

# Time Series Coursework

December 16, 2021

```
[ ]: import numpy as np
import pandas as pd
from scipy.fft import fft
from scipy.linalg import toeplitz

import matplotlib.pyplot as plt
from matplotlib import cm
plt.style.use('seaborn-white')
```

## 0.0.1 1

(a) By page 52 in notes, the spectral density function of an ARMA( $p, q$ ) process is given by

$$S_X(f) = \sigma_\epsilon^2 \frac{G_\theta(f)}{G_\phi(f)}. \quad (1)$$

where

$$G_\theta(f) = |1 - \theta_{1,q}e^{-2i\pi f} - \dots - \theta_{q,q}e^{-2i\pi fq}|^2,$$
$$G_\phi(f) = |1 - \phi_{1,q}e^{-2i\pi f} - \dots - \phi_{q,q}e^{-2i\pi fq}|^2.$$

```
[ ]: def S_ARMA(f, sigma2, phis=[], thetas=[]):
    ...
    Computes the theoretical sdf for an ARMA(p, q) process.

    param f: numpy array of frequencies at which sdf should be evaluated
    param sigma2: scalar; the variance of the white noise process
    param phis: array containing the phi parameters
    param thetas: array containing the theta parameters

    return: numpy array of sdf evaluated at f
    ...
    p, q = len(phis), len(thetas)
    xis_t = np.exp(-1j*2*np.pi*np.outer(f, np.arange(1, q+1)))
    xis_p = np.exp(-1j*2*np.pi*np.outer(f, np.arange(1, p+1)))
    G_t = 1 - (thetas * xis_t).sum(1) # G_theta
    G_p = 1 - (phis * xis_p).sum(1) # G_phi
    return (sigma2 * np.abs(G_t)**2 / np.abs(G_p)**2)
```

(b) The following code implements the burn in method to generate an ARMA(2,2) process of length  $N$ : 1. Set  $X_1 = X_2 = 0$

2. For  $i = 3$  to  $100 + N$ , set  $X_i \leftarrow \phi_{1,2}X_{i-1} + \phi_{2,2}X_{i-2} + \epsilon_i - \theta_{1,2}\epsilon_{i-1} - \theta_{2,2}\epsilon_{i-2}$
3. Return  $X_{101}, \dots, X_{101+N}$ .

```
[ ]: def ARMA22_sim(phis, thetas, sigma2, N):
    """
    Simulates a Gaussian ARMA(2,2) process.

    param phis: numpy array containing the phi parameters
    param thetas: numpy array containing the theta parameters
    param sigma2: scalar; variance of the white noise process
    N: scalar; length of the simulated process

    return: numpy array of size N; realiasation of an ARMA(2,2) process
    """
    # simulate white noise process
    eps = np.random.normal(size=100+N, scale=np.sqrt(sigma2))
    X = np.zeros(100 + N)

    for i in range(2, 100+N):
        X[i] = phis[0]*X[i-1] + phis[1]*X[i-2] + eps[i] \
            - (thetas[0]*eps[i-1] + thetas[1]*eps[i-2])

    return X[-N:]
```

(c) By page 58 in notes, the periodogram for a time series  $X$  is given by

$$\hat{S}^{(p)}(f) = \frac{1}{N} \left| \sum_{i=1}^N X_t e^{-i2\pi f t} \right|^2.$$

```
[ ]: def periodogram(X):
    """
    Computes the periodogram at the Fourier frequencies for a time series X.
    """
    N = X.shape[-1]
    S = fft(X)  # computes the sum inside the norm
    return np.abs(S)**2 / N
```

By page 62 in notes, the spectual density of a  $p \times 100\%$  cosine tapered estimato is given by

$$\hat{S}^{(d)}(f) = \left| \sum_{i=1}^N h_t X_t e^{-i2\pi f t} \right|^2,$$

where the consine taper sequence  $h_t$  is as defined on page 64.

```
[ ]: def tapered_series(X, taper):
```

```

    """Computes the series ht*Xt for p*100% cosine tapered spectral estimates.
    """
N = X.shape[-1]

# Construct the H_t sequence
H = np.ones(N,)
tt = np.floor(taper*N)
H[:int(tt//2)] = (1 - np.cos(2*np.pi*np.arange(tt//2) / (tt+1))) / 2
H[N-int(tt//2):] = (1 - np.cos(
    2*np.pi*(N+1-np.arange(N-int(tt//2), N)) / (tt+1))
) / 2
H /= np.sqrt((H**2).sum())      # normalise

return H * X

def direct(X, p):
    """
    Computes the direct spectral estimate at the Fourier frequencies using the \
    p*100% cosine taper for a time series X.
    """
    hX = tapered_series(X, p)
    return np.abs(fft(hX))**2

