Bachelor Project in Physics

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todo: Make cover page todo: Add abstract

1 Note about the project

In the beginning of February 2022, Troels Petersen (KU) and Lars Buchhave (DTU) started working together to improve computational LFC calibration methods, and Troels invited me and a few other students to come along for the ride. During the project I've been exploring real data and developing a crude method for performing LFC calibrations and subsequent radial velocity extractions. With this project report it is my intent to share and explain the process of performing such analyses, thinking of the reader as a fellow physics student perhaps interested in picking up where I left off.

2 Introduction

The radial velocity method was used to discover the first exoplanets and continues to be one of the main methods for the discovery and characterization exoplanets [5]. With new extreme-precision radial velocity (EPRV) spectrographs such as the EXtreme PREcision Spectrograph (EXPRES), data from which this project is based on, we are slowly approaching the precision necessary for the discovery of Earth-sized planets around Sun-like stars. For this to succeed, it is however necessary to understand and mitigate many effects, such as the movement of the Earth relative to the center of mass of the solar system (barycentric corrections), light scattering in the Earth's atmosphere (tellurics) and light scattering inside the spectrograph (blaze). A general wavelength calibration of the spectrograph is also needed, the quality of which of course directly influences the precision of the final radial velocities that can be obtained. For that purpose, EXPRES utilizes a Laser Frequency Comb (LFC), is a rather new technique. While to extract radial velocities (RV) we measure the apparent wavelength shift between observations over a period of time, preferably more than a year, and utilize the well-known doppler effect.

The full procedure from raw data to results, also referred to as the *pipeline* in the literature, is extensive and complex, which is why I've ignored many aspects of it during this project. I've worked directly on the LFC calibration and RV extractions and tried to orient myself about the most important corrections, thus that is what I will describe in this report. The full pipeline is described in detail in [6].

3 Theory

3.1 Radial velocity method for exoplanet detection

The radial velocity method is one of the few current methods of detecting exoplanets. Two celestial bodies in orbit around each other, such as a star and a planet, orbit their common center of mass (barycenter). This means that the star, although typically much more massive than the planet, is also in movement relative to an outside observer. The larger the planet is, compared to the star, the larger this movement will be. For now, it is not possible to directly observe exoplanets. One way of indirect detection however is thus measuring the movement of the star. We can measure the relative movement of a star through the doppler effect; the electromagnetic spectrum of the star observed on Earth will be blue shifted when the star is moving toward us and red shifted when moving away. The potential indirect signal of a planet will thus consist of a periodic doppler shift in the spectrum of the host star. For this method in general, several years of data collection is necessary. One should at least have one full orbit of observational data, and, to decrease statistical uncertainties, several orbits is preferential.

A large planet like Jupiter induces a radial velocity (RV) in the Sun of about 12.7 m/s when observed in its plane of orbit. While a small one like Earth only induces an RV of about 9 cm/s. (p. 29, [5]).

todo: Possibly describe the RV calculation in more detail and compute Earth K.

3.1.1 Doppler shift

The radial velocity method relies on the well-known Doppler effect. Ignoring terms of c^{-4} and higher, the general shift caused by a relative displacement between the source and an observer at zero gravitational potential is given by

$$\lambda = \lambda_0 \frac{1 + \frac{1}{c} \mathbf{k} \cdot \mathbf{v}}{1 - \frac{\Phi}{c^2} - \frac{v^2}{2c^2}},\tag{1}$$

which accounts for both special relativistic effects and gravitatonal doppler shift described by general relativity. Where λ is the observed wavelength, λ_0 is the emitted wavelength, Φ is the Newtonian gravitational potential at the source ($\Phi = GM/r$ at a distance r of a spherically symmetric mass M), \mathbf{k} is the unit vector pointing from the observer to the source, \mathbf{v} is the velocity of the source relative to the observer and c is the speed of light [4].

Special relativistic effects we can safely ignore, as we are dealing with velocity shifts on the order of meters or centimeters per second, and thereby cross out the third term in the denominator. The remaining terms $(1 - \Phi/c^2)$ evaluted for HD 34411 $(M = (1.08 \pm 0.14) M_{\odot}, R = (1.28 \pm 0.04) R_{\odot}$ [2]) is around 0.999998. Adding this term to the computation of RV does not change my results, so we can neglect the denominator completely. todo: How to check analytically though?

