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Time Series Analysis using Auto Regressive models

11. Task for Model Analysis I, 2023/24

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1 Introduction

In physics we often encounter time series data as we measure the evolution of a system in time. Today we'll be having a look at one of the plethora of methods in our time series analysis arsenal, called Auto Regressive (AR) models. Then we'll apply this method to a few different datasets to study its performance, applicability and behavior. AR models are a class of linear models that predict the future values of a time series based on its past values. Based on the coefficients of the model, we can infer the underlying dynamics of the system or find the signals spectrum.

An AR model of order p is defined as:

$$X_t = \sum_{i=1}^p a_i X_{t-i} + \epsilon_t, \quad (1)$$

where X_t is the time series at time t , a_i are the coefficients of the model and ϵ_t is the error term which we assume to be uncorrelated and with zero mean. In our idealized case, we'll only have a short look at how noise affects the model. For the coefficients a_i we can without loss of generality divide by the first coefficient, a_1 , which yields a model with a leading coefficient of 1. The order of the model, p , is a hyperparameter that determines how many past values of the time series we use to predict the future value. The model is commonly used in the context of machine learning but today we'll be solving for the coefficients more hands-on, with out explicit use of optimization algorithms.

A popular method to determine the AR model coefficients are the Yule-Walker equations. These are a set of p linear equations that can be solved for the coefficients which are given by the autocorrelation of the time series $\{X_t\}$. The autocorrelation is defined as:

$$r_n = \sum_{i=1}^p a_i r_{n-i}, \quad (2)$$

where r_n is the autocorrelation at lag n . Essentially we need to solve a Toeplitz system of equations. The system of equations is given by:

$$\mathbf{R}\mathbf{a} = \mathbf{r}, \quad (3)$$

where \mathbf{R} is the autocorrelation matrix, given as:

$$\mathbf{R} = \begin{bmatrix} r_0 & r_1 & \cdots & r_{p-1} \\ r_1 & r_0 & \cdots & r_{p-2} \\ \vdots & \vdots & \ddots & \vdots \\ r_{p-1} & r_{p-2} & \cdots & r_0 \end{bmatrix}, \quad (4)$$

and \mathbf{r} and \mathbf{a} are the autocorrelation vector and coefficient vector respectively, given as:

$$\mathbf{r} = \begin{bmatrix} r_1 \\ r_2 \\ \vdots \\ r_p \end{bmatrix} \quad \text{and} \quad \mathbf{a} = \begin{bmatrix} a_1 \\ a_2 \\ \vdots \\ a_p \end{bmatrix}. \quad (5)$$

Stability of the model is ensured by the roots of the characteristic polynomial of the model which must lie inside the unit circle on the complex plane. This is something we'll have to keep in mind when analyzing the results, as any roots outside the unit circle will make the model unstable.

2 Task

2.1 Spectra of Signals

The instructions want us to plot the spectra of a few different signals that have been provided. The main comparison should be between the spectra given by the AR model and the spectra given by the Fourier transform. The FFT algorithm is truly a marvel of modern mathematics and computing so it will definitely be a tough competitor. The signals that have been provided are plotted in Figure 1.

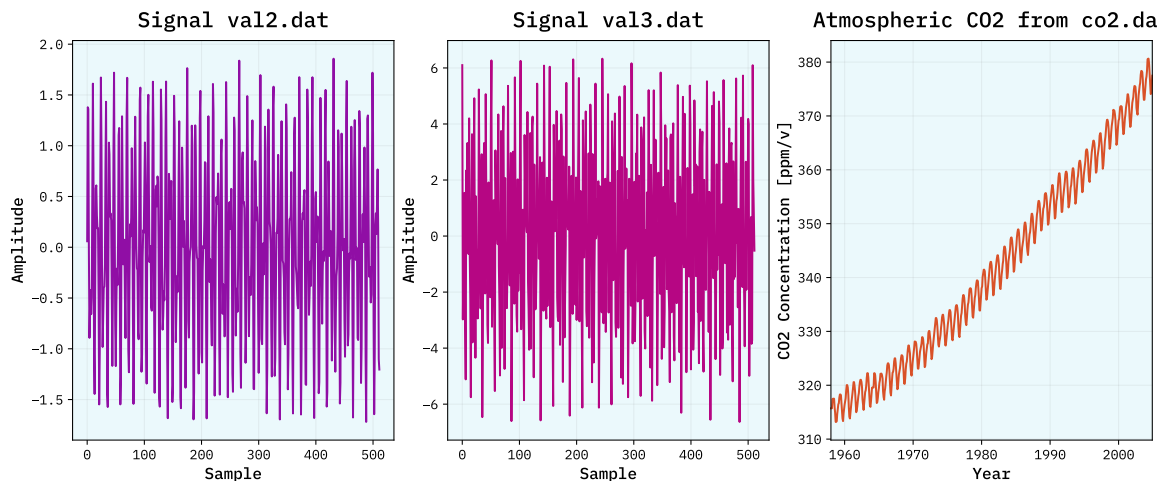


Figure 1: The signals that we'll be analyzing in the first subtask.

