

Bachelor Project in Physics

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todo: Make cover page

todo: Add abstract

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.

1 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].

2 Theory

2.1 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.

2.1.1 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

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

which accounts for both special relativistic effects and gravitational 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 ($1 - \Phi/c^2$) evaluated 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 inclination angle, a possible simplification is to omit \mathbf{k} and treat the resulting \mathbf{v} as a minimum radial velocity.

Thus we are left with

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

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 .

2.2 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 $10,560 \times 10,560$ array containing $9\mu\text{m} \times 9\mu\text{m}$ 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 measurements.

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 $\Delta\lambda = 5000\text{Å}/150,000 \approx 0.033\text{Å}$, 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 lines are separated 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 exposure-meter, which is essentially a smaller, less precise spectrograph. Described in detail in [1]. EXPRES as a whole is described in technical detail in [3].

2.3 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.

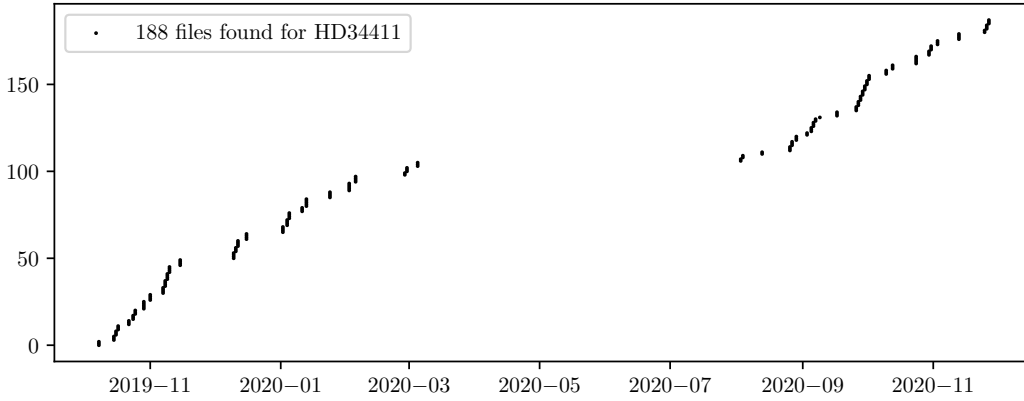


Figure 1: EXPRES observations of HD 34411

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

2.3.1 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 measurements 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 elongated 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_excalibur`, `excalibur_mask`, `spectrum`, `uncertainty` and `continuum`.

2.3.2 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 empirically, details of which will not be discussed, but is assumed to be consistent throughout each night of observation.

2.3.3 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 measuring the photon count in between orders.

2.3.4 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>].

2.3.5 Correction: Barycentric corrections

todo:

3 Data Analysis

todo now: Write this: the data analysis of this project is to calibrate and to extract rv.

3.1 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 LFC generates a series of 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 lines are given by the relation

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

for integers n . The repetition rate v_{rep} and offset frequency v_{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 1s. (p. 8, [6]). The values I have used in the calibration, $v_{\text{rep}} = 14\text{e}9$ and $v_{\text{offset}} = 6.19\text{e}9$, were provided by Lars Buchhave. 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 lines on the CCD: 1) Find peaks using scipy peak finder, 2) make data slices around each peak with the size of the average distance between peaks, 3) using iminuit do a chi2 minimisation fit to each peak with a super-gauss plus a linear background. See figure 2 left side.

A super-gauss, defined in equation 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.

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

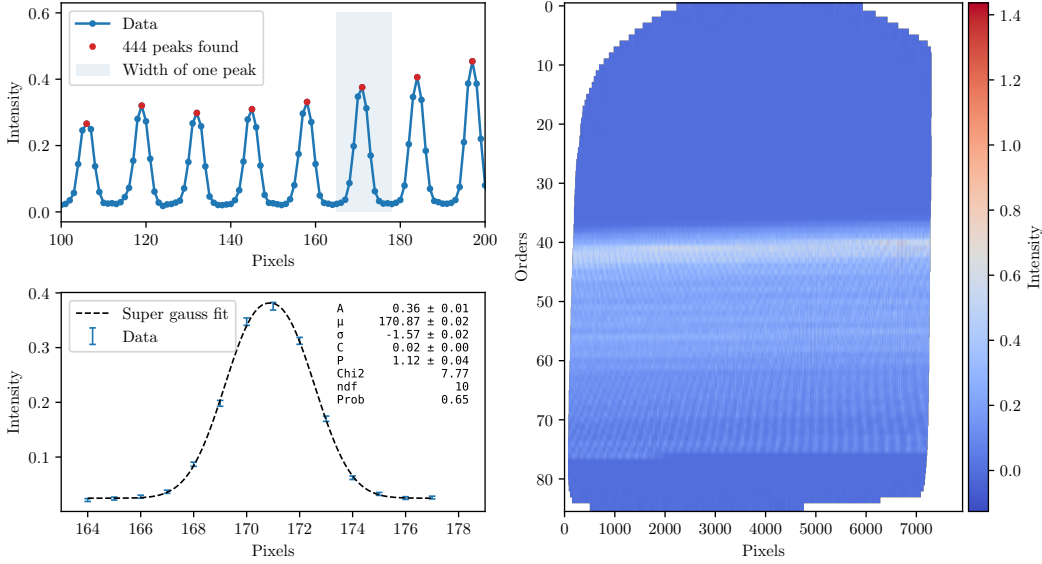


Figure 2: Right: Measured intensities for the LFC across the CCD (unitless). Upper left: illustration of a few LFC lines in order 65. Peaks are identified with scipy peak finder. Lower left: each peak is fitted with a super gauss to find the exact top of the peak with uncertainties.

The fit is then a standard chi2 minimization:

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

Where N is the number of data points, x is pixel-space, y_i and σ_i is the measured intensity and uncertainty respectively. The fit returns the values and uncertainties for the parameters A, B, C, P, μ, σ when the chi2 is minimized.