If the unit vector \mathbf{k} were pointing directly toward us, it would mean that we were observing the system in the plane of orbit. This is however unlikely. Since we don't know the inclication angle, a possible simplification is to omit \mathbf{k} is treat the resulting \mathbf{v} as a minimum radial velocity.

Thus we are left with

$$\lambda = \lambda_0 \times \left(1 + \frac{v}{c}\right),\tag{2}$$

which is to say that the observed wavelength is simply the emitted wavelength scaled by a factor (1 + v/c). This formula allows us to compute the minimum relative velocity shift, v, between two observations, λ and λ_0 .

3.2 Description of the instrument

The EXtreme PREcision Spectrograph or EXPRES is an extreme-precision spectrograph situated at the Lowell Observatory's 4.3m Lowell Discovery Telescope (LDT) near Flagstaff, Arizona, USA. The LDT allows for up to 280 partial nights of observation per year.

Like in many spectrographs, at the heart of EXPRES is a Charge Coupled Device (CCD). A CCD is a silicon-based multi-channel photon detector consisting of a large number of small light-sensitive areas called pixels. The CCD is EXPRES an STA1600LN CCD backside illuminated image sensor with a 10,560×10,560 array containing 9µm×9µm pixels, designed to with a wavelength range of 3800–7800Å. When a photon hits a pixel it is converted into a charge, and each pixel can thus supply independent measurements. Since a one dimensional sensor would be impractical, EXPRES is constructed in such a way, that it wrap the spectrum inside the CDD, meaning that the spectrum starts in the top row of the sensor, and continues in the second row. Short wavelengths are thus to be found in the top of the CCD and long wavelengths at the bottom. EXPRES is housed in a vacuum enclosure to minimize changes in temperature and pressure, which can otherwise cause the spectra to change position on the CCD and thus lead to errors in the RV mesaurements.

Calibration device: Wavelength calibrations are performed with the use of a Laser Frequency Comb (LFC), produced by Menlo Systems, which is a laser source whose spectrum consists of a series of discrete, equally spaced frequency lines. The LFC however also needs calibration for which a Thorium Argon (ThAr) lamp with known frequencies is used.

Spectral resolution: EXPRES has a spectral resolution of R=150'000, where R is defined as $R=\lambda/\Delta\lambda$. Inverting this we get what's called the resolution element of the Instrumental Spread Function (ISF) (or Line Spread Function (LSF)), $\Delta\lambda=\lambda/R$. For a wavelength of say of $\lambda=5000\text{Å}$, this comes out to $\Delta\lambda=5000\text{Å}/150,000\approx0.033\text{Å}$, and it describes the "blurring" of monochromatic beams on the detector. Absorption features narrower than $\Delta\lambda$ can, in a well-behaved spectrograph, be approximated as a normalized, symmetric Gaussian function with FWHM = $\Delta\lambda$. Specifically for EXPRES, a super-gauss is however a better fit. LFC lines being monochromatic and thus very narrow will appear on the detector as a super-gaussian with $\Delta\lambda$ ranging from 3.9-5 pixels across the detector. By fitting, the center of the peak can be identified to a fraction of a pixel. The LFC peaks are seperated by about 10 pixels to remain distinct after this blurring. For star-spectra however some emission and absorption lines will be too close together and will appear "blended" on the detector.

Barycentric correction: Barycentric corrections are derived from the EXPRES exposuremeter, which is essentially a smaller, less precise spectrograph. Described in detail in [1]. EXPRES as a whole is described in technical detail in [3].

3.3 Description of the data

EXPRES data are meant to serve as an example of the data being produced by next-generation spectrographs.

The data used in this project was supplied by Lily Zhao and is by no means raw data, but data that has already gone through a lot of processing.

For development of RV extraction method, observations from four stars were used:

- HD 101501 (45 observations, 22 nights, Feb. 10, 2019 Nov. 26, 2020)
- HD 26965 (114 observations, 37 nights, Aug. 20, 2019 Nov. 27, 2020)
- HD 10700 (174 observations, 34 nights, Aug. 15, 2019 Nov. 27, 2020)
- HD 34411 (188 observations, 58 nights, Oct. 08, 2019 Nov. 27, 2020)

The observations for HD34411 are plotted in figure 1. Most days have 3-4 observations, and there are significant gaps in the data as well.