The instructions also want us to explore the various properties of the AR model and how it behaves when we change the order of the model as well as the number of points we use to evaluate the spectra. The spectra can be calculated using the following formula:

$$S(f) = \frac{1}{|1 - \sum_{i=1}^p a_i e^{-2\pi i f i}|^2}, \quad (6)$$

however I used the `scipy.signal.freqz` for ease of use and to avoid any mistakes.

2.2 Forecasting Signals

In the second subtask we've been provided some additional signals that we'll be using in combination with the signals from the first subtask to study the performance of the AR model in forecasting. The supplied signals are plotted in Figure 2.

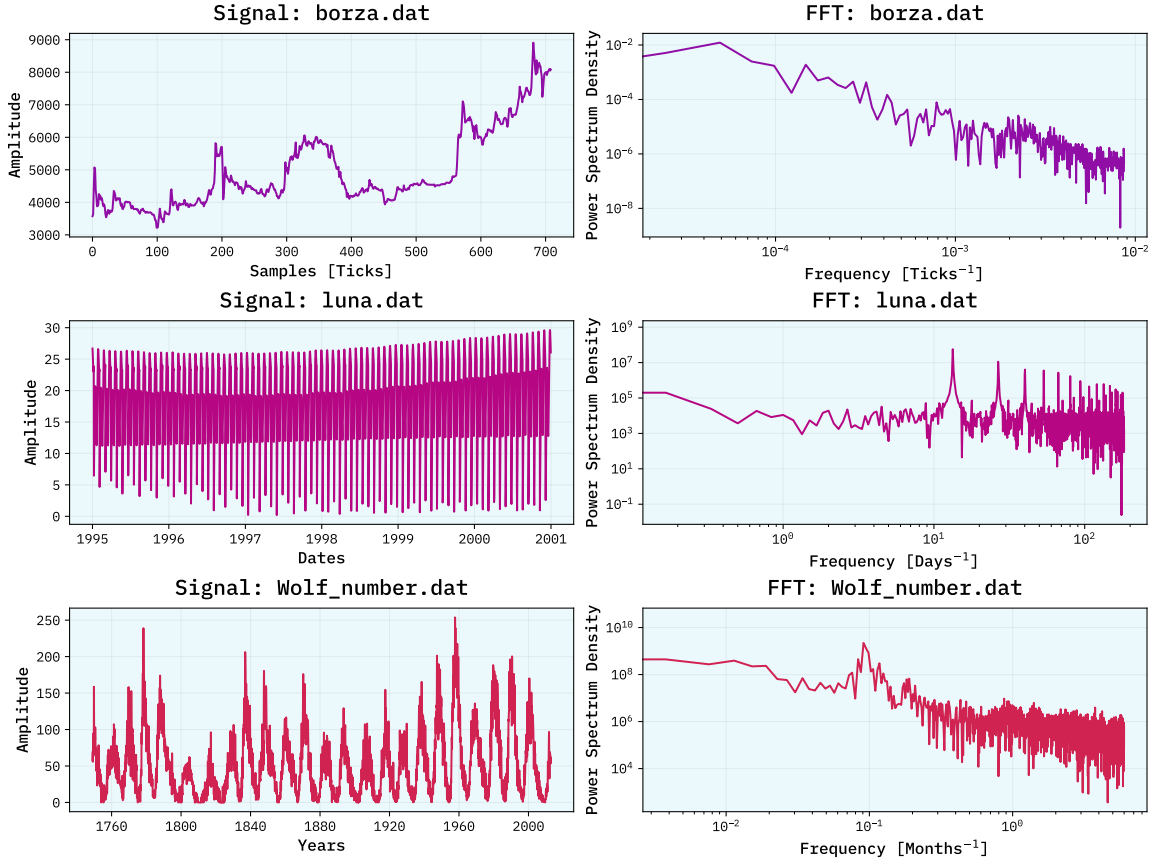


Figure 2: The signals that we'll be adding to our analysis in the second subtask.

They represent a stock's value, ephemeris data for the moon and Wolf sunspot numbers. The instructions want us to forecast the signals using the AR model and compare the results with the actual data. We can do so by providing half of the data to the model and then comparing the forecasted values with the actual values of the other half. We should also have a look at how noise affects the model and the forecasted values.

3 Solution Overview

Due to my speedrunning of these tasks this task lacks any extra bells and whistles. I implemented the Yule-Walker method using `numpy` and `scipy` by making a class `YuleWalker` that can be used to fit the model and evaluate PSDs. It takes inputs of the time series and the order of the model. I really like classes in Python. They really serve well as containers for variables and their associated functions. The class also has a method to forecast the time series using the AR model. The results are presented in the following sections. Before we proceed we need to detrend some of the signals. This was most commonly done by subtracting the mean of the signal, however I encountered two special cases. The first was the CO2 atmospheric concentration signal to which I fit a square polynomial and subtracted the fit from the signal. The second signal was the stock value signal which I detrended by the linear fit of the signal. The signals with their respective trends are plotted in Figures 3 and 4.

