long range

MT

+

scaling

Wavelet

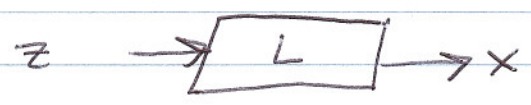
4/2

Another approach, specific to time series

"Autoregressive Models"

(Stationary case first)

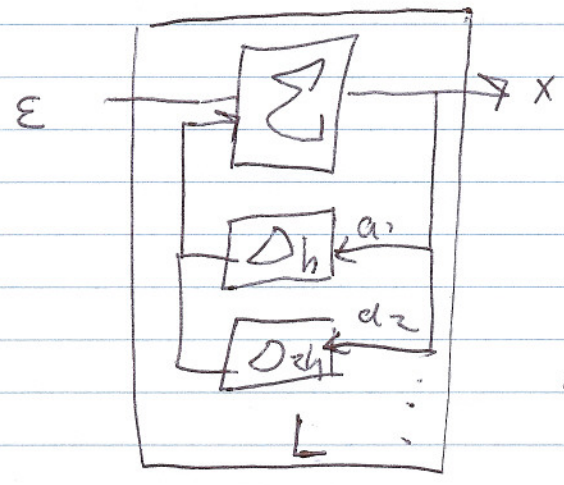
basic idea is to model the spectrum as the result of white noise acted on by a linear filter



z is unobservable. So we attempt to write

$$x(t) = a_1 x(t-h) + a_2 x(t-2h) + \dots + a_r x(t-rh) + \epsilon(t)$$

and minimize $\langle \epsilon(t) \rangle^2$



Power spectrum of $X = P_\epsilon(\omega) |L|^2$

$P_\epsilon(\omega)$ assumed to be flat, $= \langle \epsilon^2 \rangle$

$$L = \frac{1}{1 - (a_1 e^{-i\omega h} + a_2 e^{-2i\omega h} + \dots)}$$

Put $\theta = e^{i\omega h}$

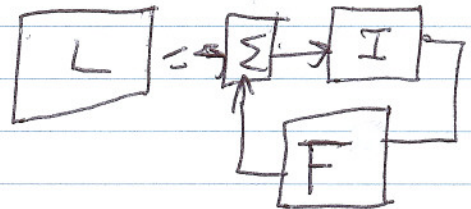
$$P_x(\omega) \approx \langle \epsilon^2 \rangle \frac{1}{|1 - \sum a_k \theta^k|^2}$$

(43)

Note that the model samples x at intervals h .

So it will only capture frequencies $\ll \frac{2\pi}{h}$. Equivalently, $L(\omega) = L(\omega + \frac{2\pi}{h})$.

Also, recall that for a feedback system,

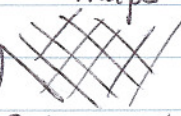


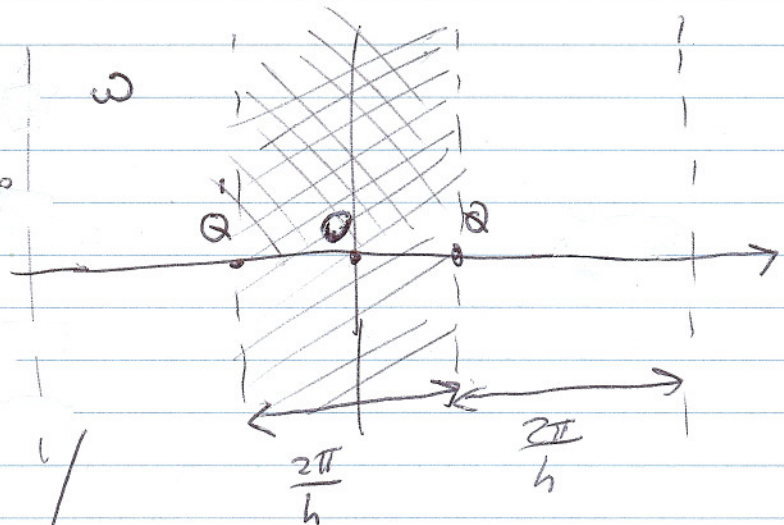
$$\tilde{L}(\omega) = \frac{L}{1 - F(\omega)}$$

is stable only if $|F(\omega)| < 1$ for all ω on real axis.

(More generally, on complex plane, $\tilde{L}(s)$ must be $< \infty$ for $\text{Im } s < 0$.)