We are most interested in μ , which gives the position of the LFC peak on the CCD (in pixel-space). With the initial rough wavelength solution derived 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. We would however like to have a wavelength solution for all pixels on the CCD, so we need to find a way to estimate the wavelength solution in between LFC lines. To do that I have explored two approaches: polynomial fitting and cubic-spline interpolation. Before moving on though, it is worthwhile to study the chi2 values from the many fits we just performed, as they can help us evaluate whether the errors of the original data are correct.

3.1.1 Errors in the calibration data

According to Lily Zhao et al, the line spread function of EXPRES can best be represented by a super-gaussian [7], and so as we fit a super-gaussian to the peaks, the chi2 will grow if the measured values do not agree with the predicted values within the uncertainties. For a given data point, if the measured value differs by exactly the uncertainty, that data point would contribute 1 to the sum. However, fitting with the super-gaussian (including a linear background) we also have 6 fitting parameters that allow some wiggle room for the model to fit the data, which we must compensate for. For a given data series, such as an LFC peak, we can expect the chi2 roughly equal the number of data points in the sum *minus* the number of parameters in the fit, i.e. the number of the degrees of freedom:

$$N_{\text{dof}} = N_{\text{data-points}} - N_{\text{fit-parameters}} = 13 - 6 = 7, \quad (6)$$

as I use roughly 13 data points in each fit (the LFC line spacing does vary a bit across the CCD).

So if we plot all chi2 values we get from fitting the LFC lines, as I've done in figure 3, we should see a typical chi2 distribution with a peak around 7. Using the uncertainties as provided in the data file, the chi2 distribution is however very flat (red curve). It peaks somewhere around 25, which suggests that the uncertainties are about a factor $\sqrt{3}$ too small (square-root because the chi2 of course is squared and $25/3 \sim 8$), and scaling up the uncertainties by $\sqrt{3}$ and fitting the peaks again does in fact give a chi2-distribution with

a peak roughly around 8. additionally we can also look at the p-value distribution. With proper uncertainties, we should expect neither too many awful fits ($p=0$) nor too many perfect fits ($p=1$), but rather a roughly flat distribution. Although the chi2-distribution check is more clear-cut, the p-value distribution for uncertainties scaled by $\sqrt{3}$ (green) is better. Scaling by $\sqrt{6}$ is too much: the chi2 peaks around 4 and the p-distribution shows more perfect fits.

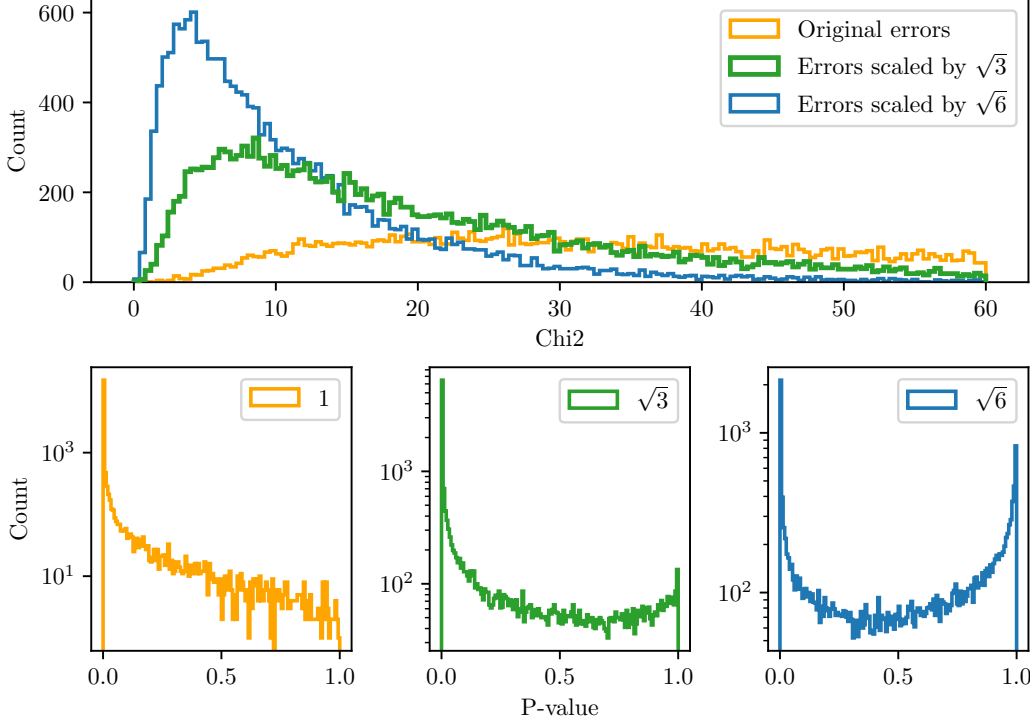


Figure 3: Chi2-values and p-values from individual LFC peak super-gauss fits with photon count (spectrum) errors multiplied by different scale-factors (1 , $\sqrt{3}$ and $\sqrt{10}$). See text for more details.

3.1.2 Poly-fit calibration

Since the LFC peak positions, as seen in the right plot in figure 2, appear to exhibit a certain periodic behaviour, 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 line positions with polynomials of increasing degree revealed smaller and smaller periodic variations, until reaching 5th degree, see figure 4. The p-value drops to 0 for 6th degree, and the per line rms (**todo now: 1**) explodes. Only LFC lines with a chi2 smaller than 100 was used for the analysis and errors were scaled by $\sqrt{3}$ locating LFC lines.