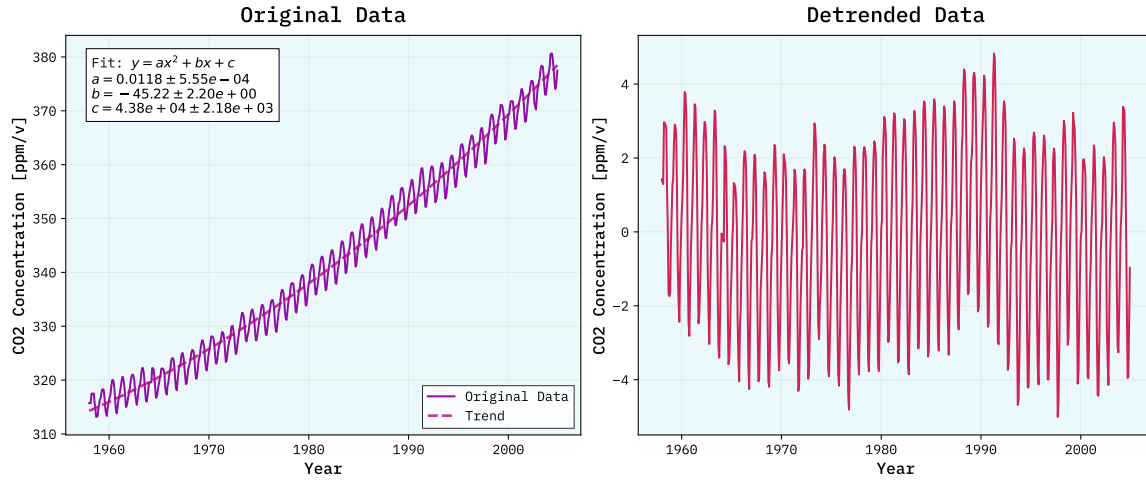


Figure 3: The CO2 atmospheric concentration signal with a square polynomial fit.

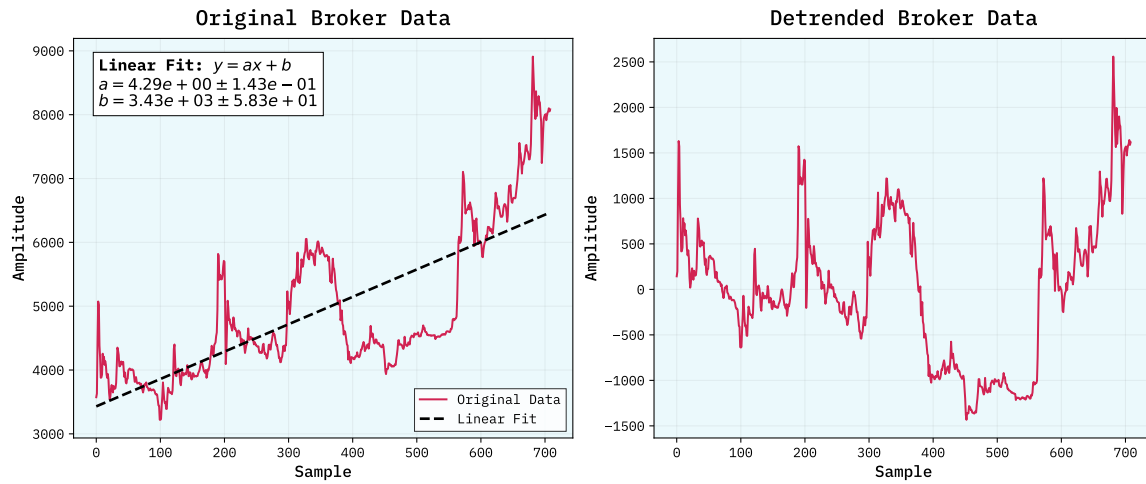


Figure 4: The stock value signal with a linear fit.

4 Results

4.1 Spectra of Signals

As described above I used the `scipy.signal.freqz` function to evaluate the spectra of the signals. The results are plotted in Figures 5, 6 and 7. The AR model seems to do a fantastic job at approximating the spectra of the signals. The absolute scale of the spectra are larger than expected however they match up relatively well with the FFT spectra. What is definitely visible are the wider peaks in the AR model spectra compared to the FFT spectra. I think it makes sense to study the width of the peaks a bit more in depth.

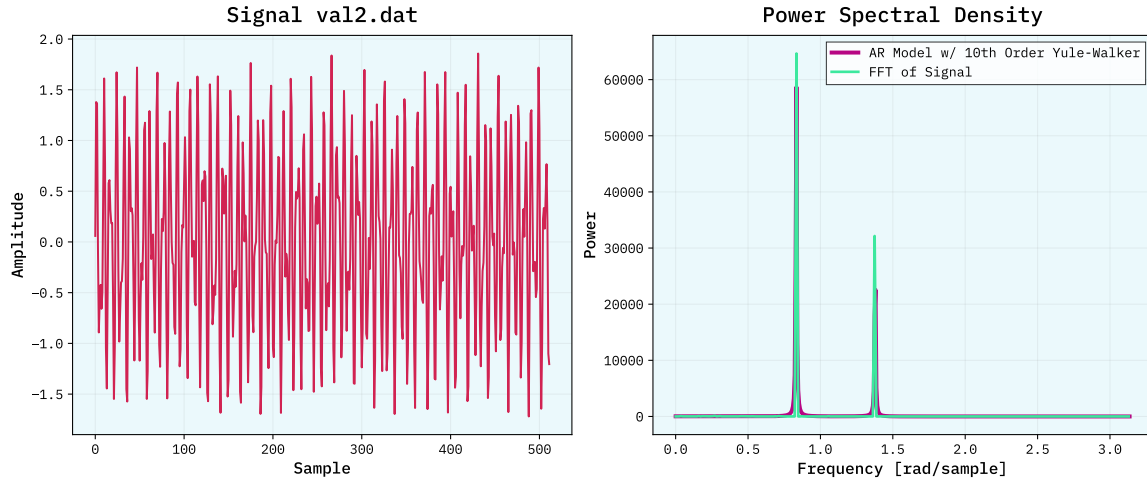


Figure 5: The spectra signal val2.dat.