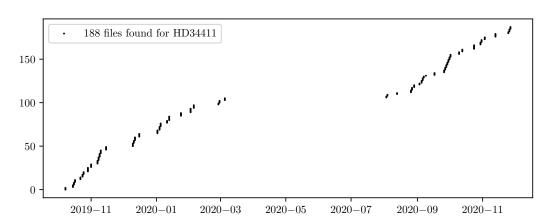


Figure 1: EXPRES observations of HD 34411

todo: Specify which data was used for LFC calibration method development $\,$

3.3.1 Data structure

The data used in this project consists of already packaged FITS (Flexible Image Transport System) files, which is a portable file standard widely used in the astronomy community to store images and tables. There is a FITS file for each observation, containing a variety of mesaurements for each pixel on the CCD. The rows of the CCD data are referred to as orders. There are 86 orders each of which has values from 7920 pixels. Drawing a coordinate system on the CCD, we are thus moving through pixels as we go along the x-axis and through orders as we go along the y-axis.

This would give the CCD the very elonganted dimensions of 86×7920 , but as mentioned earlier, the CCD is actually square. The orders however hit the CCD at an angle and for this reason *order tracing* is necessary. Order tracing reduces each order from 2d array to a 1d array, which means the final image comes out much shorter in the vertical/order dimension. Described in detail in section 3.2.1 of [6].

Furthermore, the CCD is not equally sensitive everywhere, and there are areas along the edges that are deemed useless. The data comes with a mask which shows which pixels are should be used.

To perform the calibration the following variables are used: spectrum, uncertainty, wavelength and continuum. To perform the RV extraction on the pre-calibrated data the following variables are used: bary_excalbur, excalibur_mask, spectrum, uncertainty and continuum.

3.3.2 Noise

Photon noise and read noise are the two largest contributors to the noise on a given pixel on the EXPRES CCD. These two quantities are measured and summed in quadrature for each pixel. Photon noise is assumed to be poisson distributed and the standard deviation is then the square root of photon counts. Read noise is calculated emperically, details of which will not be discussed, but is assumed to be consistent throughout each night of observation.

3.3.3 Correction: Scatter / blaze

Although manufactures have tried their best to limit it, the CCD still gets hits by scattering light, being the strongest in the center of the detector. It is modeled and subtracted by measureing the photon count in between orders.

3.3.4 Correction: Tellurics

Tellurics (general definition being originating from the Earth) in this context refers to the contamination that ground based spectrographs have to deal with, which occurs as the light passes through the Earth's atmosphere, encountering molecules such as oxygen and water vapor. On EXPRES the technique used is called SELENITE [https://arxiv.org/pdf/1903.08350.pdf].

3.3.5 Correction: Barycentric corrections

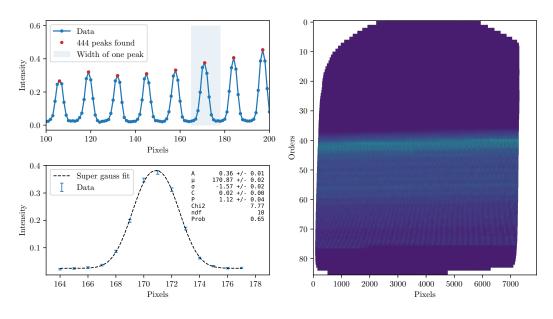
todo:

4 Data Analysis

4.1 Calibration

The calibration is needed to map each pixel on the CCD to a specific wavelength. Such a map is referred to as a wavelength solution. To do this, we need a light source with known frequencies, preferably many discrete peaks. EXPRES uses a Thorium Argon lamp for an initial trial wavelength solution and then a laser frequency comb (LFC) for an more precise solution.

The Thorium Argon lamp produces 4,000 lines across 82 orders, which can be identified and mapped to a wavelength through a *line atlas*. An intial wavelength solution for all pixels is then produced by linear interpolation. (In this project I have not done this calibration).



the CCD (unitless). Upper left: illustration of a few LFC peaks in order 65. Peaks are identified with scipy peak finder. Lower left: each peak is fitted with a super gauss to find the exact top of the peak with uncertainties.