The residuals plotted show the difference between the theoretical line position given by equation 3 and the wavelength solution evaluated at the line position of

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$$\text{RMS/line} [\text{ms}^{-1}] = \sqrt{\sum_{n=1}^N \sum_{p=1}^P \frac{\left[\frac{(\lambda_{n,p, \text{pred.}} - \lambda_{p, \text{theory}})}{\lambda_{p, \text{theory}}} \times c \right]^2}{N \times P}}$$

- and refer to it from todo 1, and in the caption

The residuals of a wavelength solution represent the difference between the wavelength solution evaluated at the line position of a calibration line and the assigned theoretical wavelength (i.e., that from Equation (4) for the LFC lines) on a line-by-line basis in every exposure. The reported rms of a wavelength solution is therefore the per-line rms, i.e.,

3.1.3 Interpolation calibration

A cubic-spline interpolation would force all the residuals to be zero, so in order to evaluate the quality of the method, we can omit for instance every second peak from the interpolation and then compute the residual between the omitted peaks and the resulting interpolation

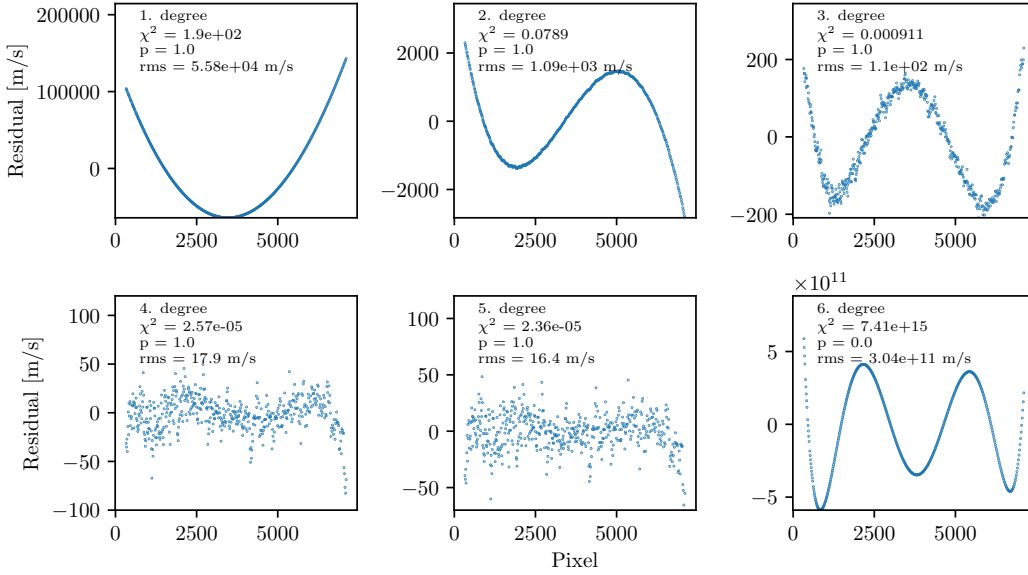


Figure 4: Residuals from fitting LFC peak positions with polynomials of increasing degree, in order 44. Pixels on the x-axis and residuals in m/s on the y-axis. Errors used for determining peak locations have been scaled by $\sqrt{3}$.

function. Then we can flip it around and interpolate the peaks we left out before and compute the remaining residuals.

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.

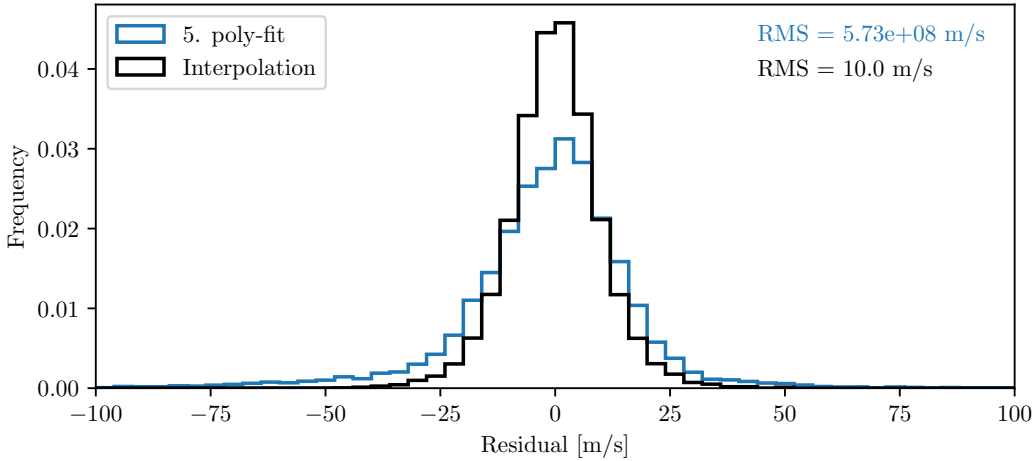


Figure 5: Residuals from calibrations performed through a 5th degree poly-fit and a cubic-spline interpolation. The per-line rms as defined in Equation (5) is given in the top-right corner in the color of the corresponding method.

3.2 RV extraction

To extract radial velocities we need to measure the doppler shift between spectra from different days of observation. One way to do that is to compute the cross-correlation, which is a measure of the similarity of two data series as a function of the displacement of one relative to the other.

We can do this either for individual absorption features, chunks of the spectrum a few angstroms wide or entire orders at a time. I've chosen primarily to work with the individual features.

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 [8].

The data is visualized in 6, where the top left 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 in the bottom left. On the right side is plotted all continuum normalized data within the EXCALIBUR mask, i.e. data marked as having a proper calibration.