```

(d)

(A) By page 30 in notes, an ARMA(2,2) process is stationary if all the roots of

$$\Phi(z) = 1 - \phi_{1,2}z - \phi_{2,2}z^2$$

lie outside the unit circle. Using the representation of roots  $z_1 = \frac{1}{r}e^{-i2\pi f'}$  and  $z_2 = \frac{1}{r}e^{i2\pi f'}$ , by similar arguments in page 51, we have that

$$\phi_{1,2} = 2r \cos(2\pi f'), \quad \phi_{2,2} = -r^2.$$

```
[ ]: r = 0.8
N = 128
fp = 12/128
phis = np.array([2*r*np.cos(2*np.pi*fp), -r**2])
thetas = np.array([-0.5, -0.2])

# Simulate 10000 ARMA(2,2) realisation
np.random.seed(7)
X = np.array([ARMA22_sim(phis, thetas, 1, N) for _ in range(10000)])
```

per\_single, dse\_p1\_single, dse\_p2\_single, dse\_p3\_single and dse\_p4\_single store the spectral estimates at  $f = 12/128, 32/128, 60/128$  based on the periodogram, 5%, 10%, 25% and 50% cosine tapers respectively.

```
[ ]: per_single = periodogram(X).T[[12, 32, 60]]      # periodograms

# cosine tapers
ps = [0.05, 0.1, 0.25, 0.5]
dse_p1_single = direct(X, 0.05).T[[12, 32, 60]]
dse_p2_single = direct(X, 0.1).T[[12, 32, 60]]
dse_p3_single = direct(X, 0.25).T[[12, 32, 60]]
dse_p4_single = direct(X, 0.5).T[[12, 32, 60]]
```

(B) The sample bias of a sequence of estimates  $\hat{s}_1, \hat{s}_2, \dots, \hat{s}_N$  at frequency  $f$  is given by

$$\frac{1}{N} \sum_{i=1}^N \hat{s}_i - s$$

where  $s$  is the true spectral density at frequency  $f$ .

```
[ ]: fs = np.array([12/128, 32/128, 60/128])

estimates = np.dstack([per_single, dse_p1_single, dse_p2_single, dse_p3_single,
                     dse_p4_single])
true_sdf = S_ARMA(fs, sigma2=1, phis=phis, thetas=thetas)
sample_bias_vals = estimates.mean(1).T - true_sdf

# Sample bias table for each estimator when r=0.8
sample_bias = pd.DataFrame(sample_bias_vals, columns=['f=' + str(int(f*128)) +
                                                       '/128' for f in fs])
sample_bias.rename_axis(columns='freq', inplace=True)
sample_bias.set_axis(['periodogram'] + [f'{int(p*100)}% cosine taper' for p in
                                         ps], axis='index', inplace=True)
print('Sample bias of different spectral density estimators (r=0.8):')
sample_bias
```

Sample bias of different spectral density estimators (r=0.8):

```
[ ]: freq          f=12/128  f=32/128  f=60/128
periodogram      -2.561065  0.129317  0.051919
5% cosine taper  -2.318748  0.068238  0.005019
10% cosine taper -2.041909  0.018517  0.000321
25% cosine taper -1.721829 -0.004065  0.000002
50% cosine taper -1.513529 -0.007609  0.000101
```

(C)

```
[ ]: N = 128
fp = 12/128
thetas = [-0.5, -0.2]      # unchanged

rs = np.arange(0.8, 1.0, 0.01)
```

```

phis = np.array([2*rs*np.cos(2*np.pi*fp), -rs**2]).T    # 20x2 array, recording
    ↳phi values for each r

# sample 10000 realisations for each sets of phi
np.random.seed(7)
XX = np.array([[ARMA22_sim(phi, thetas, 1, N) for _ in range(10000)] for phi in
    ↳phis])      # shape=(20, 10000, 128)

# Compute the spectral density estimates using the same 5 estimators as above
perXX = periodogram(XX)[..., [12, 32, 60]]      # shape=(20, 10000, 3)
dse_p1 = direct(XX, 0.05)[..., [12, 32, 60]]
dse_p2 = direct(XX, 0.1)[..., [12, 32, 60]]
dse_p3 = direct(XX, 0.25)[..., [12, 32, 60]]
dse_p4 = direct(XX, 0.5)[..., [12, 32, 60]]

# Compute the true spectral density
true_sdf = np.array([S_ARMA(
    fs, sigma2=1, phis=phi, thetas=thetas
) for phi in phis])  # 20x3 array

# Compute the sample bias
sample_mean = np.stack((perXX, dse_p1, dse_p2, dse_p3, dse_p4)).mean(2)
sample_bias = sample_mean - true_sdf      # shape=(5, 20, 3)