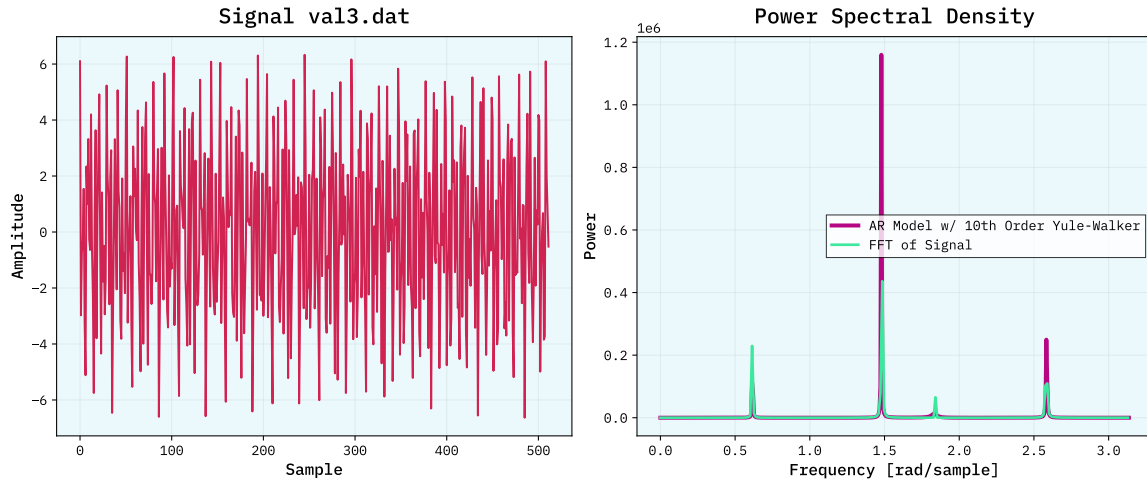


Figure 6: The spectra signal val3.dat.

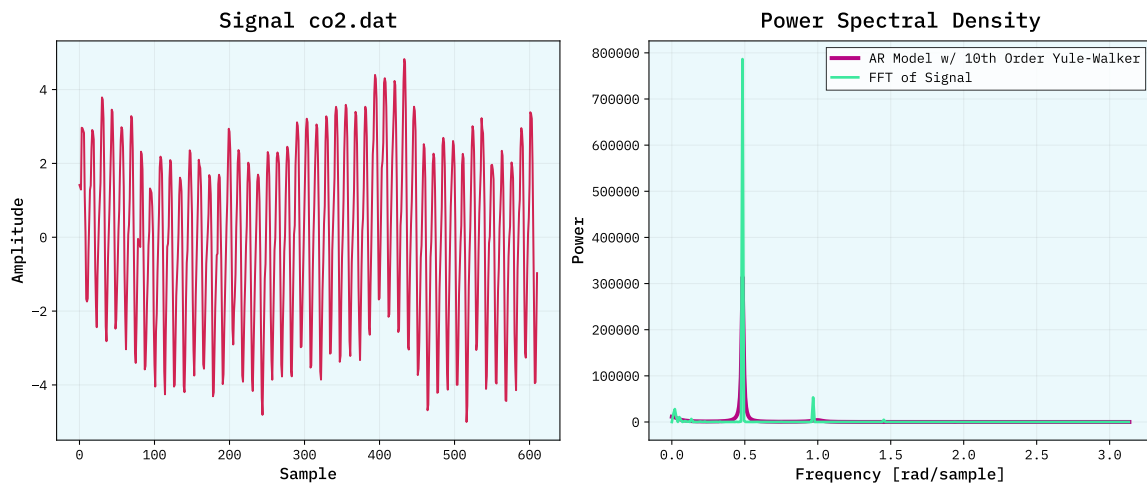


Figure 7: The spectra signal co2.dat.

As stated above let's have a look at the width of peaks in the spectra. I evaluated the widths with the help of `scipy.signal.find_peaks` and `scipy.signal.peak_widths` on signals `val2.dat` and `val3.dat`. The results are plotted in Figures 8 and 9. Plotted alongside are also the roots of the characteristic polynomial for each model.