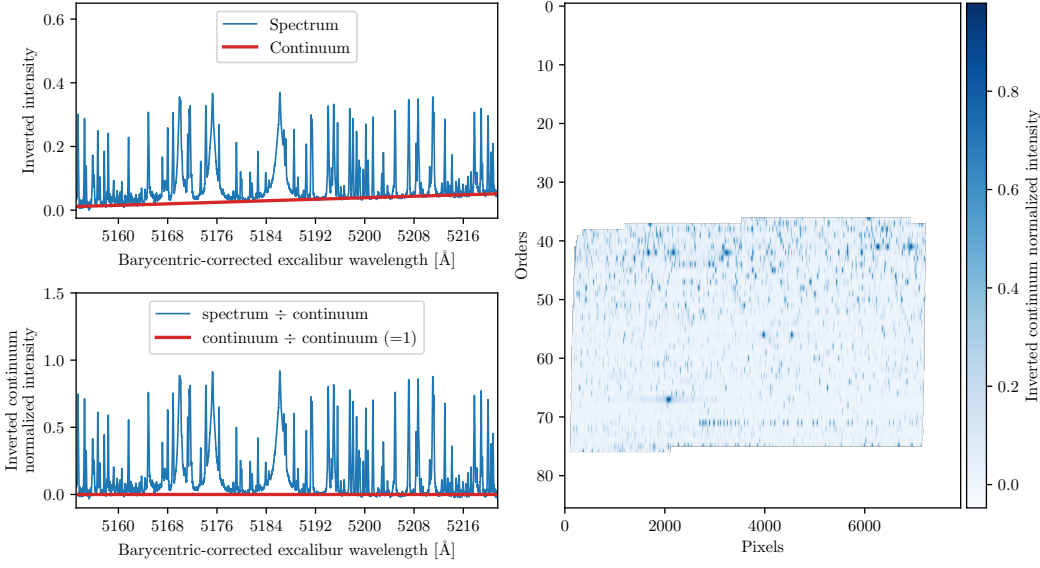


Figure 6: Overview of ex-calibur calibrated data from an observation of HD 34411. Upper left: extract of wavelength solution vs. intensity. Lower left: continuum normalized spectrum. Right: all continuum normalized data within the excalibur mask.

3.2.1 Finding and matching features across observations

In order to measure how much individual absorption features move in between observations the first challenge is to find the "same" features in both observations. What follows is a quick rundown of the procedure I've devised:

- Load intensities from the data column "spectrum" and errors from "uncertainty" as well as excalibur calibrated barycentric-corrected wavelengths from "bary_excalibur", all masked by "excalibur_mask".
- Normalize intensities and errors with the continuum function from "continuum".
- Invert intensities to turn absorption features into positive peaks by $y = 1 - y$.
- Locate peaks using `find_peaks` from `scipy.signal` with minimum peak distance of 5 pixels and a minimum peak prominence of 0.25 (unitless).
- Finally slice data around each peak with a width of 30 data points.

And then to match features/peaks between two observations:

- Iterate through the peaks of observations1 and find the closest peak in observation2. With excalibur calibrated data, peaks should not shift so much that they overlap. However the algorithm laid out so far does sometime match peaks that are far apart or do not resemble each other in shape at all. To bypass such matches we can add two filters:
 - Maximum peak distance: We could filter out all matches where the distance between the peaks is equivalent to a radial velocity greater than 12.5 m/s (the RV Jupiter induces in the Sun). However, since we are dealing with discrete data, the difference sometimes comes out much larger than it actually is and a narrow cut of 12.5 m/s would remove many good matches. Instead setting a very generous cut of 0.5 Å, equivalent to about 20-30 km/s depending on the wavelength, filters out the few very bad matches, but leaves the rest. When analyzing non-barycentric-corrected data, I set this up to 1 Å.
 - Maximum difference between the areas under the graph of two features (the sum of the intensity values in the feature): Peaks with similar shapes will give a low difference. This filter is most useful when analyzing non-barycentric-corrected,

where features often move so much that they overlap. In figure 7 is plotted an example of a bad and a good match, where the bad match can be avoided by setting the max area difference down to 0.1 (unitless).

- I've not yet devised any formal way to determine the best choice of these filters.

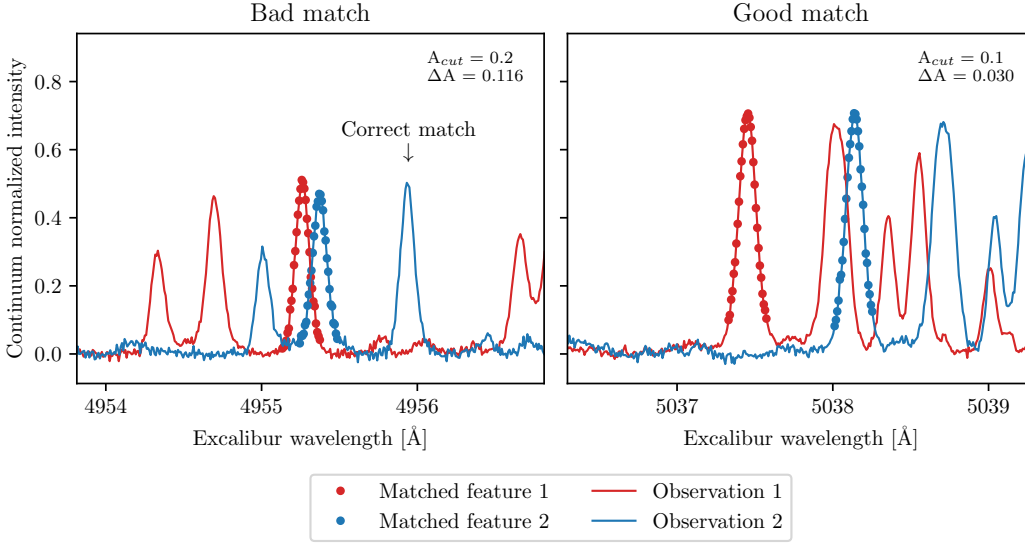


Figure 7: Example of match filtering with non-barycentric-corrected data, where features move a lot. The maximum area difference cut was set to 0.2 and 0.1 on the left and right plot respectively. The cut on the left plot was not strict enough to avoid the bad match that was made, with an area difference of 0.116. While on the right plot, the correct feature was selected spite another feature being closer in wavelength.

3.2.2 Computing velocity shift as the cross correlation

At this point, I have a list of matching features in different observations. The cross-correlations will be performed as a chi2 minimization fit to find the radial velocity that for each match correctly shifts one feature onto the other one.