Figure 2: Right: Measured

intensities for the LFC across

The LFC generates a series of equadistant (evenly spaced) spectral lines, typically 20,000 lines across 50 orders. The range of the LFC is thus shorter, and for this reason the ThAr exposures can also be used for a rough calibration outside the LFC range. The frequencies of the LFC peaks are given by the relation

$$v_n = v_{\text{rep}} \times n + v_{\text{offset}}$$
 (3)

for integers n. The repetition rate $v_{\rm rep}$ and offset frequency $v_{\rm offset}$ are referenced against a GPS-disciplined quartz oscillator, providing calibration stability corresponding to a fractional uncertainty of less than 8×10^{-12} for integration times greater than 1 s. (p. 8, [6]). The values I have used in the calibration, $v_{\rm rep}=14e9$ and $v_{\rm offset}=6.19e9$, were provided by Lars Buchhave, but may be outdated. See figure 2 right side for a plot of the intensities measured across the CCD.

The following procedure is followed to determine the location of the LFC peaks on the CCD: 1) Find peaks using scipy peak finding algorithm 2) make data slices around each peak with the size of the average distance between peaks, 3) using iminuit do a χ^2 minimisation fit to each peak with a super-gauss plus a linear background. See figure 2 left side.

A super-gauss, defined in eq. (4), is a regular gaussian but with an extra parameter, here denoted P, that allows the top of the gaussian to be flattened. The last two terms here add a linear background and an offset.

$$f(x; A, B, C, P, \mu, \sigma) = A \exp\left(-\left(\frac{(x-\mu)^2}{2\sigma^2}\right)^P\right) + B(x-\mu) + C$$
 (4)

The fit then is a minimisation of

$$\chi^2 = \sum_{i=1}^{N} \left[\frac{y_i - f(x; A, B, C, P, \mu, \sigma)}{\sigma_i} \right]^2$$
 (5)

Where N is the number of data points, x is pixel-space, y_i and σ_i is the measured photon count and uncertainty respectively. The fit returns the values and uncertainties for the parameters A, B, C, P, μ, σ when the value of χ^2 is minimized.

We are most interested in μ , which gives the position of the LFC peak on the CCD (in pixel-space). With the intial rough wavelength solution dervied from the ThAr lamp (pre-calculated in the data set that I've used) I can determine what the approximate wavelength of the LFC peak should be. To find the better wavelength solution I then go look up the closest frequency given by eq. 3. And we now have a map of 20,000 points on the CCD with a good wavelength solution.

Of course we need to have a wavelength solution for all points on the CCD and to do that I have explored two approaches: polynomial fitting and cubic interpolation.

4.1.1 Poly-fit calibration

Since the LFC peak positions, as seen in the right plot in figure 2, appear to exhibit a certain periodic behavour, my initial approach to compute a wavelength solution across the whole CCD was to fit the LFC peak positions with a polynomial. Looking at the residuals of fitting the LFC peaks with polynomial of increasing degree revealed smaller and smaller periodic variations, until reaching 5th degree, see figure 3.

But, as the orders are far from idential, this turns out not to work very well across the CCD, see figure 4. So I decided to turn to interpolation.

4.1.2 Interpolation calibration

A cubic interpolation would force all the residuals to be zero, so in order to evaluate the quality of the method, we can for instance omit every second peak from the interpolation and then compute the residual between the omitted peaks and the resulting interpolation function. We can additionally flip the data set and interpolate the peaks we left out before and compute residuals for the rest, thus ending up with an array of residuals equal to the length of the results from the poly-fit method, allowing us to compare the two, as is done in figure 5.

The RMS of the residuals from the interpolation come out much smaller than that of the polyfit (values specified in figure 5), in this example, suggesting that the interpolation method is superior. It is also worth noting that because the interpolation was done on only half the data points at a time, it will be even better when performed on all data points, as it would be, when used for calibrating data before an RV analysis.

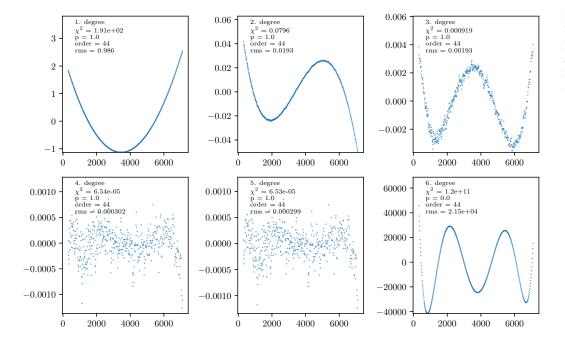


Figure 3: Residuals from fitting LFC peak positions with polynomials of increasing degree. Pixels on the x-axis and residuals in ångstrom on the yaxis