$L(\omega)$ on the complex plane is periodic and must be bounded in lower half plane

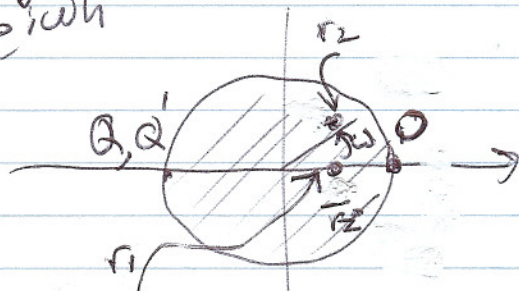
$\theta = e^{s\omega h}$ maps the region marked by  into the interior of the unit circle



$$P_x(\omega) = \frac{KE^2}{|1 - A(\theta)|^2}$$

$$\theta = e^{s\omega h}$$

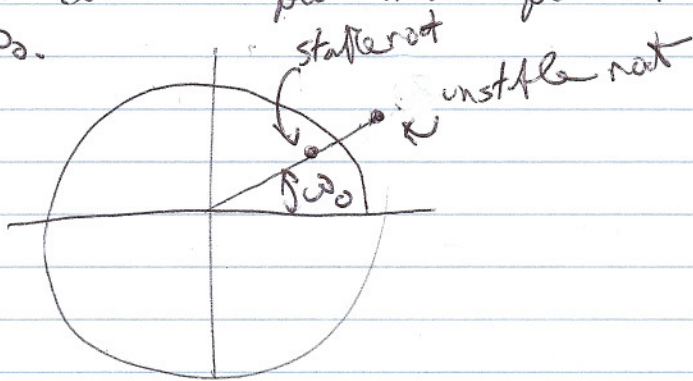
$A(\theta)$ is a polynomial, τ



$A(\theta) = 1$ can't have roots outside the unit circle. Roots can come singly (as r_1) or in cc pairs (r_2, r_2^*).

(11)

A root near the unit circle corresponds to a peak in the spectrum at the corresponding frequency ω_0 .



Bottom line - only stereotypical spectra can be well-modeled for low-degree $A(\theta)$ - peak width - peak height are coupled.

But if the spectrum is of the form $\frac{\langle \epsilon^2 \rangle}{|1 - A(\theta)|^2}$ then the procedure works well.

How to determine r and the a_r ?

"The-Walker" equation: Minimize

$$R = \sum_t \left(x(t) - (a_1 x(t-h) + a_2 x(t-2h) + \dots + a_r x(t-rh)) \right)^2$$

(= $\sum \epsilon(t)^2$)

This is a linear system: $x(t)$ is regressed against lags of $x(t)$

Criterion for adding another term: R will decrease when another term is added, but is this decrease more than one would expect if the $\epsilon(t)$'s are independently drawn from a Gaussian?

(43) Minimize R : put $\frac{\partial R}{\partial a_k} = 0$, this leads to

$$a_1 c_0 + a_2 c_1 + a_3 c_2 + \dots + a_r c_{r-1} = c_1$$

$$a_1 c_1 + a_2 c_0 + a_3 c_1 + \dots + a_r c_{r-2} = c_2$$

$$a_1 c_2 + a_2 c_1 + a_3 c_0 + \dots + a_r c_{r-3} = c_3$$

⋮

$$a_1 c_{r-1} + a_2 c_{r-2} + a_3 c_{r-3} + \dots + a_r c_0 = c_{r-1}$$

Note diagonal structure.

Several strategies for assessing tradeoff of # of params and red. R -

- reduced chi-squared

- Akaike criteria $2r + N \ln R$ (N points)

Nonstationary

fit spectral estimates to moving windows.

Multivariate

fit to multiple time series

$$x(t) = \begin{pmatrix} x_1(t) \\ \vdots \\ x_k(t) \end{pmatrix}$$

a 's are each $k \times k$ matrices

Autoregressive-moving average:

can add "moving average" terms to model, e.g.,

$$x(t) = a_1 x(t-h) + \dots + a_r x(t-rh) + b_0 \varepsilon(t) + b_1 \varepsilon(t-h) + \dots + b_s \varepsilon(t-sh)$$

46

"Granger Causality" is just multivariate autoregression --
how well can one predict $X(t)$ from the past history
of X and Y , vs how well can one predict $X(t)$ from
 $X(t)$ alone.