Before moving on though, we have to cubic-spline interpolate the spectra data. I do this for two reasons: 1) the shifts we are looking for are much smaller than the individual pixels on the CCD, so we need to be able to shift by sub-pixel amounts, and 2) in order to compute the difference in intensity values between peaks, the intensity values must have the same wavelength solution, but, since EXPRES is calibrated independently for each observation, the wavelength solutions are different.

So I cubic-spline interpolate the spectra data from the first observation in the match, but before interpolating the second observation, I shift the wavelength solution by multiplying the shift factor from equation 2. Now I can evaluate the two interpolation functions on a common wavelength range, using $N = 1000$ steps¹, and I am ready to compute the chi2:

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

where y_i are the unshifted interpolated intensity from the first observation, σ_i are the errors on the intensity of the first observation also sampled through a cubic-spline interpolation, and the function $f(x; v)$ is the cubic-spline interpolation function given by interpolating the intensity values of the second observation with wavelength values shifted by equation 2, evaluated on the wavelength range common to both features:

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

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

¹1000 steps appeared as the best balance between run time and resulting uncertainty. See figure 12 in appendix A

3.2.3 Computing velocity shifts for all features

We can now find and compute the cross-correlation for all feature matches between two observations. In general I find between 500-1000 matches between two observations, which results in a fair amount of statistics. The computed radial velocities do build up a normal distribution around a central value, but there are many outliers. This is due to primarily two things: 1) Bad matches making it through the selection filters, and 2) stellar activity, which are various processes on the star that cause absorption features to shift independently.

This is one place where there is definite room for improvement. A better matching algorithm could supposedly be developed, but when it comes to sorting through lines affected by stellar activities, one could imagine that machine learning might be applicable. More on this in the discussion.

For now, I've experienced with settings different cuts both in radial velocity and chi2 to weed out the bad ones, but what I've found to work the best, is simply taking the median instead of the weighted average or the mean, as the median is much less affected by outliers than both. And as there is no reason to believe that stellar activity would cause more RV outliers in one particular direction, the spread can be assumed to be symmetric and thus the median is a good approximation. In figure 8 is plotted an example, which also contains at least one bad match on the far left and a constellation of features likely to be affected by stellar activity on the far right.

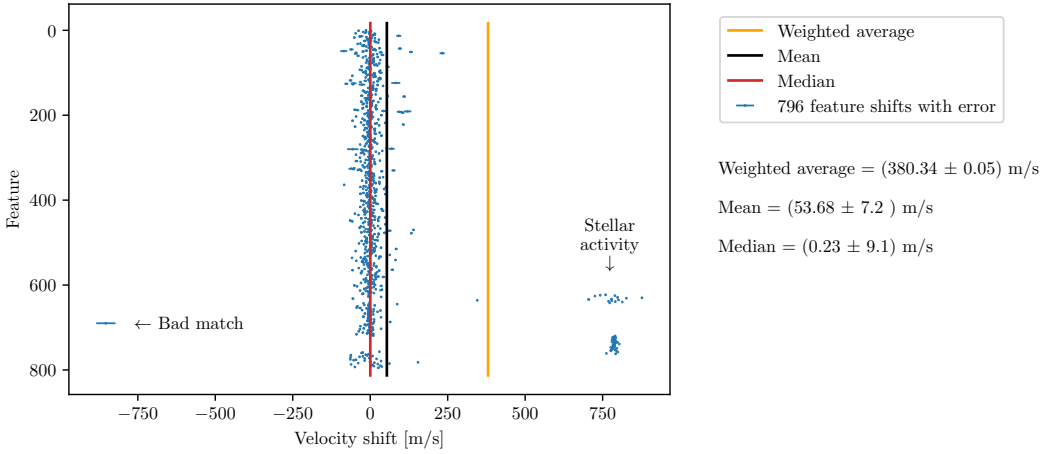


Figure 8: Comparison of weighted average, mean and median for computed rv shifts for two observations of HD 34411. The feature on the left side turns out to be a very broad feature, for which the standard data slice is too narrow and therefore yields bad results, (see figure 13 in appendix A) for details. While the features on the right side are likely to be caused by some kind of stellar activity.

The errors however is where my method is falting. Also plotted in figure 8 are the respective errors on the weighted average, the mean and the median. As it is the median that I've decided to work with, it would make sense to use the median error

$$\sigma_{\text{median}} = \frac{\sigma}{\sqrt{N-1}} \times \sqrt{\pi/2}$$

for large N where σ is the standard deviation, however, it not only comes out very large, but also breaks the propagation of the original errors from the spectrograph, as it essentially is a scaled standard deviation. For the sake of continuing with the analysis I go on with the weighted error, although that in contrast is very small. More on errors in the discussion.

3.2.4 Extracting relative shifts from over an overconstrained system

With the devised method we can now compute the relative radial velocity shift between two observations and get out one number with an uncertainty. The next most obvious step would be to compute the shift between observations 1 and 2, 2 and 3, etc. Doing this however leads to correlated results, as the difference between say observation 1 and 10 will depend on all the observations in between, and if there is one bad one, this will affect all the rest. To circumvent this, we can first compute the relative shift between all observations. This will give us an *overconstrained system*, in the sense that there is more information than necessary. All these differences can then be reduced down to a single array, where each shift is relative to all the rest, not only the neighbor.

Computing the shifts between all observations yields an $N \times N$ upper triangular matrix, where each cell is the shift between observations i and j , and thus with a diagonal of zero. I will call this matrix ΔV_r^{ij} , see figure 9 for an example.