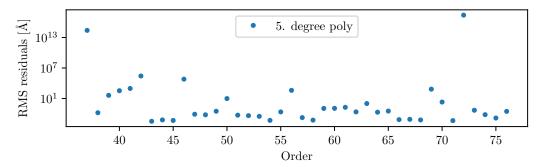


Figure 4: RMS of residuals from fitting LFC peak positions with 5. degree polynomial. Pixels on the x-axis and residuals in ångstrom on the y-axis

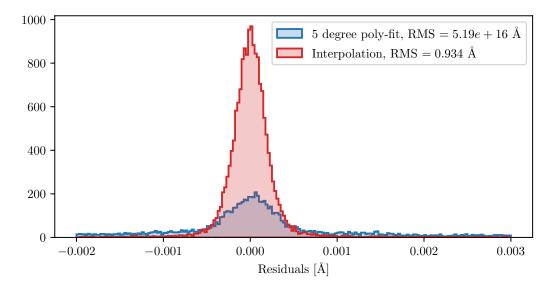


Figure 5: Residuals from calibrations performed through a 5th degree poly-fit and a cubic interpolation. Both results contain approximately the same amount of points.

todo: Can we convert this RMS to an RV error?

todo: add graph comparing residuals using gauss vs super gauss

todo: perhaps add plot of changes in parameters across the CCD

4.1.3 Errors in the calibration data

The LFC fits files come with an uncertainty on the spectrum-variable. It appears however that this uncertainty might be a bit underestimated. We can see this by plotting the χ^2 -

and P-values for the LFC peak super-gauss fits, as done in figure 6. The χ^2 value should be roughly equal to the number of degrees of freedom in the fit (eq. 4), which is:

$$N_{\text{dof}} = N_{\text{data-points}} - N_{\text{fit-parameters}} = 13 - 5 = 8.$$
 (6)

I set the rounded average distance between peaks in a given order to be the number of data points per fit. Although the LFC should generate equadistant peaks, this does vary between 13 and 18 points as we go through orders, most fits having 13 data points (see figure 13 in appendix A). We should therefore see a spike in the χ^2 distribution roughly around 8, falling of at a slower pace to the right side. Looking again at figure 6. This is not the case for the errors as provided (scale-factor 1, red). The distribution is very flat but peaks somewhere around 25, which suggests that the errors are about a factor $\sqrt{3}$ too small (square-root because χ^2 of course is squared and $25/3 \sim 8$). Another check to make is whether the p-value distribution is approximiately flat. For a

is whether the p-value distribution is approximiately flat. For a Also in figure 6 I've plotted the resulting χ^2 distributions if we scale up the errors, here by $\sqrt{3}$ (green) or $\sqrt{10}(blue)$. Looking at several other factors in between (see figure 14) I would argue that scaling the errors by $\sqrt{3}$ is appropriate as this gives a peak in the χ^2 distribution at 7.5 and furthermore

 $\sqrt{10}$ is overdoing it, however. I looked at several other scaling factors between $\sqrt{3}$ and $\sqrt{10}$, see figure 14 in appendix A for details, and $\sqrt{3}$ comes closest to giving a peak at 8.

Another check is that the probability distribution should be roughly flat. todo: explain why, help.

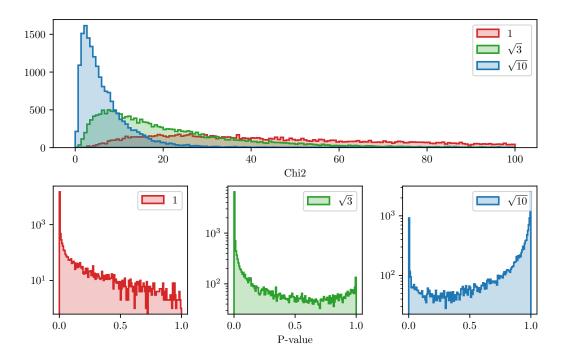


Figure 6: Chi2-values and P-values from individual LFC peak super-gauss fits with photon count (spectrum) errors multiplied by different scale-factors $(1, \sqrt{3} \text{ and } \sqrt{10})$. See text for more details.

