



The CPL Wavelength Calibration method using Iterative Cross-Correlation

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Introduction

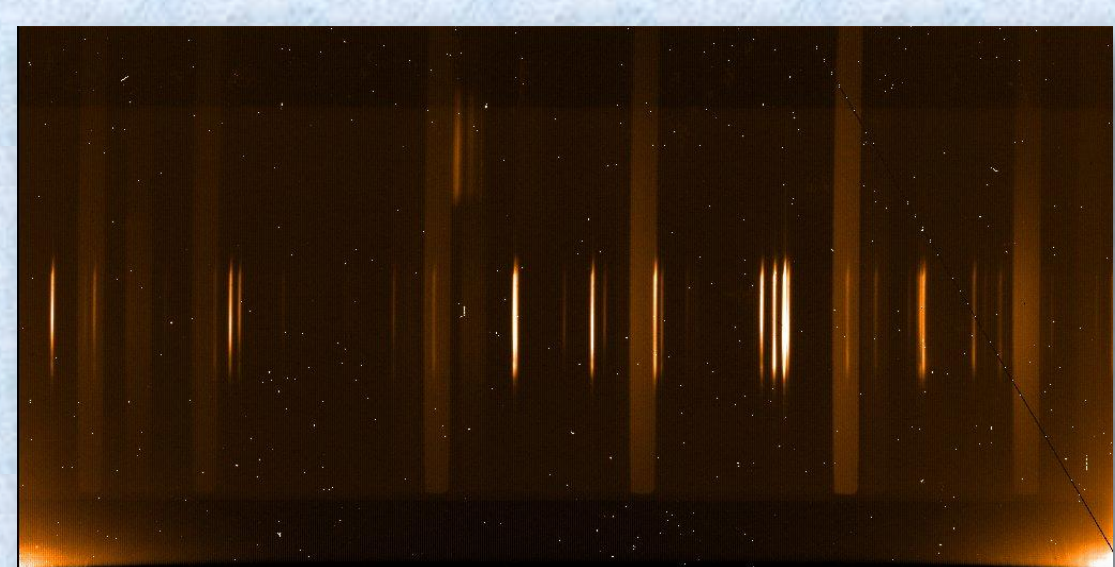
Most of the ESO VLT data reduction pipelines need to calibrate the instruments in wavelength (wavelength = P(pixel)). The generic method described here allows to search iteratively this relation, starting from a first guess, going through a well defined search space identified by wavelength bounds around this first guess.

The way the inputs are changed during the successive iterations depend on the precision of the first guess, the accuracy of the requested calibration, the requested degree of the polynomial solution, the precision of the catalog, etc...

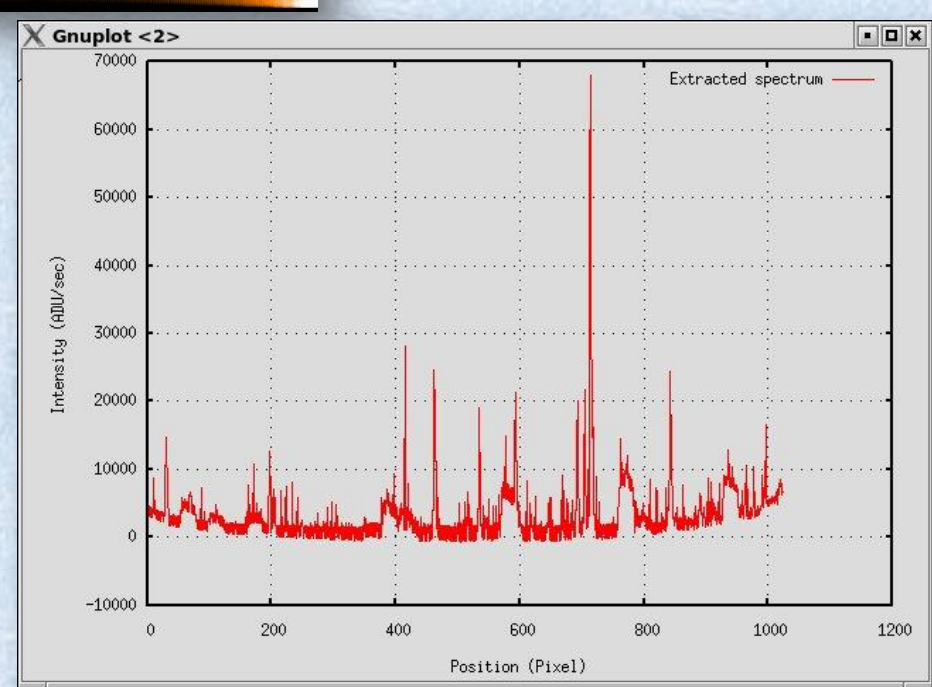
The **CPL** (Common Pipeline Library) function called iteratively has been designed to have a clear and meaningful interface, making it useable in any context.