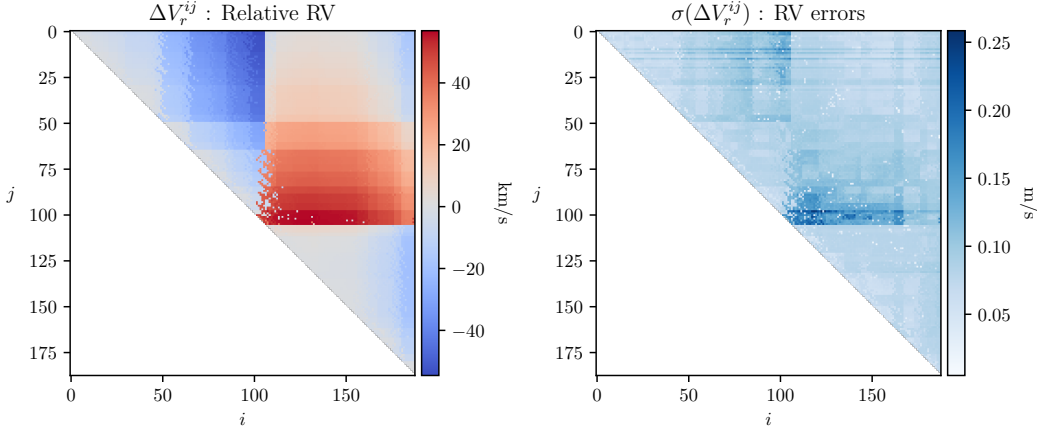


Figure 9: Radial velocity shifts matrix computed for 188 observations for HD34411 using excalibur calibrated but non-barycentric-corrected data (column `excalibur`). Each cell shows the median radial velocity shift for all features found between observations i and j .

To reduce the matrix to one array, we can perform another chi2 minimization fit, defined below in equation 9, in which we fit an array of parameters we can call V_r^i of length N (the number of observations), initialized to zero. The chi2 will be at its minimum when it has found an array of velocities V_r^i that best describe all the differences in the matrix ΔV_r^{ij} and each of the resulting velocities are thus relative not only to its neighbors but to all the other observations as well, thereby avoiding the correlation.

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

For the sake of illustrating the method, I've analysed non-barycentric-corrected data, that is to say, data in which we should be able to see the movement of the Earth around the center-of-mass of the solar system. That is the data plotted in the matrix in figure 9 and the resulting extracted relative radial velocities (the fit parameters V_r^i) are plotted in figure 10 (black), where we see a clear signal of Earth's movement. I've fitted the data with a periodic function and found a period of $(366.482 \pm 3e-6)$ days and an amplitude (the orbital speed of the Earth really) of $(28.132 \pm 9e-7)$ km/s. Here there are three things to notice: 1) the errors of the period and amplitude do not cover the discrepancy with the actual values of 365.24 days and 29.78 km/s, 2) the very high chi2 value, and 3) the p-value of zero. These suggest very clearly that my errors are wrong. Nevertheless, the signal is clear and the period and orbital speed found are definitely in the right order of magnitude, from which I confirm that my method in general is working. Comparing with the direct differences between observations 1 and 2, 2 and 3, etc, plotted in blue in figure 10, it is also clear that the last step of computing the relative shift between all observations is vital.

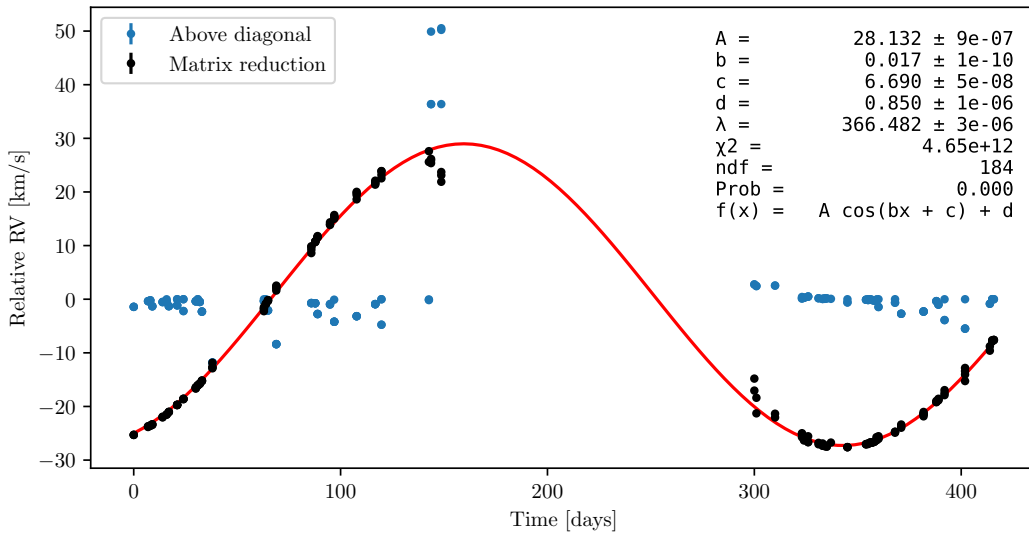


Figure 10: Final relative radial velocity results for 188 observations of HD 34411 using excalibur calibrated but non-barycentric-corrected data (data column `excalibur`). Black: values computed through the overconstrained system approach. Blue: above diagonal of V_r^{ij} .

Although the amount of computations necessary for the described analysis is high, it is possible to execute for 188 observations on a personal computer in about 2 hours. More

details are listed in appendix A.1.

4 Results

Figure 11 shows my extracted relative radial velocity results for HD 34411 compared to those of Lily Zhao et al (also available upon request [7]). As the Earth’s movement has been removed, we are now on the scale of m/s instead of km/s. The method used by Lily Zhao et al is a “chunk-by-chunk” method, where each spectrum is split into 2\AA chunks for each of which an RV is found by shifting a template spectrum to match the observed spectrum. That is to say, a different method from mine. Our results however coincide within the same order of magnitude and with a similar rms of 1.67 m/s and 1.78 m/s, mine and hers respectively. The residuals however do jump around quite a bit, and, of course, my mean error of 0.8 cm/s is not correct.