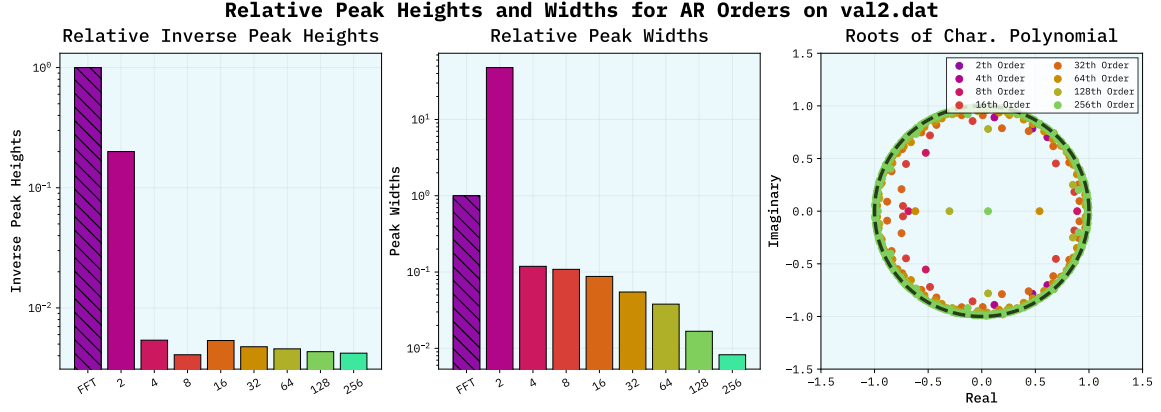


Figure 8: The peak widths/heights of the spectra of signal `val2.dat`.

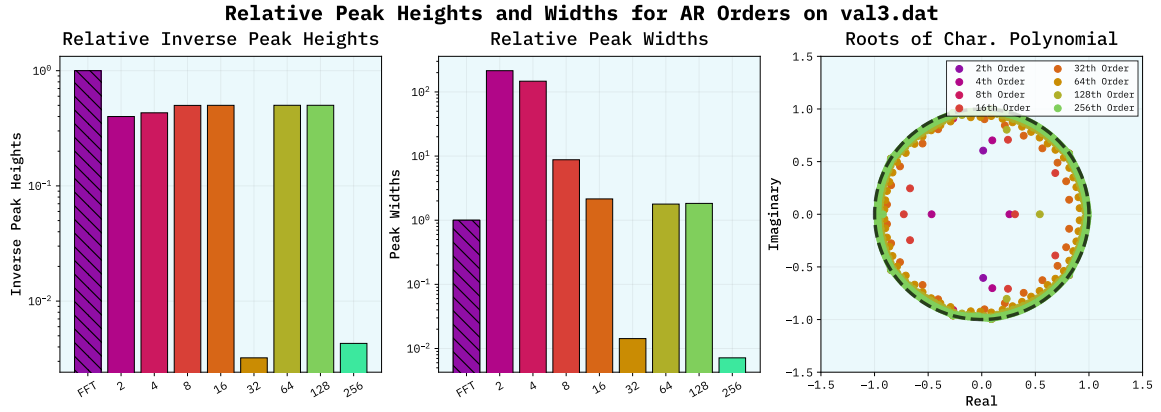


Figure 9: The peak widths/heights of the spectra of signal `val3.dat`.

These plots are a bit more information dense so it makes sense to give some explanation. The left-most plot shows the relative inverse peak height as compared to the FFT spectra. The inverse height meaning that the values plotted present $\frac{1}{\text{height}}$. This was done due to a massive range in the heights of the peaks which generally depend on the order of the filter as we can also see from the plot. There seems to be a relatively steady downwards trend in the inverse peak height, which means that they are growing. A similar trend can be seen in the middle subplot which represents the relative peak widths. Do note, that there is no inverse in play here. We can see that the widths of the peaks narrow as the order of the AR model rises. For low values of the AR model order the peaks are much wider than the FFT peaks. In fact they more resemble bumps rather than sharp peaks. The right-most plot shows the roots of the characteristic polynomial of the AR model. The main focus here was to check that all of the roots are inside the unit circle which is also plotted. The roots seem to be stable. I find the way the roots spread around in the complex plane incredibly fascinating.

For a better comparison between the AR model and the FFT spectra I plotted all the spectra on the same plot. Do note here that the AR and FFT spectra have been normalized such that we look at relative differences in scale, not absolute. The result is presented in Figure 10.