1. Instrument specific pre-processing

The algorithm inputs (the spectrum to calibrate, the catalog, **degree**, **Po**, **WL_ERR**, **nsamples**) first need to be prepared.



Spectrum Extraction



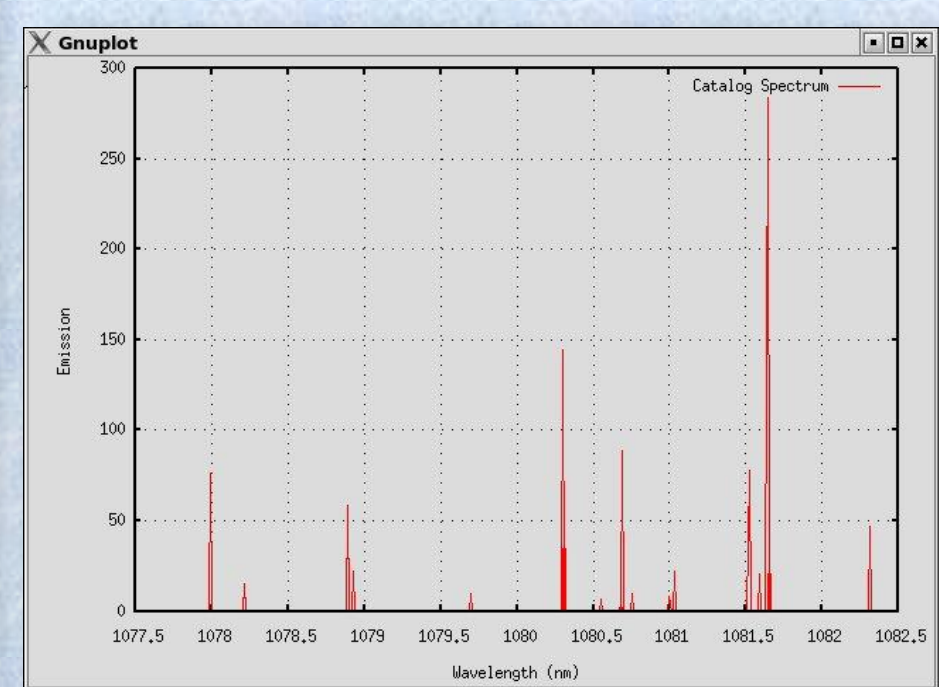
Wavelength range estimation (Physical model, Instrument setting)
wavelength = **Po**(pixel)

Lines catalog (OH, Argon lines, HITRAN model, Absorption lines)

Po(1)

| Wavelength | Emission |
|------------|----------|
| 1077.541 | 16.74 |
| 1077.621 | 93.61 |
| 1077.731 | 4.862 |
| 1077.991 | 76.46 |
| 1078.211 | 15 |
| 1078.891 | 58.45 |
| 1079.931 | 22.13 |
| 1079.71 | 9.439 |
| 1080.21 | 144.8 |
| 1080.311 | 79.28 |
| 1080.251 | 6.445 |
| 1080.481 | 5.031 |
| 1080.691 | 89.11 |
| 1080.751 | 9.661 |
| 1081.051 | 8.051 |
| 1081.051 | 22.46 |
| 1081.821 | 77.64 |
| 1081.891 | 20.3 |
| 1081.641 | 283.7 |
| 1081.851 | 44.8 |
| 1082.311 | 47.13 |

Po(1024)



Wavelength Error **WL_ERR** of the estimation

Precision of the wished solution:

- number of samples **nsamples** for the solution space sampling
- **degree** of the final polynomial solution

2. Instrument Independent Iterative Wavelength Calibration

<http://www.eso.org/cpl>

Function Documentation

```
cpl_polynomial* cpl_wcalib_xc_best_poly ( const cpl_vector * spectrum,
const cpl_bivector * lines_catalog,
int degree,
const cpl_polynomial * guess_poly,
double wl_error,
int nsamples,
double * xc )
```

Find the best polynomial in a given range.