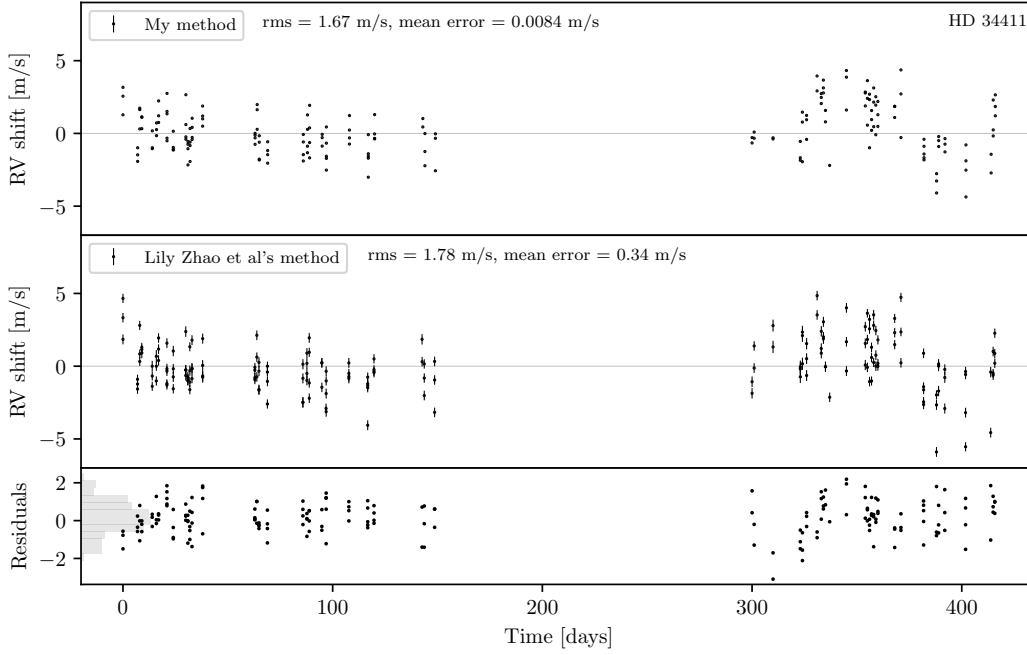


Figure 11: Top: my final extracted RV shifts for HD 34411 using excalibur calibrated, barycentric-corrected data (data column `bary_excalibur`). Center: Lily Zhao et al’s results using a chunk-by-chunk (CBC) analysis technique [7]. Bottom: residuals from subtracting Lily’s results from mine. 188 observations made between 2019-10-8 and 2020-11-27.

I’ve analysed data from three more stars and again compared the results with those of Lily et al. As with HD 34411, my rms values come within a few m/s, the residuals are not small, and my errors are too small. See figures 14, 15 and 16 in appendix B for plots.

5 Discussion

I would have liked to have performed RV extractions on spectra data that I had calibrated myself. But, it was not possible for me get access to spectra data with associated LFC spectra. So the two parts of the project, calibration and rv extractions, have been somewhat separated.

I should have liked to compare my

Although the LFC calibration initially was the main aspect of the project, I ended up spending much more time on the RV extractions.

Future ideas:

- Errors
- Errors from the interpolation, the more points we sample the smaller the error?
- Filters: chauvenet’s, rv cut, z-score cut, always having the barycentric data to confirm if it works is good. Weighting of lines/features based on the stability. Just like Lily et al do in the chunk-by-chunk method [paper](#)..

Future ideas:

- Auto encoder

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

6 Conclusion

7 Acknowledgements

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A RV extractions

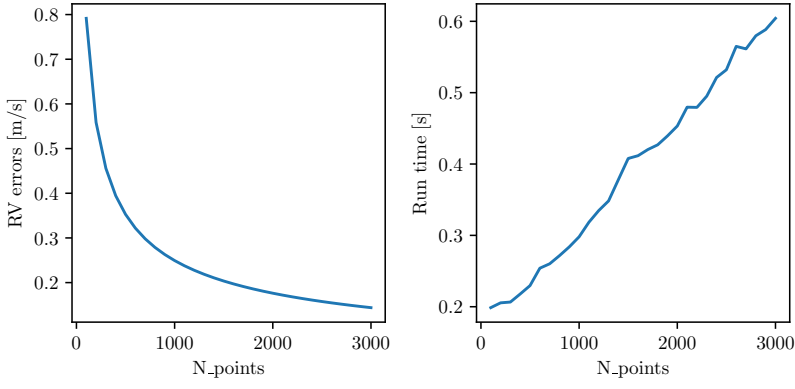


Figure 12: Number of data points in each LFC peak fit, determined by the average distance between peaks in each order.

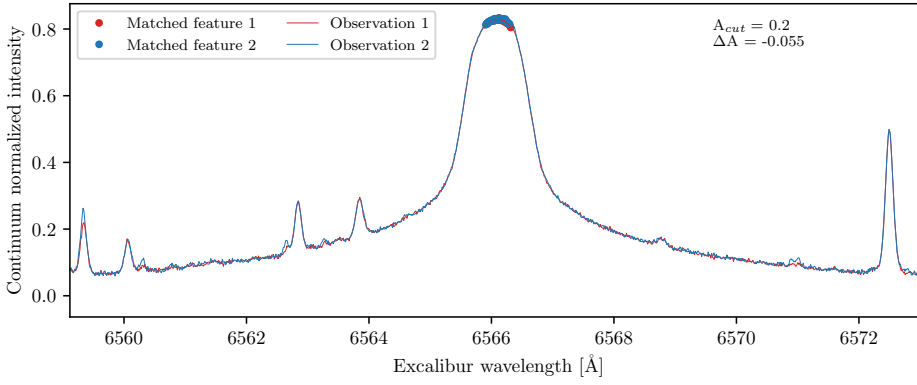


Figure 13: Example of a bad match that made it through the filters. This match is actually correct, but yields a bad result because it is too broad for the standard data slice size.

A.1 Computational run times

todo: add number of function calls for the fitting

todo: add number of peak analyzed and so on

todo: add run times

B Results

RV extracts from three more stars: HD 101501, HD 10700 and HD 26965. In each figure the top shows my final extracted RV shifts using excalibur calibrated, barycentric-corrected data (data column `bary_excalibur`), the center shows Lily Zhao et al's results using a chunk-by-chunk (CBC) analysis technique [7], and the bottom shows residuals from subtracting Lily's results from mine.

