# Fitting Polynomial Baselines to Spectra

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#### Introduction

This document discusses baseline correction methods that can be used with *hyperSpec*. *hyperSpec* provides two fitting functions for polynomial baselines, spc.fit.poly and spc.fit.poly.below.

In contrast to many other programs that provide baseline correction methods, these functions do least squares fits. However, the baselines can be forced through particular points, if this behaviour is needed.

The main difference between the two functions is that spc.fit.poly returns a least squares fit through the complete spectrum that is given in *fit.to* whereas spc.fit.poly.below tries to find appropriate spectal regions to fit the baseline to.

## Syntax & parameters

apply.to: hyperSpec object giving the spectral range, on which the baselines should be evaluated.

If apply is NULL, a hyperSpec object with the polynomial coefficients is returned instead

of evaluated baselines.

poly.order: polynomial order of the baselines

npts.min: minimal number of data points per spectrum to be used for the fit.

npts.min defaults to the larger of 3 times (poly.order + 1) or  $\frac{1}{20th}$  of the number of

data points per spectrum.

If npts.min ≤ poly.order, a warning is issued and npts.min <- poly.order + 1 is used.

noise: a vector giving the amount of noise, see below.

short, user, date: are handed to logentry

#### **General Use**

Both functions fit the polynomial to the spectral range given in *hyperSpec* object *fit.to*. If *apply.to* is not NULL, the polynomials are evaluated on the spectral range of *apply.to*. Otherwise, the polynomial coefficients are returned.

Substracting the baseline is up to the user, it is easily done as hyperSpec provides the - (minus) operator.

## Fitting polynomial baselines using least squares

Commonly, baselines are fit using (single) support points that are specified by the user. Also, usually n+1 support point is used for a polynomial of order n. This approach is appropriate for spectra with high signal to noise ratio.

Such a baseline can be obtained by restricting the spectra in *fit.to* to the respective points (see figure 1):

```
> bl <- spc.fit.poly (chondro [c (1,3),, c(630, 1750)], chondro [c (1,3)])
> plot (chondro [c (1,3)], plot.args = list (ylim = c(200, 600)), col = 1 : 2)
> plot (chondro [c (1,3),, c(630, 1750)], add = TRUE, lines.args = list (type = "p", pch = 20), col = 1:2)
> plot (bl, add = TRUE, col = 1 : 2)
```

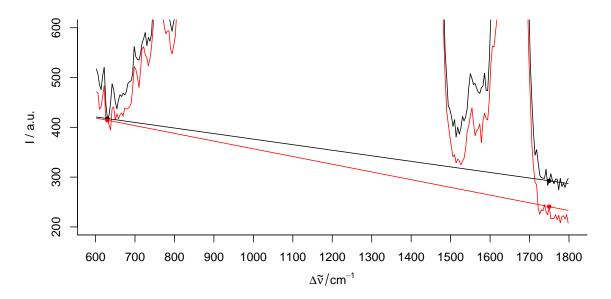


Figure 1: Fitting a linear baseline through two points. If the signal to noise ratio is not ideal, wavelengths that work fine for one spectrum (black) may not be appropriate for another (red).

However, if the signal to noise ratio is not ideal, a polynomial with n+1 supporting points (i.e. with zero degrees of freedom) is subject to a considerable amount of noise. If on the other hand, more data points consisting of baseline only are available, the uncertainty on the polynomial can be reduced by a least squares fit.

Both spc.fit.poly and spc.fit.poly.below therefore provide least squares fitting for the polynomial.

As spc.fit.poly fits to the whole spectral region of fit.to. Thus, the spectra need to be cut to appropriate wavelength ranges that do not contain any signal.

In order to speed up calculations, the least squares fit is done by using the Vandermonde matrix and solving the equation system by qr.solve.

This fit is not weighted. A spectral region with many data points therefore has greater influence on the resulting baseline than a region with just a few data points. It is up to the user to decide whether this should be corrected for by selecting appropriate numbers of data points (e.g. by using replicates of the shorter spectral region).

## The mechanism of automatically fitting the baseline in spc.fit.poly.below

spc.fit.poly.below tries to automatically find appropriate spectral regions for baseline fitting. This is done by excluding spectral regions that contain signals from the baseline fitting. The idea is that all data points that lie above a fitted polynomial (initially through the whole spectrum, then through the remaining parts of the spectrum) will be treated as signal and thus be excluded from the baseline fitting.

The supporting points for the baseline polynomials are calculated iteratively:

- 1. A polynomial of the requested order is fit to the considered spectral range, initially to the whole spectrum given in *fit.to*
- 2. Only the parts of the spectrum that lie below this polynomial plus the noise are retained as supporting points for the next iteration.

These two steps are repeated until either

- no further points are excluded, or
- the next polynomial would have less than npts.min supporting points.

The baselines and respective supporting points for each iteration of spc.fit.poly.below (chondro [1], poly.order = 1) are shown in figure 2.