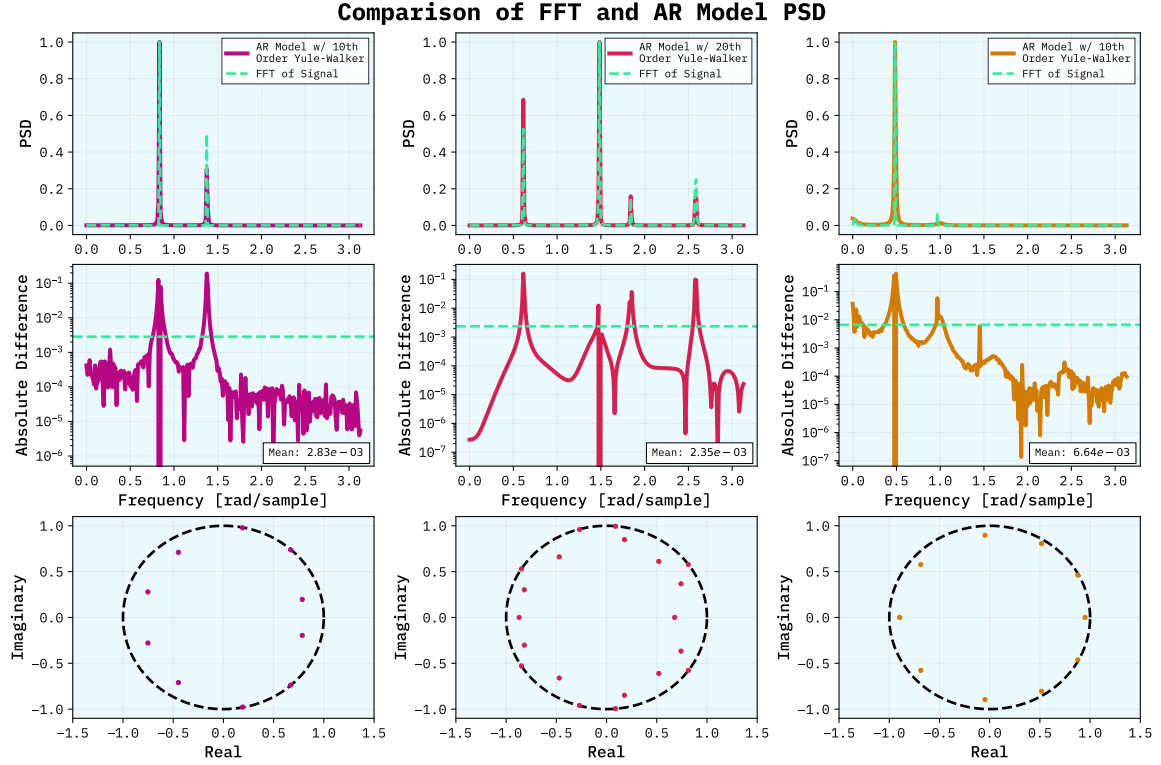


Figure 10: All the spectra of the signals on the same plot.

Again for the sense of safety and to check our methods I've also included the roots of the polynomials on the same plot. I'd like the reader to note that the AR models are not of the same order. The AR model order was chosen based on what I felt empirically gave the best results. We see that the spectra line up very nicely. There are some differences but we can see that for all the provided signals the mean absolute error is in the order of 10^{-3} .

Lastly let's have a look at the effects filter order and number of evaluation points (evaluation density) have on the spectra. This time I've also included `co2.dat` in the analysis. The results are plotted in Figures 11, 12 and 13.

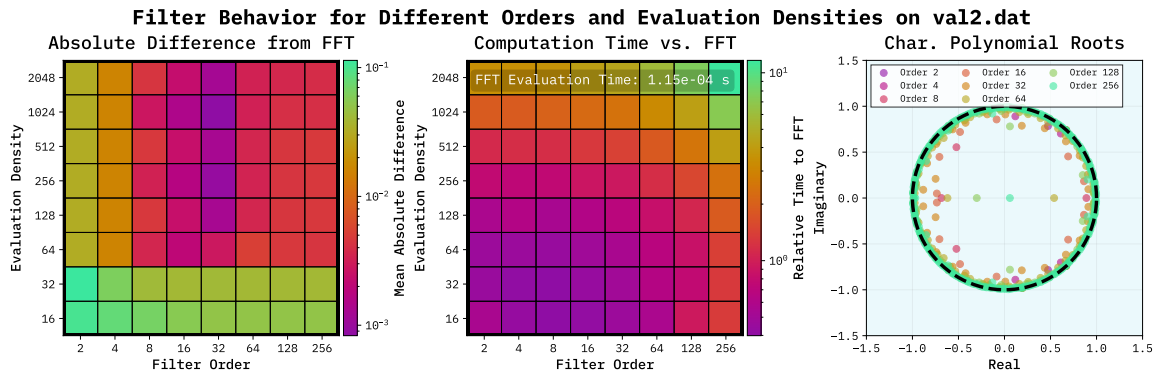


Figure 11: The effects of filter order and evaluation density on the spectra of signal `val2.dat`.

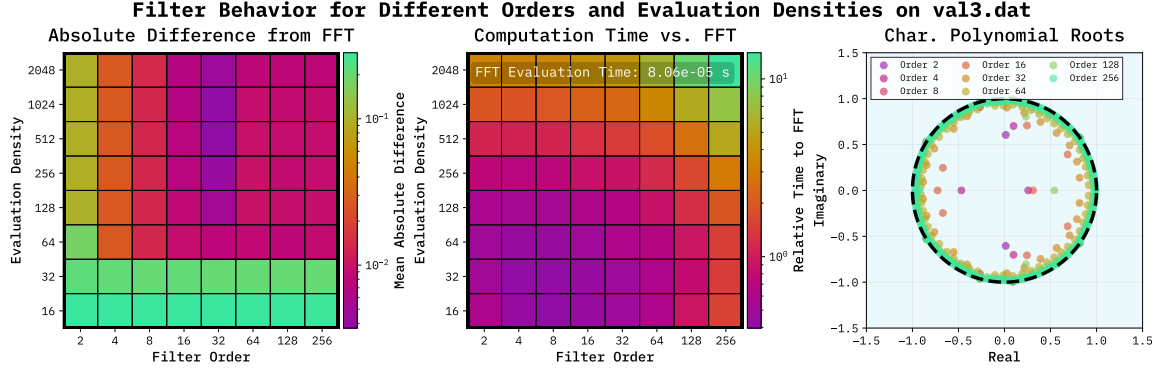


Figure 12: The effects of filter order and evaluation density on the spectra of signal `val3.dat`.