```

(D)

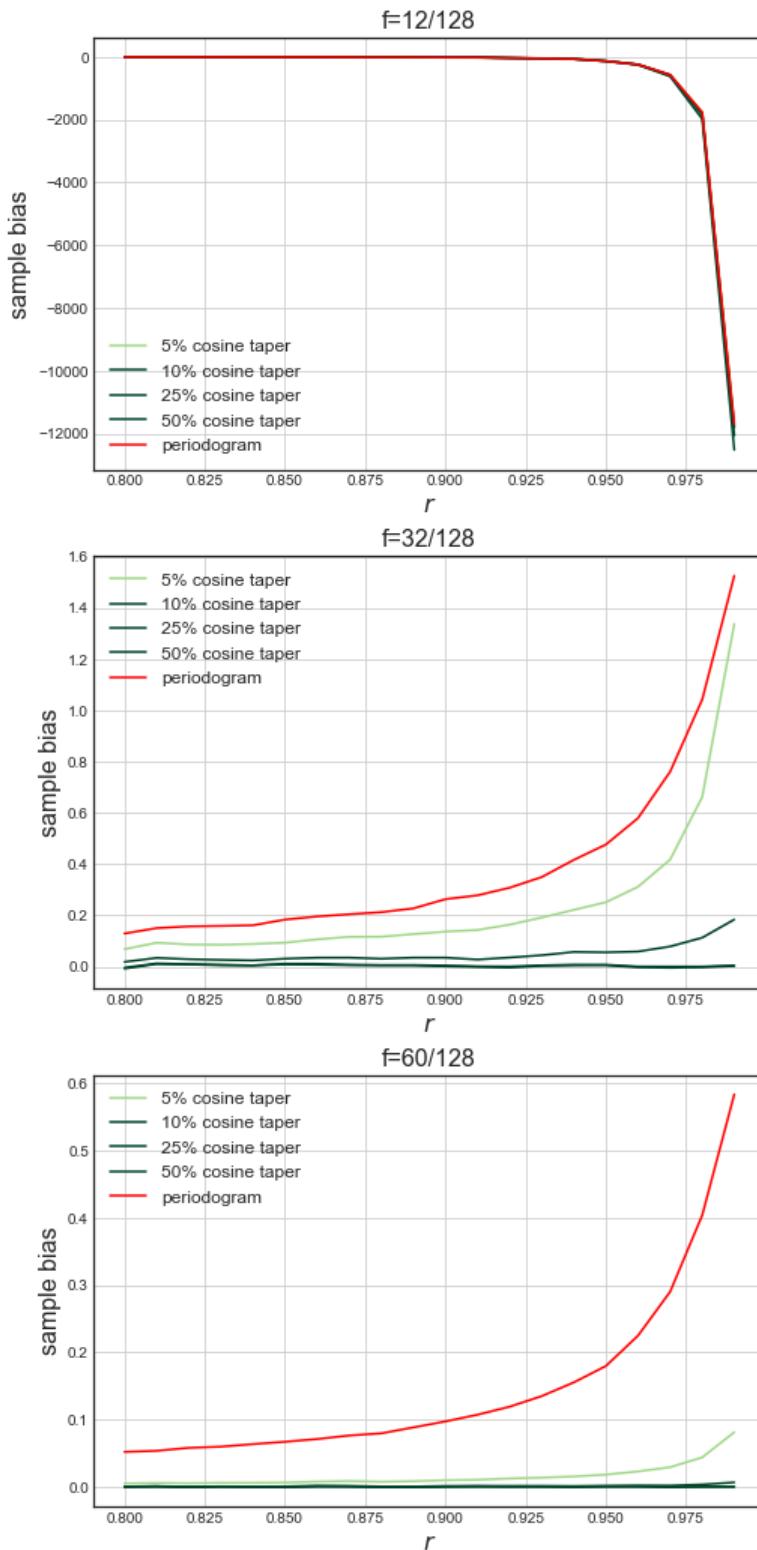
```

[ ]: fig = plt.figure(figsize=(8, 18))

for i in range(3):
    plt.subplot(3, 1, i+1)
    for ip in range(4):
        plt.plot(rs, sample_bias[ip+1, :, i], c=cm.YlGn(1.25*ip+0.4), ↳
            ↳label=f'{int(ps[ip]*100)}% cosine taper')
        plt.plot(rs, sample_bias[0, :, i], 'r-', label='periodogram')
        plt.xlabel('r', fontsize=16)
        plt.ylabel('sample bias', fontsize=16)
        plt.title(f'f={int(fs[i]*128)}/128', fontsize=16)
        plt.grid()
        plt.legend(fontsize=12)
    # fig.suptitle('Comparision of sample bias for different spectral estimators', ↳
    ↳fontsize=20)
    txt = 'Fig 1: Comparision of sample bias for different spectral estimators'
    plt.figtext(0.5, .92, txt, wrap=True, horizontalalignment='center', fontsize=25)
plt.show()

```

Fig 1: Comparision of sample bias for different spectual estimators



(e) - As shown in Figure 1, at the oscillating frequency  $f = 12/128$ , spectral density is underestimated by all five estimators, while at the two frequencies away from the oscillating frequency, spectral density is overestimated. This suggests the existence of sidelobe leakage.