Effects on calibration:

The errors used during the production of the calibration residuals shown in figure 5 I have already multiplied by $\sqrt{3}$. This gave a $\sigma = 0.934$, without this correction I got $\sigma = 2.34$.

todo: Computing the sigma for different error scaling factors gave very jumpy results.. why?

4.2 RV extraction

In signal processing, cross-correlation is a measure of similarity of two series as a function of the displacement of one relative to the other.

To extract radial velocity we need to measure the doppler shift between spectra from different days of observation. The most straight-forward way to do this is to compute the cross-correlation, since, in signal processing in general, the cross-correlation is exactly a measure of the similarity of two data series as a function of the displacement of one relative to the other.

Due to a lack of access to data consisting of star spectra with associated LFC captures, I've worked on RV extractions using already calibrated data provided by Lily Zhao. This data has been calibrated using a technique called excalibur [7].

Top left of figure 7 shows an extract of wavelength vs. intensity data from an observation of HD 34411. The file also includes a model of the continuum function, with which we can normalize the spectrum through division, shown bottom left. On the right side is plotted all normalized data within the EXCALIBUR mask, i.e. data marked as having a proper calibration.

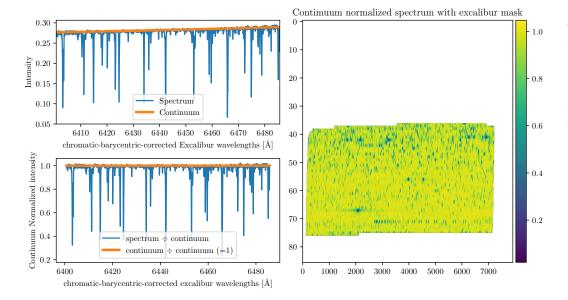


Figure 7: Overview of excalibur calibrated data from an observation of HD 34411. Upper left: extract of wavelength solution vs. intensity. Lower left: continuum normalized spectrum. Right: all continuum normalized data within the excalibur mask.

4.2.1 Finding and matching features across observations

In order to measure how much individual peaks move in between observations the first challenge is to find the "same" peak in both observations. The following procedure is followed:

First I invert the spectrum so absorption lines becomes peaks, and then I use scipy peak finder to find individual peak locations.

- Load intensity/spectrum from spectrum and err from uncertainty as well as excalibur calibrated barycentric-corrected wavelengths from bary_excalibur, all masked by excalibur_mask.
- Normalize spectrum and error with continuum from continuum.
- Then invert spectrum to turn absorption lines into peak by y = 1 y.
- Locate peaks find_peaks from scipy.signal

And to match features/peaks between two observations:

- Iterate through the peaks of file1 and find the closest peak in file2. With excalibur calibrated data, peaks should not shift so much that they overlap. However we can add at least two filters to bypass bad matches:
- Maximum peak distance: 0.5 Å is a very generous cut, equivalant to about 20-30 km/s depending on the wavelength. Since the peak location we are using here however is not the actual peak but the neartest data-point, setting the cut smaller, cuts away many good matches.
- Maximum difference between the area under the graph of two peaks: i.e. the sum of the intensity values. Peaks with similar shapes will give a low difference. Some bad matches can be avoided with this filter, but it also excludes a lot of good matches. Ideally I would keep the filter as strict as possible while maintaining a generous amount

¹To increase computation run time, I'm not sad about cutting away some good matches though.

of matches. In figure 8 is plotted the number of matches found between all files as a function of the maximum area difference. From that I conclude that 0.1 or 0.2 are good options, as the number of matches doesn't rise much with a looser restriction.

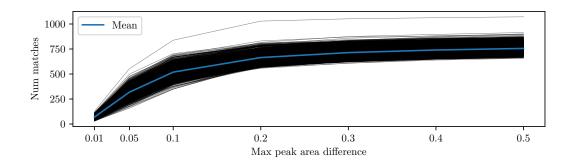


Figure 8: Number of matches found as a function of max peak area difference for all features in all files for HD 34411.

todo: Add feature feature shift plot (median, mean??)