# Specifying the spectral range

It is possible to exclude spectral regions that do not contribute to the baseline from the fitting, while the baseline is used for the whole spectrum. This selection of appropriate spectral regions is essential for <code>spc.fit.poly</code>. But also <code>spc.fit.poly.below</code> can benefit from narrower spectral ranges: the fitting gains speed. The default value for <code>npts.min</code> depends on the number of data points per spectrum. Thus one should consider using more support points than the default value suggests.

```
> system.time (spc.fit.poly.below (chondro[], NULL, npts.min = 20))
   user system elapsed
2.120  0.030  2.149
> system.time (spc.fit.poly.below (chondro [,, c (min ~ 700, 1700 ~ max)], NULL, npts.min = 20))
   user system elapsed
1.100  0.000  1.098
```

The choice of the spectral range in fit.to influences the resulting baselines to a certain extent, as becomes clear from figure 3.

## Fitting polynomials of different orders

Figure 4 shows the resulting baseline polynomial of spc.fit.poly.below (chondro [1], poly.order = order) with order = 0 to 3 for the first spectrum of the chondro data set.

#### The noise level

Besides defining a minimal number of supporting points, a "noise level" may be given. Consider a spectral range consisting only of noise. The upper part of figure 5 illustrates the problem. As the baseline fitting algorithm cannot distinguish between noise and real bands appearing above the fitted polynomial, the resulting baseline (black) is too low if the noise parameter is not given.

Setting the noise level to 10 (2 standard deviations), the fitting converges immediately with a much better result. The resulting baselines for spc.fit.poly.below (chondro [1], poly.order = 1, noise = 12) of the whole spectrum are shown in the middle and lower part of figure 5

noise may be a single value for all spectra, or a vector with the noise level for each of the spectra separately.

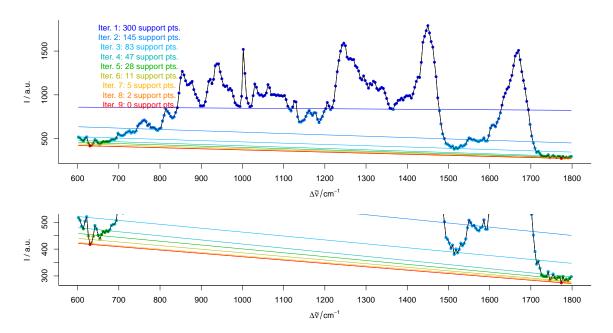


Figure 2: Iterative fitting of the baseline. The dots give the supporting points for the baselines in the same colour. The lower part is a magnification of the intensity axis.

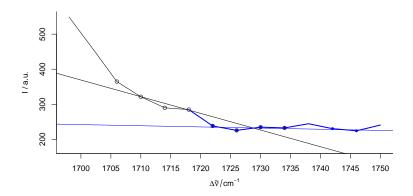


Figure 3: Influence of fit.to on the baseline polynomial. The black baseline is fit to the spectral range  $1700-1800~{\rm cm^{-1}}$ , the blue to  $1715-1800~{\rm cm^{-1}}$  only (dots & circles: supporting points).

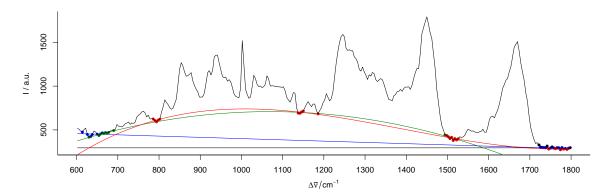


Figure 4: Baseline polynomial fit to the first spectrum of the chondro data set of order 0 (black), 1 (blue), 2 (green), and 3 (red). The dots indicate the points used for the fitting of the polynomial.

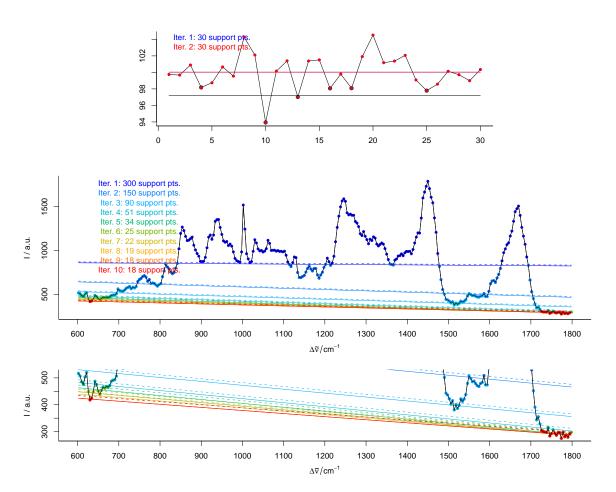


Figure 5: Iterative fitting of the baseline with noise level. Upper part: effects of the noise parameter on the baseline of a spectrum consisting only of noise and offset: without giving noise the resulting baseline (black) is clearly too low. A noise level of 10 results in the red baseline. The middle and lower part show the baseline fitting with noise level on the complete spectrum. Colour: iterations, dots/circles: supporting points for the respective baselines. Dashed: baseline plus noise. All points above this line are excluded from the next iteration.