- From the two Figure 2 and 4 below, one can see that for larger  $r$  the process has larger dynamic range. As a result, for each estimator there are more sidelobe leakage as  $r$  gets close to 1, resulting in significantly larger (absolute) sample bias, as shown in Figure 1.
- At the oscillating frequency  $f = 12/128$ , the periodogram and the tapered estimators have similar sample bias.
- At the two frequencies away from the oscillating frequency, the tapered estimators have smaller bias than periodogram. In particular, more tapering (larger  $p$ ) results in larger bias reduction.
- At  $f = 32/128$ , the 25% and 50% taperered estimators have almost zero bias. At  $f = 60/128$  (even further away from the oscillating frequency), all four cosine tapered estimates have very low bias for  $r$  that is not extremely close to 1.

```
[ ]: ## True sdf
fss = np.arange(128) / 128
sdf = np.array([S_ARMA(fss, sigma2=1, phis=phi, thetas=thetas) for phi in phis])

plt.figure(figsize=(19, 5))
plt.subplot(121)
for k in range(20):
    plt.plot(fss, sdf[k], c=cm.Reds(k/20), lw=1.5)
sm = plt.cm.ScalarMappable(cmap=cm.Reds)

for f in fs:
    plt.axvline(x=f, alpha=0.6, c='k', linestyle='--',
    label=r'$f=$'+f'{int(f*128)}/128')
cbar = plt.colorbar(sm)
cbar.set_label(r'$r$', rotation=0, fontsize=16, loc='bottom')

plt.xlabel(r'$f$', fontsize=14)
plt.ylabel(r'$S(f)$', fontsize=14)
plt.figtext(0.26, 0.001, '\nFig 2: True spectral density', wrap=True,
horizontalalignment='center', fontsize=20)
# plt.title('True spectral density', fontsize=16)
plt.legend(loc=9, bbox_to_anchor=(0.7, 1), fontsize=12)
plt.grid()

## Dynamic range
plt.subplot(122)
dynamic_range = 10 * np.log10(sdf.max(-1) / sdf.min(-1))
plt.plot(rs, dynamic_range)
```

```

# plt.title(r'Dynamic range of the process with different $r$', fontsize=14)
plt.figtext(0.55, 0.001, '\nFig 3: Dynamic range of the process with different',
            r'$r$', wrap=True, fontsize=20)
plt.xlabel(r'$r$', fontsize=14)
plt.grid()
plt.show()