todo: Add examples of bad matches

4.2.2 Cross correlation

To perform a smooth minimization of a shift parameter, it is however necessary to have continuous data, so the first step is to perform a cubic interpolation of the spectra data. In practice I do this inside my chi2 minimization function. Taking the spectra data and wavelength solution from two different observations, but multiplying one of the wavelength with the doppler shift factor from eq 2 before performing the interpolation. The interpolation functions are then evaluated in the range of wavelength solutions common to both observations, using N=1000 steps.²

$$\chi^2 = \sum_{i=1}^{N} \left[\frac{y_i - f(x; A)}{\sigma_i} \right]^2 \tag{7}$$

where y_i are the unshifted interpolated photon counts for one file and the function f(x; A) is an cubic interpolation function given my the intensity values of the second observation with wavelength values shifted by equation 2, then evaluated on the common wavelength range:

$$f(x; A) = \mathbf{interp}[x \times (1 + A/c), y] (x_{\text{common}})$$
(8)

The errors, σ_i , are also computed through interpolation.³

I then compute the cross-correlation and obtain the radial velocity, A, as a minimization of eq (7) using iminuit.

When it comes to selecting data I've tried three approaches: order by order, chunk by chunk (chunks of 2Å, inspired by Lily Zhao's approach [6]) and feature by feature. So far, I've only managed to get good results fitting feature by feature, which I will go through in more detail below.

4.2.3 Extracting relative shifts from over constrained system - matrix reduction:

The end desired end result is a plot of the relative radial velocities as a function of time or observation. If we were to simply compare each observation with the following one, our results would become correlated in the sense that the difference between day 1 and day 10, let's say, will depend on the quality of all observations in between. To circumvent this, we can compute the relative shift between all observations, yielding an $N \times N$ upper triangular

²1000 steps was chosen as the best balance between run time and resulting uncertainty. See figure 15 in appendix B. todo: Are errors correct?

 $^{^3}$ The errors, σ_i , which of course also need to be continuous, are computed by: 1) computing another two cubic interpolations for the non-shifted observation, one with photon counts plus photon count errors and another one with photon counts minus photon count errors. Both interpolations are then evaluted on the same common wavelength range with 1000 steps. The error for each data point is then computed as the mean difference between the measured value and the upper and lower error.

matrix, where each cell is the shift between observations i and j, and thus with a diagonal of 0. I will call this matrix ΔV_r^{ij} , see figure 9, here non-barycentric-corrected data is used for the sake of illustration.

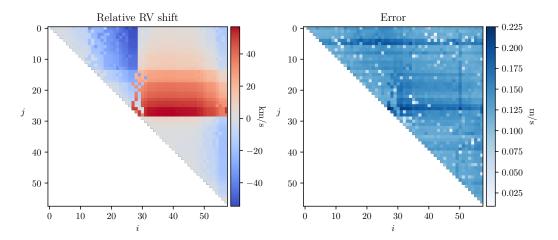


Figure 9: RV shifts matrix computed for HD34411 using non-calibrated, barycentric-corrected data (column wavelength). Each cell shows the median relative radial velocity all features found between observations i and j. diagonal should be zero and has been omitted for the sake of computational speed. I only used one out of the several exposures from each day.

The following chi2 minimization with the fit parameters V_r^i (an array of length N) will then find a list of values that best describe the whole matrix. In other words, it finds the relative shift for each observation with all other observations.

$$\chi^{2} = \sum_{i,j=0}^{N} \left[\frac{\Delta V_{r}^{ij} - (V_{r}^{i} - V_{r}^{j})}{\sigma(\Delta V_{r}^{ij})} \right]^{2} : i < j.$$
 (9)

The resulting fit parameters are plotted in figure 10, and we see a clear signal of the Earth movement around the barycenter of the solar system. Considering the resulting chi2- and p-value, it is possible that my errors are much too small. But I do get a wavelength/period close to that of one earth year.

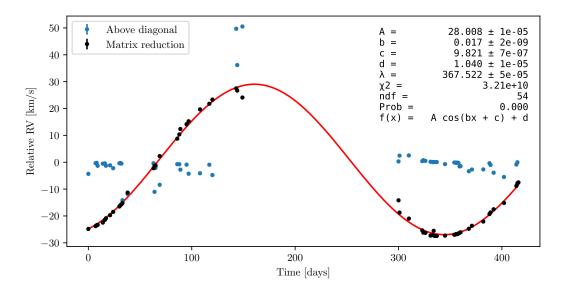


Figure 10: Computed relative wavelength shifts for HD34411 using non-calibrated, non-barycentric-corrected data (column wavelength). Blue: the shifts from day to day. Black: values computed through the chi2 minimization.

todo: add run times

5 Results

In figure 11 is plotted the ΔV_r^{ij} matrix for barycentric-corrected excalibur calibrated data for HD 34411. We are now on the order of m/s instead of km/s.