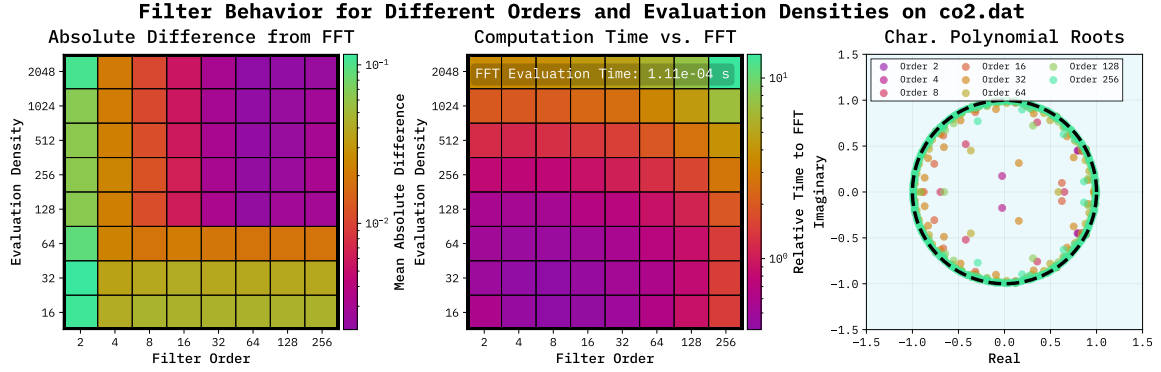


Figure 13: The effects of filter order and evaluation density on the spectra of signal `co2.dat`.

The bright colors in combination with the bolded grid really remind me of a 80s dance floor, I hope the reader enjoys the visual experience. The plots show that both the filter order and evaluation density have a significant impact on the spectra. One would naively expect that the absolute difference from the FFT spectra would decrease with increasing filter order, however that is not the case. In the plot for `val2.dat` and `val3` we see that the absolute difference is lowest for the AR model of order 32. The evaluation density has a lesser effect on the spectra, only coming into play when the density is very low and the peaks are not well resolved. In addition to absolute differences I've also plotted the computational times in comparison to the FFT spectra, where we can see that the AR model is comparable in speed to the FFT algorithm at the filter orders that yielded the best results. The right-most plot is there again for a sanity check to confirm that all of the roots are inside the unit circle.

It would be interesting to what the resolving power of an AR model is. What I mean by this is how close can to peaks be in the spectra for them to still be resolved. I did a little parameter scan for a couple of different values of filter order, peak separation distance and spectra evaluation density. The results of the scan are plotted in Figure 14. The FFT spectra from the scan test signals are plotted in the Figure 15.