```

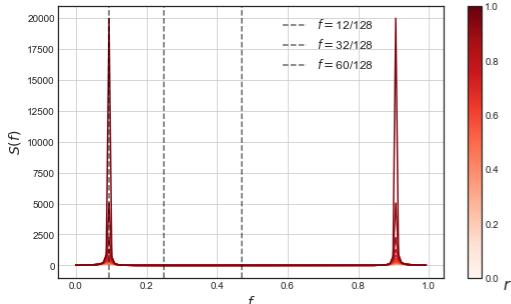


Fig 2: True spectral density

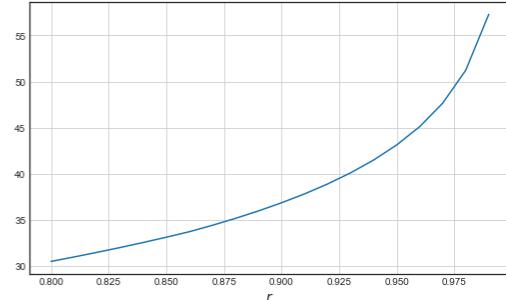


Fig 3: Dynamic range of the process with different  $r$

## 0.0.2 2

(a)

```

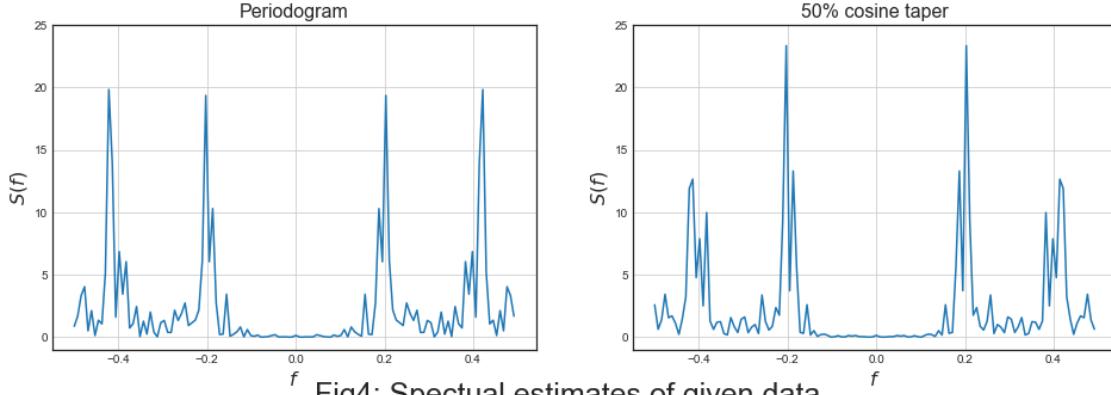
[ ]: # Load data
X2 = np.genfromtxt('154.csv', delimiter=',')
N = X2.shape[-1]

# Compute periodogram and 50% cosine tapered estimate shifted to (-1/2, 1/2]
per = np.fft.fftshift(periodogram(X2))
dse = np.fft.fftshift(direct(X2, 0.5))
titles = ['Periodogram', '50% cosine taper']

# Plot the estimates
fig = plt.figure(figsize=(16, 5))
for i, est in enumerate([per, dse]):
    plt.subplot(1, 2, i+1)
    plt.plot(np.arange(-1/2, 1/2, 1/128), est)
    plt.xlabel(r'$f$', fontsize=16)
    plt.ylabel(r'$S(f)$', fontsize=16)
    plt.title(titles[i], fontsize=16)
    plt.ylim(-1, 25)
    plt.grid()

txt = '\nFig4: Spectual estimates of given data'
plt.figtext(0.5, 0.01, txt, wrap=True, horizontalalignment='center',
            fontsize=25)
plt.show()

```



(b) Using the notations as page 65-66 in notes, for **Yule-Walker method (untapered)**, I will construct the arrays and then solve for

$$\hat{\phi}_p = \hat{\Gamma}_p^{-1} \hat{\gamma}_p,$$

where  $\hat{\Gamma}_p^{-1}$  is the symmetric Toeplitz matrix with  $(\hat{s}_0, \dots, \hat{s}_{p-1})$  as its first row,  $\hat{\gamma}_p = (\hat{s}_1, \dots, \hat{s}_p)^T$ , and  $\{\hat{s}_\tau\}$  denotes the sample autocovariance sequence of the data  $\{X_t\}$ .

For **Yule-Walker method with 50% cosine taper**, compute the autocovariance sequence of the tapered series  $\{h_t X_t\}$ , let this be  $\{\hat{s}_\tau\}$ , then proceed the same as untapered Yule-Walker.

For **approximate maximum likelihood**, using the same notations as section 5.2 and 5.3 in notes, I will solve for

$$\hat{\phi} = (F^T F)^{-1} F^T \mathbf{X}, \quad \sigma_\epsilon^2 = \frac{(\mathbf{X} - F\hat{\phi})^T (\mathbf{X} - F\hat{\phi})}{N - p}$$

where  $\mathbf{X} = (X_{p+1}, \dots, X_N)^T$ ,  $F$  is the Toeplitz matrix with  $(X_p, X_{p-1}, \dots, X_1)$  as its first row and  $(X_p, X_{p+1}, \dots, X_{N-1})^T$  as its first column.

```
[ ]: def autocov_seq_hat(X, symmetry=False):
    """Compute the sample autocovariance for a time series."""
    N = X.shape[-1]
    taus = np.arange(-N, N) if symmetry else np.arange(N+1)
    Ts = np.abs(taus).astype(int)
    return np.array([(X[:N-T]*X[T:]).sum() / N for T in Ts])

### Yule-Walker (untapered)
def estimators_YW(X, p):
    """
    Fit an AR(p) model to a given times series using Yule-Walker.

    returns phi_hat: numpy array of length p, containing the phi estimates
    returns sigma2_hat: scalar, estimated variance of the white noise process
    """