Performing the matrix reduction chi2 fit yields the results plotted in figure

todo: Add plot with all files todo: Add plots for the rest of the stars

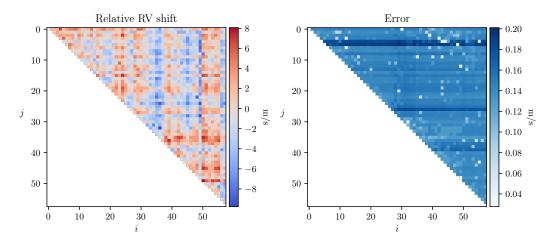


Figure 11: RV shifts matrix computed for HD34411 using excalibur calibrated, barycentric-corrected data (column bary_excalibur). Each cell shows the median relative radial velocity across all features found between observations i and j. The diagonal should be zero and has been ommited for the sake of computational speed.

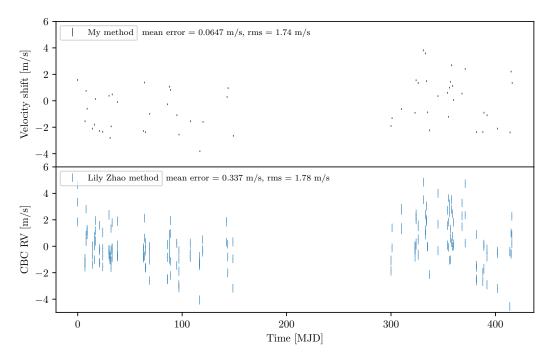


Figure 12: Computed relative wavelength shifts for HD34411 using excalibur calibrated, barycentriccorrected data(column bary_excalibur). Top: my results todo: Errors are too small. Bottom: Lily Zhao's results using a chunk-by-chunk analysis. todo: Cite?

6 Discussion

Future ideas:

- Auto encoder
- Lunch LFC into orbit around earth to have a base truth
- Better understand the stellar activities

7 Conclusion

8 Acknowledgements

References

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A LFC photon count errors

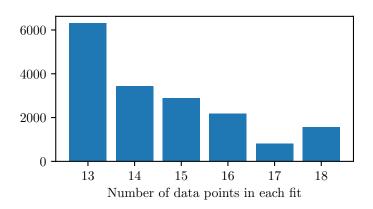


Figure 13: Number of data points in each LFC peak fit, determined by the average distance between peaks in each order.

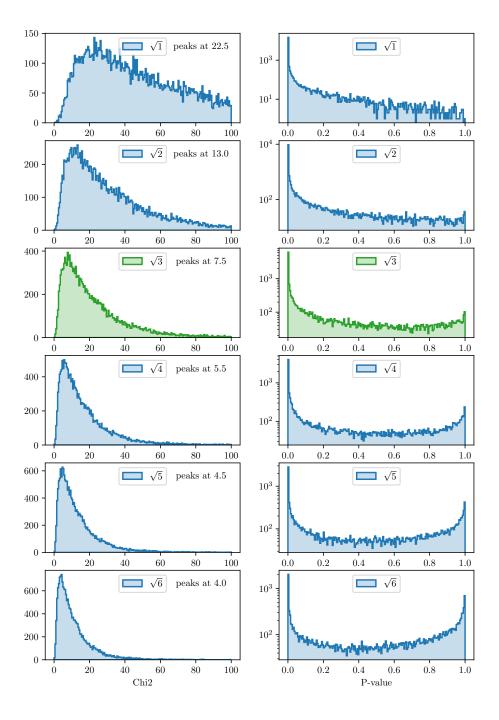


Figure 14: Chi2-values and P-values from individual LFC peak super-gauss fits with photon count (spectrum) errors multiplied by different scale-factors.

B RV extractions

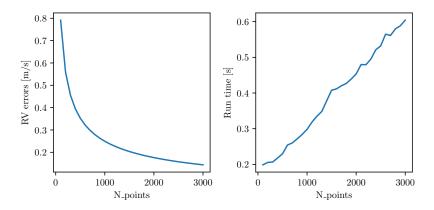


Figure 15: Number of data points in each LFC peak fit, determined by the average distance between peaks in each order.