Resolution Power of AR vs. FFT

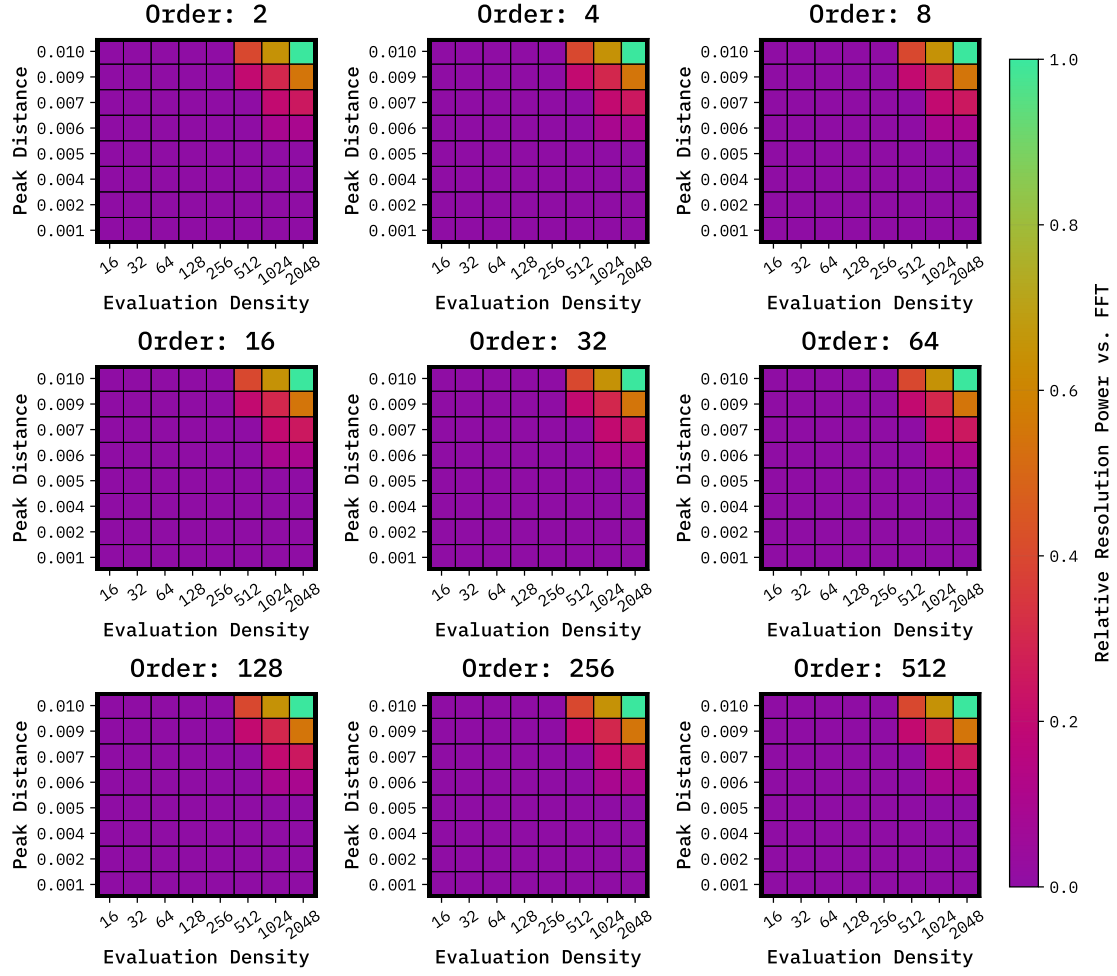


Figure 14: The 'resolving power' of the AR model. Unresolved is colored in purple.

Spectra of Test Signals with Different Peak Distances

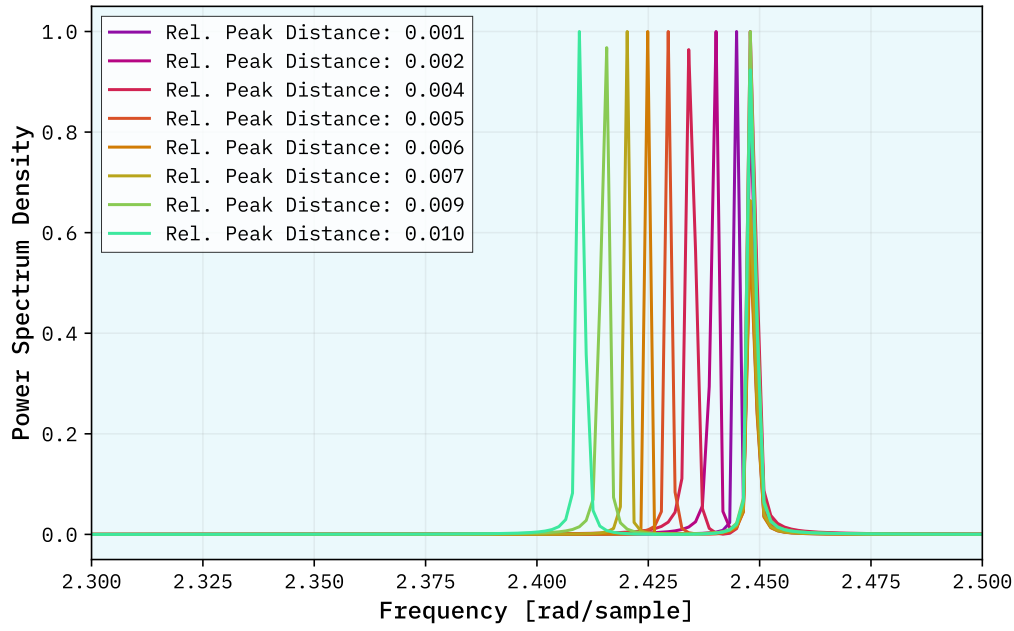


Figure 15: The test spectra by FFT from the resolving power scan.

All resolving powers are compared to how far apart the peaks are in the FFT spectra. This plot represents one of my major disappointments during this task in that, there is no dependence on the filter order, at least not in the range that I scanned. The results are the same for all filter orders. We can however see that in the range that I've provided only a small section of the parameter space contains peaks that actually managed to resolve. The majority of the peaks were not resolved to my great disappointment. I inspected the spectra for any possible errors but they truly do seem to not resolve. The subset of the resultant spectra are plotted in Figure 16.

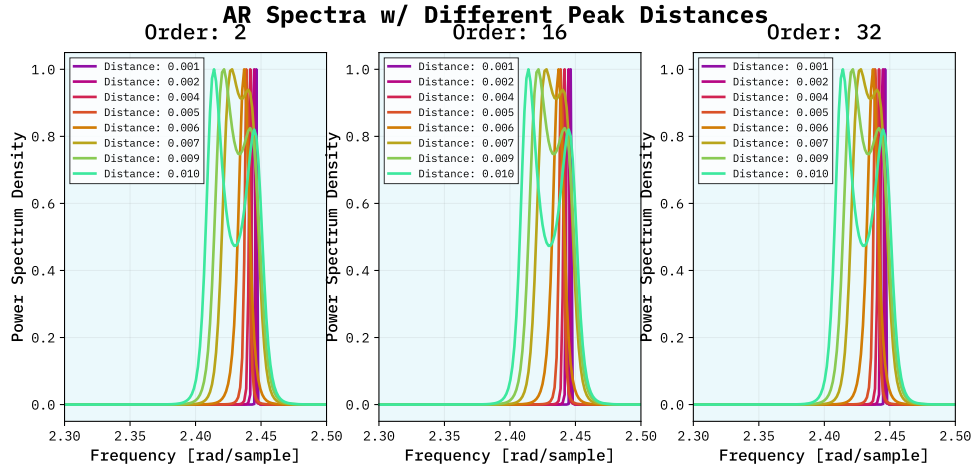


Figure 16: The subset of the spectra, where we can see most of the peaks are unresolved.

5 Conclusion and Comments