```

```

stau_hat = autocov_seq_hat(X)
Ghat = toeplitz(stau_hat[:p])
gamma_hat = stau_hat[1:p+1]
phip_hat = np.linalg.inv(Ghat).dot(gamma_hat)
sigma2_hat = stau_hat[0] - phip_hat.dot(stau_hat[1:p+1])
return phip_hat, sigma2_hat

### Yule-Walker with 50% cosine taper
def estimators_YW_tapered(X, p, taper=0.5):
    """
    Fit an AR(p) model to a given times series using Yule-Walker with \
    p*100% cosine taper.
    """
    hX = tapered_series(X, taper)
    N = hX.shape[-1]
    stau_hat = autocov_seq_hat(hX) * N
    Ghat = toeplitz(stau_hat[:p])
    gamma_hat = stau_hat[1:p+1]
    phip_hat = np.linalg.inv(Ghat).dot(gamma_hat)
    sigma2_hat = stau_hat[0] - phip_hat.dot(stau_hat[1:p+1])
    return phip_hat, sigma2_hat

### Approximate MLE
def estimators_MLE(X, p):
    """
    Fit an AR(p) model to a given times series using approximate \
    maximum likelihood.
    """
    N = X.shape[-1]
    F = toeplitz(X[p-1:N-1], X[p-1::-1])
    XX = X[p:]
    phat = np.linalg.inv(F.T @ F) @ F.T @ XX
    A = XX - F.dot(phat)
    sigma2_hat = A.T @ A / (N-p)
    return phat, sigma2_hat

```

(c)

```
[ ]: def reduced_AIC(X, p, method='YW'):
    """Compute the AIC score for the three methods above."""
    fmap = {
        'YW': estimators_YW,
        'YW taper': estimators_YW_tapered,
        'MLE': estimators_MLE,
    }
```

```

N = X.shape[-1]
s2 = fmap[method](X, p)[-1]
return 2*p + N * np.log(s2)

methods = ['YW', 'YW taper', 'MLE']
YW_data = np.array([reduced_AIC(X2, pp) for pp in range(1, 21)])
YW_taper_data = np.array([reduced_AIC(X2, pp, 'YW taper') for pp in range(1, 21)])
MLE_data = np.array([reduced_AIC(X2, pp, 'MLE') for pp in range(1, 21)])

df = pd.DataFrame(
    {'YW': YW_data, 'YW taper': YW_taper_data, 'MLE': MLE_data},
).rename_axis('p').rename(lambda x: x+1, axis='index')
print('AIC scores:')
df

```

AIC scores:

```
[ ]:          YW      YW taper        MLE
p
1  94.783802  101.207951  95.700781
2  88.539991  91.802297  90.444720
3  58.736744  57.057308  59.008583
4  1.434667  -13.417359  -2.080725
5  3.402804  -11.420202  -1.019231
6  3.883213  -10.309071  -0.916525
7  5.722645  -8.343541  0.483330
8  7.525434  -7.086085  3.029029
9  7.697199  -5.566404  1.567323
10 9.214986  -4.194696  -0.494642
11 11.158646  -2.194762  1.981352
12 12.929766  -0.436913  4.685874
13 12.199077  -0.184707  4.186692
14 14.188579  1.791134  6.299942
15 15.620487  3.224196  8.361404
16 17.616934  5.189858  5.611818
17 19.614910  7.126054  8.018493
18 19.785123  7.764564  9.227051
19 20.869850  6.129727  8.630923
20 22.256291  8.124802  8.982258
```

(d)

```
[ ]: def best_p(ps, AICs):
    return ps[np.argmin(AICs, -1)]

ps = np.arange(1, 21)
best_ps = best_p(ps, [YW_data, YW_taper_data, MLE_data])
```

```

for i in range(3):
    print(f'Best choice of p based on {methods[i]} method = {best_ps[i]}')

YW_coeffs = estimators_YW(X2, best_ps[0])
YW_taper_coeffs = estimators_YW_tapered(X2, best_ps[1])
MLE_coeffs = estimators_MLE(X2, best_ps[2])

print('----- Estimated parameter values based on the best p -----')
print(f'Yule-Walker: Phi = {YW_coeffs[0]}, sigma^2 = {YW_coeffs[1]}')
print(f'Yule-Walker with 50% taper: Phi = {YW_taper_coeffs[0]}, sigma^2 = '
    f'{YW_taper_coeffs[1]}'')
print(f'Approximate MLE: Phi = {MLE_coeffs[0]}, sigma^2 = {MLE_coeffs[1]}')

```

Best choice of p based on YW method = 4  
 Best choice of p based on YW taper method = 4  
 Best choice of p based on MLE method = 4  
 ----- Estimated parameter values based on the best p -----  
 Yule-Walker: Phi = [-0.81623113 -0.71385968 -0.79215113 -0.60892904], sigma^2 = 0.9500015475954233  
 Yule-Walker with 50% taper: Phi = [-0.82177909 -0.7716625 -0.82390842 -0.65751518], sigma^2 = 0.8459262307072268  
 Approximate MLE: Phi = [-0.83730632 -0.72914215 -0.81022676 -0.63924543], sigma^2 = 0.9242657257687535