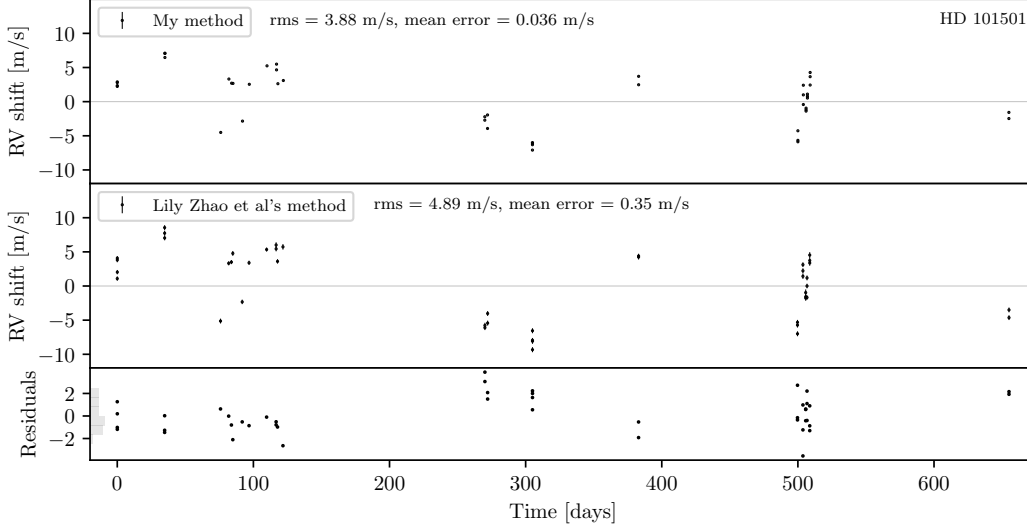


Figure 14: HD 101501
45 observations made between
2019-2-10 and 2020-11-26.

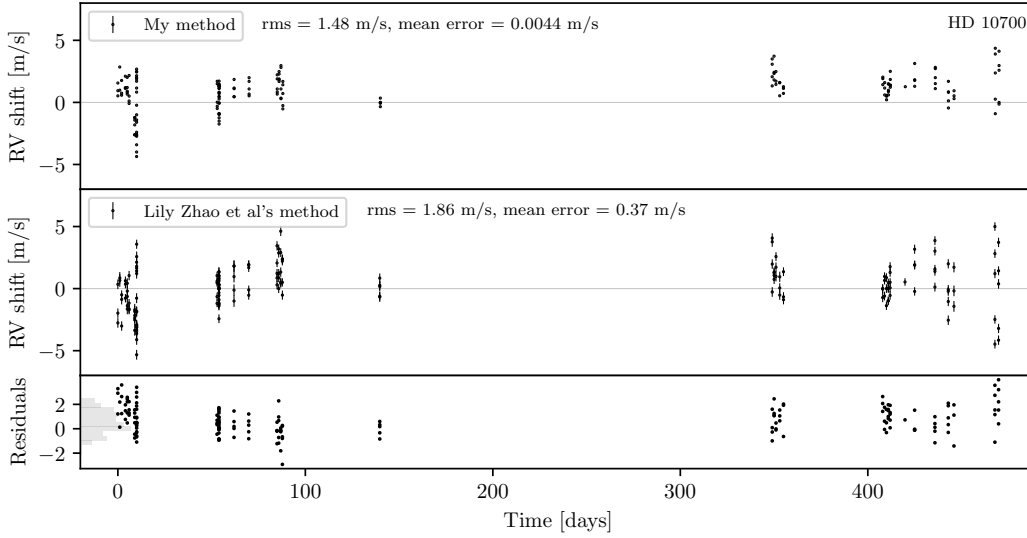


Figure 15: HD 10700
174 observations made between
2019-8-15 and 2020-11-27.

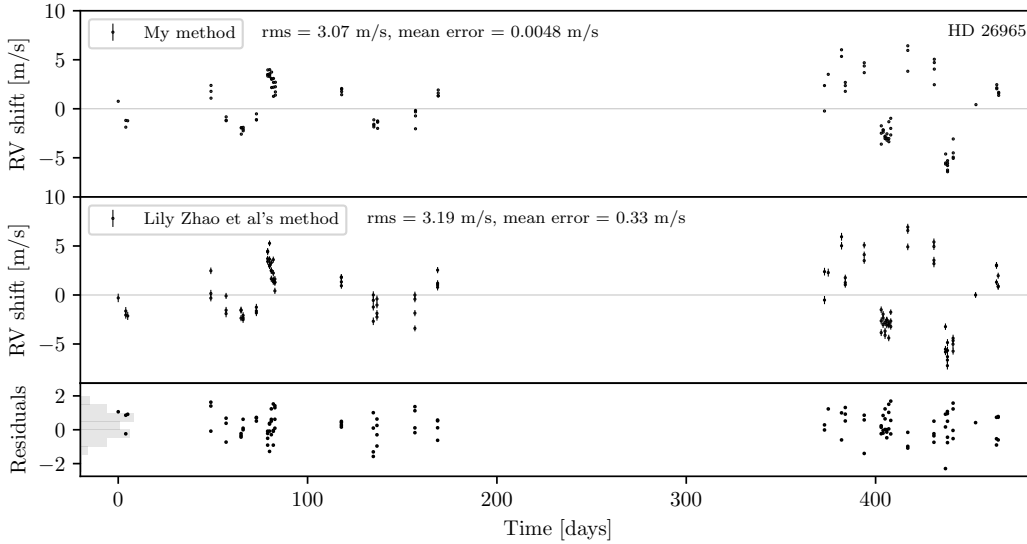


Figure 16: HD 26965
114 observations made between
2019-8-20 and 2020-11-27.