(e)

```

[ ]: plt.figure(figsize=(7, 5))
for i, data in enumerate([YW_coeffs, YW_taper_coeffs, MLE_coeffs]):
    plt.plot(
        np.arange(-1/2, 1/2, 1/128),
        S_ARMA(f=np.arange(-1/2, 1/2, 1/128), sigma2=data[1], phis=data[0]),
        label=methods[i],
        c=cm.gist_rainbow(i/3)
    )
plt.xlabel(r'$f$', fontsize=14)
plt.ylabel(r'$S(f)$', fontsize=14)
plt.legend(fontsize=14)
plt.xlim(-1/2, 1/2)
txt = '\nFig5: Spectual density functions of three fitted models'
plt.figtext(0.5, 0.01, txt, wrap=True, horizontalalignment='center',
            fontsize=20)
plt.grid()
plt.show()

```

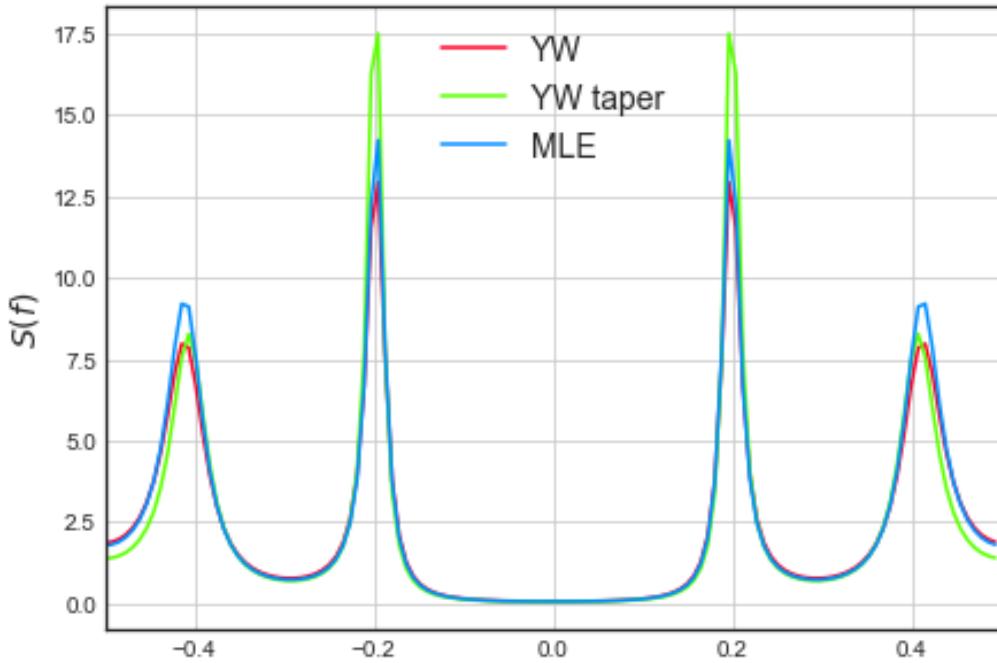


Fig5: Spectral density functions of three fitted models

### 0.0.3 3

(a) The forecasting problem is to solve for  $X_t(1), \dots, X_t(10)$  using the AR( $p$ ) model selected by approximate maximum likelihood method based on  $X_1, \dots, X_{118}$ .

Using results on page 92 in the notes, we may forecast the values by setting the future innovation terms to zero, then solving the equations:

$$\begin{aligned} X_t(1) &= \phi_{1,p}X_t + \phi_{2,p}X_{t-1} + \dots + \phi_{p,p}X_{t-p+1} \\ X_t(2) &= \phi_{1,p}X_t(1) + \phi_{2,p}X_t + \dots + \phi_{p,p}X_{t-p+2} \\ &\vdots \\ X_t(10) &= \phi_{1,p}X_t(9) + \phi_{2,p}X_t(8) + \dots + \phi_{p,p}X_{t-p+10} \end{aligned}$$

```
[ ]: T = 118
tX = X2[:T]
AICs = np.array([reduced_AIC(tX, pp, 'MLE') for pp in range(1, 21)])
p = best_p(range(1, 21), AICs)
phis, sigma2 = estimators_MLE(tX, p)
print(f'Best p value based on first {T} data points = {p}')

# In each iteration, append the newest predicted value to the end of \
# the truncated series, the iteratively using the most present p data \
```

```

# points in the series.
for t in range(N-T):
    np.append(tX, phis.dot(tX[-1:-(p+1):-1]))

df = pd.DataFrame(
    {'Predicted data': tX[-(N-T):], 'True data': X2[-(N-T):]}
).rename(lambda x: 'X_'+str(x+1+T), axis='index')
df

```

Best p value based on first 118 data points = 4

```
[ ]: Predicted data  True data
X_119      2.20780   0.44689
X_120     -0.52338  -1.16920
X_121     -1.86930   0.80533
X_122     -1.58830  -0.52534
X_123      1.12930   0.11856
X_124      1.37470  -1.28770
X_125     -0.10985   3.54240
X_126     -1.29020  -1.16590
X_127      1.33350  -0.82991
X_128     -0.82480   0.27411
```

(b)

```
[ ]: ls = N-T
m = 999
eps = np.random.normal(scale=np.sqrt(sigma2), size=(m, ls))

np.random.seed(60046)
Xsims = np.zeros((m, N))
Xsims[:, :T] = np.tile(tX, (m, 1))
for t in range(ls):
    Xsims[:, T+t] = Xsims[:, -1:-(p+1):-1].dot(phis) + eps[:, t]
```

```
[ ]: pred_interval = np.sort(Xsims[:, -10:], axis=0)[[49, 949]]
indices = np.arange(T+1, N+1)

plt.figure(figsize=(8, 6))
plt.plot(indices, X2[-ls:], 'ko', markersize=6, label='True data')
plt.plot(indices, pred_interval.T, 'b-', alpha=0.4)
plt.fill_between(indices, pred_interval[0], pred_interval[1], alpha=0.2, label='90% prediction interval')
plt.ylabel(r'$X_t$', fontsize=16)
plt.xlabel(r'$t$', fontsize=16)
txt = '\nFig7: X119 to X128: true data with 90% prediction interval'
plt.figtext(0.5, 0.01, txt, wrap=True, horizontalalignment='center', fontsize=20)
```

```
plt.grid()  
plt.legend(loc='upper left', fontsize=16)  
plt.show()
```

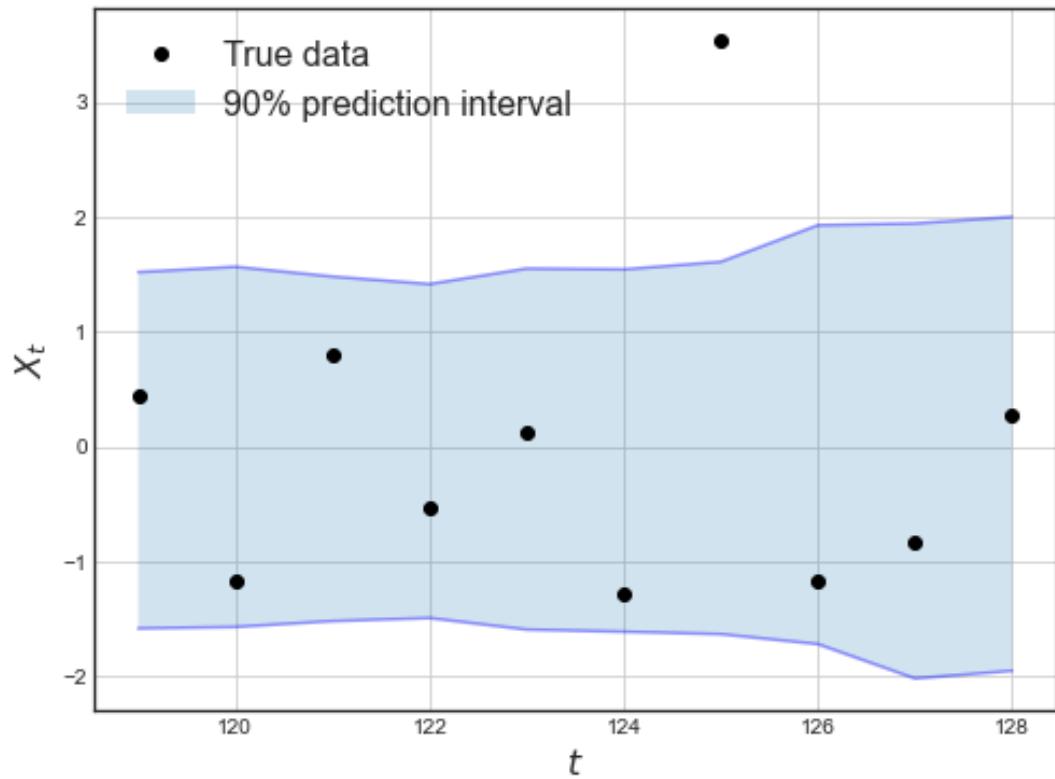


Fig7: X119 to X128: true data with 90% prediction interval