Parameters:

- spectrum The spectrum vector
- lines_catalog The lines catalog
- degree The polynomial degree
- guess_poly Guess Dispersion Polynomial
- wl_error Search range around the anchor points
- nsamples Number of samples around the anchor points
- xc Cross-correlation factor (returned)

Returns:

the best polynomial or NULL in error case

Algorithm

The CPL function first defines the solutions space. This space is defined using **Po**, **WL_ERR**, **degree** and **nsamples**. As described by the drawing, a number ($\text{nsamples}^{\text{degree}+1}$) of candidate polynomials of degree **degree** are defined passing through the **nsamples** positions around the points A1, ..., A_{degree+1}.

Each of these candidates is used to assign wavelengths to the input extracted spectrum. The result is compared to the catalog and a correlation factor is computed. The closer the spectrum and the catalog, the higher the correlation factor, the better the solution.

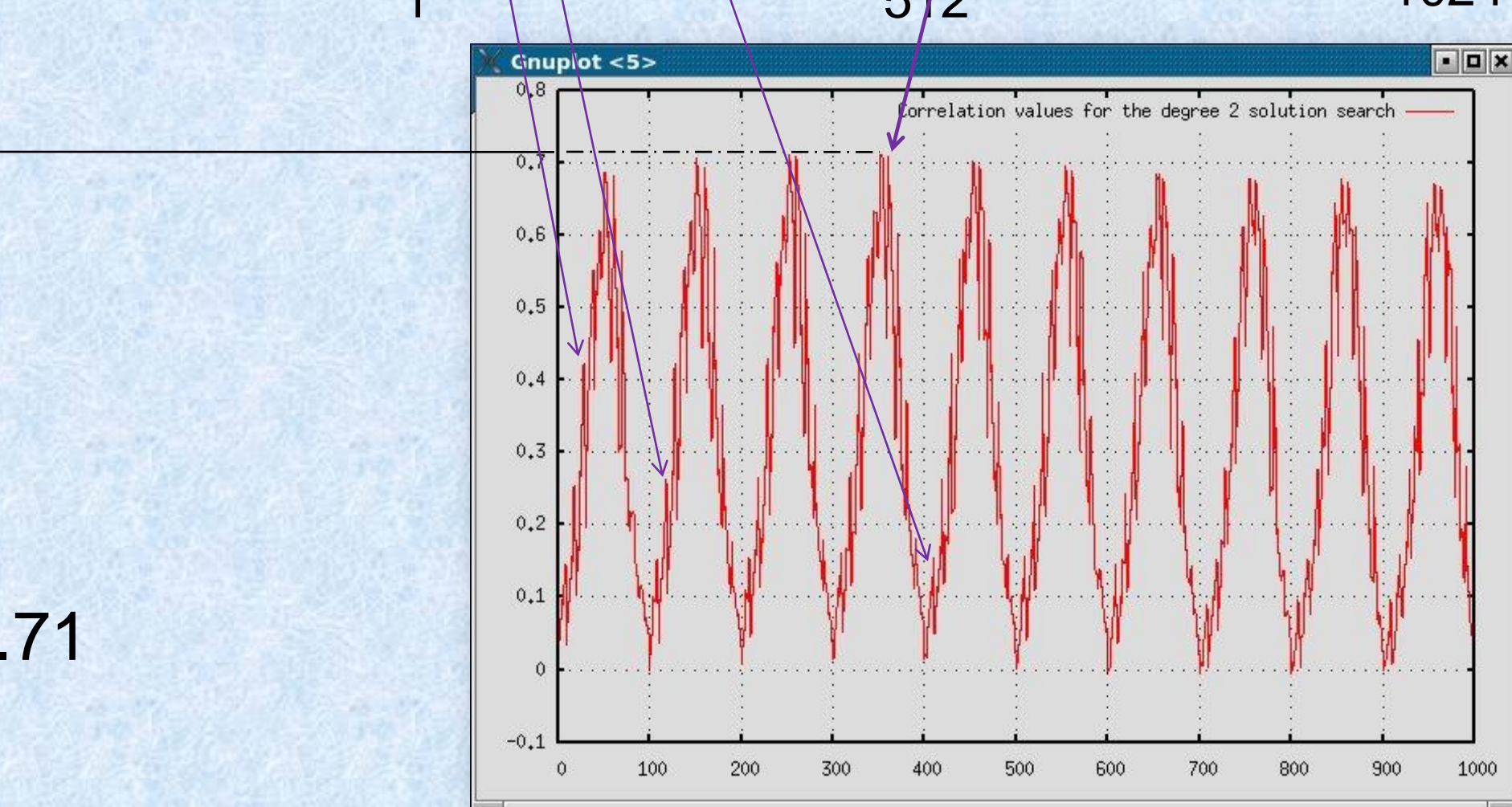
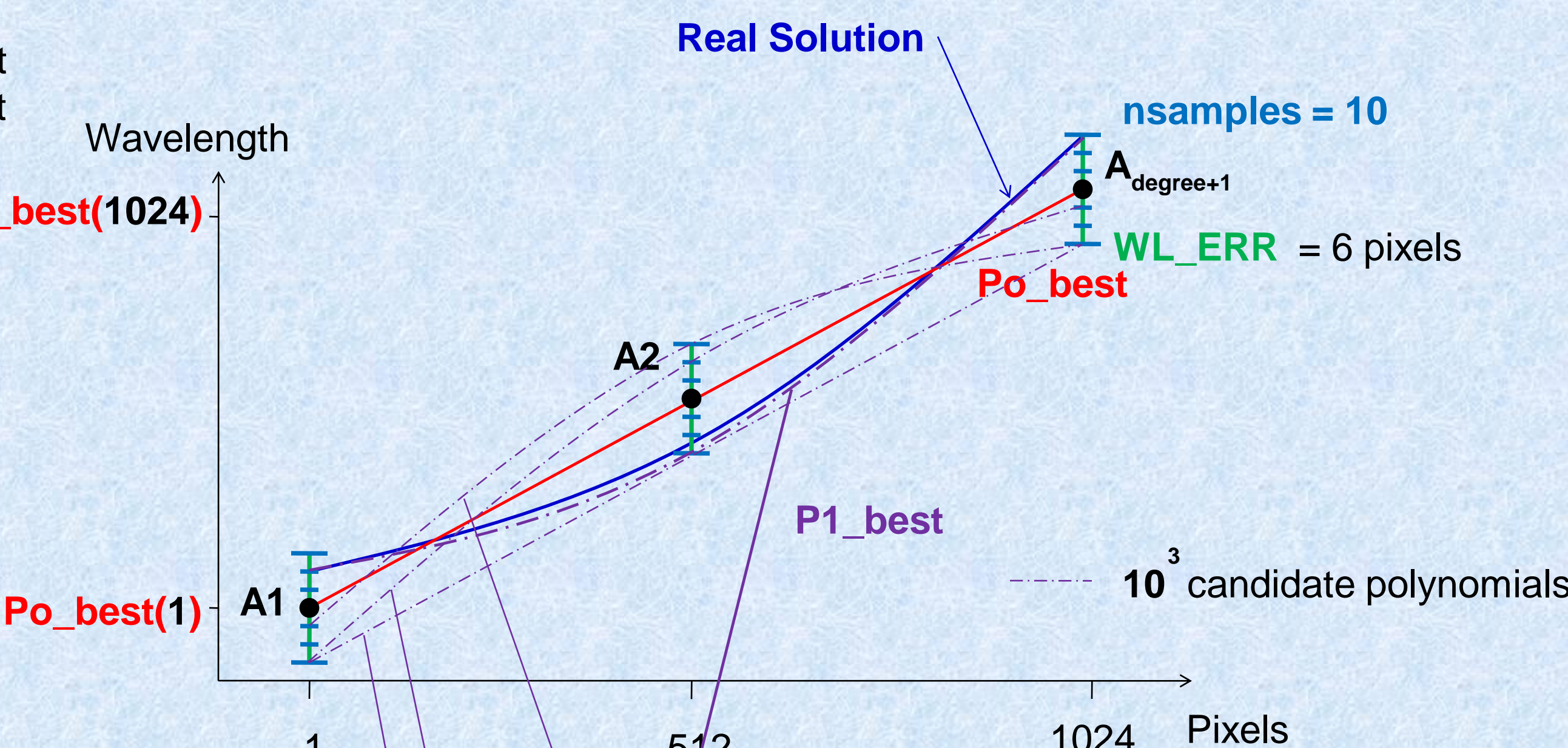
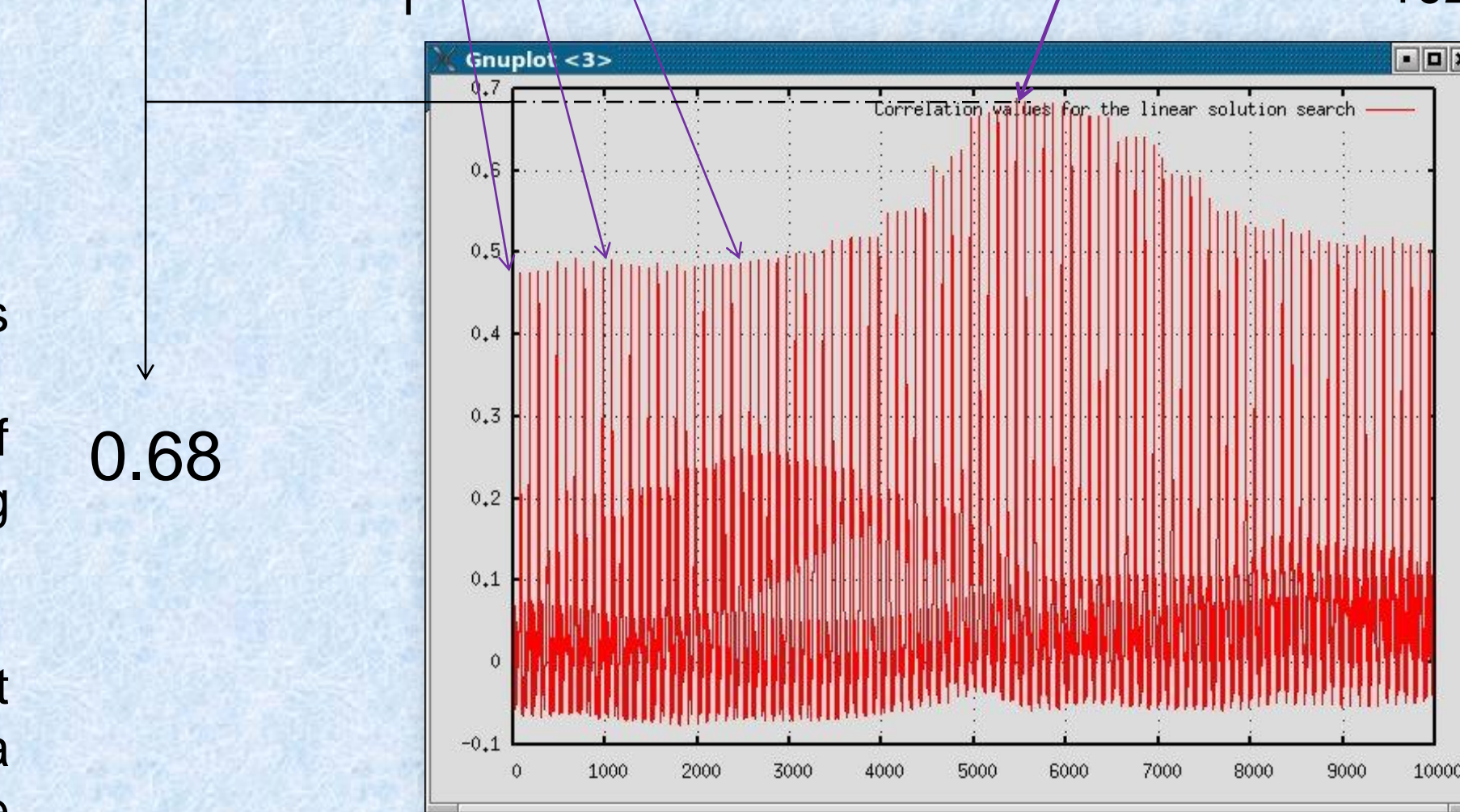
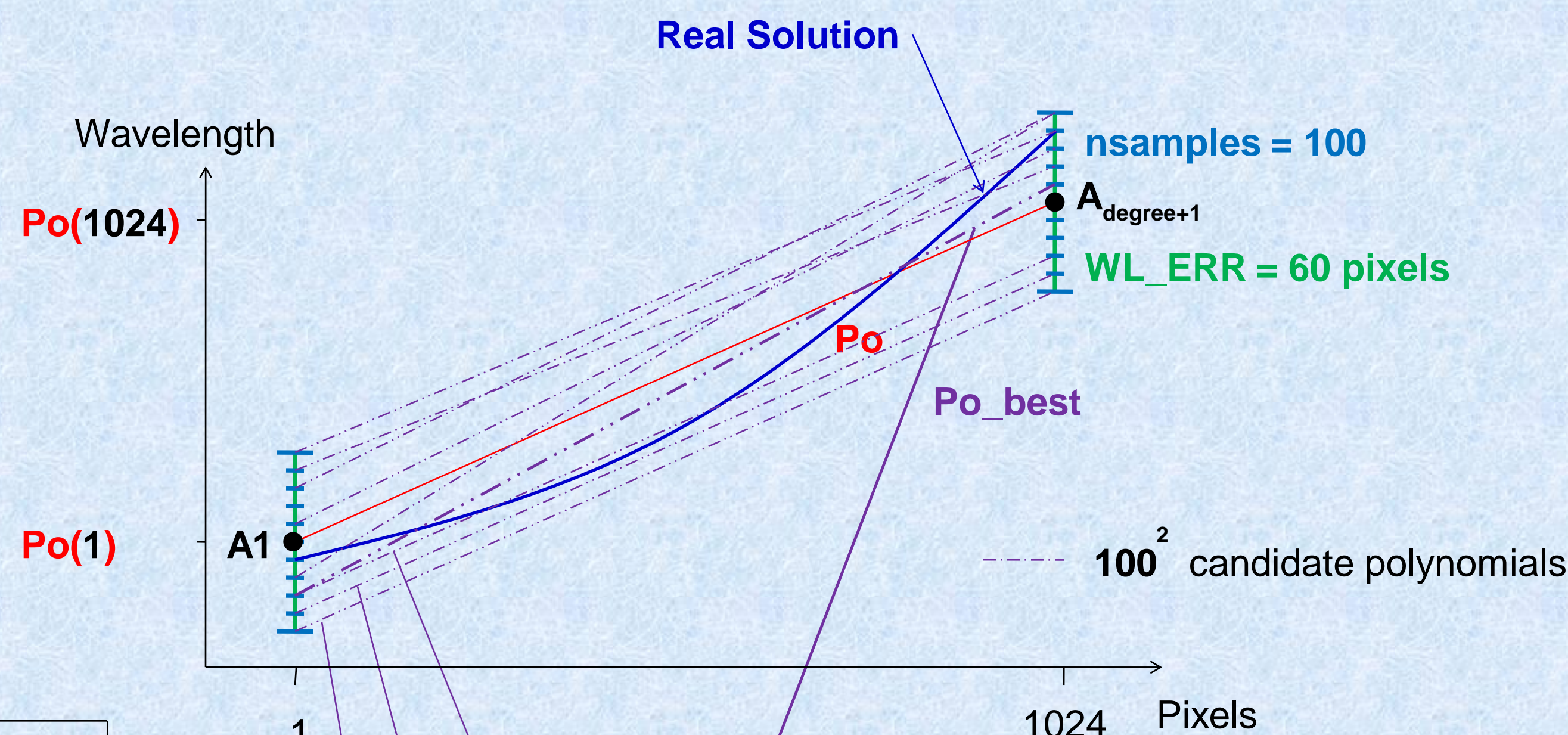
The solution is the polynomial **Po_best** corresponding to the highest correlation value. This polynomial is used as first guess in the next iteration.

Iteration

The number of candidate polynomials that are evaluated can easily explode ($\text{nsamples}^{\text{degree}+1}$) especially if a high precision calibration (with a high **degree**) is requested. To keep this number reasonable, we need to proceed iteratively. The first step refines the guess polynomial (high **nsamples** but low **degree**). This refining allows to reduce the **WL_ERR** in the second step (and thus the **nsamples**) and to increase the **degree**.

cpl_wcalib_xc_best_poly(spectrum, lines, 2, **Po_best**, **WL_ERR**/10, **nsamples**/10, &xc);

cpl_wcalib_xc_best_poly(spectrum, lines, **degree**, **P1_best**, **WL_ERR**/100, **nsamples**/10, &xc);

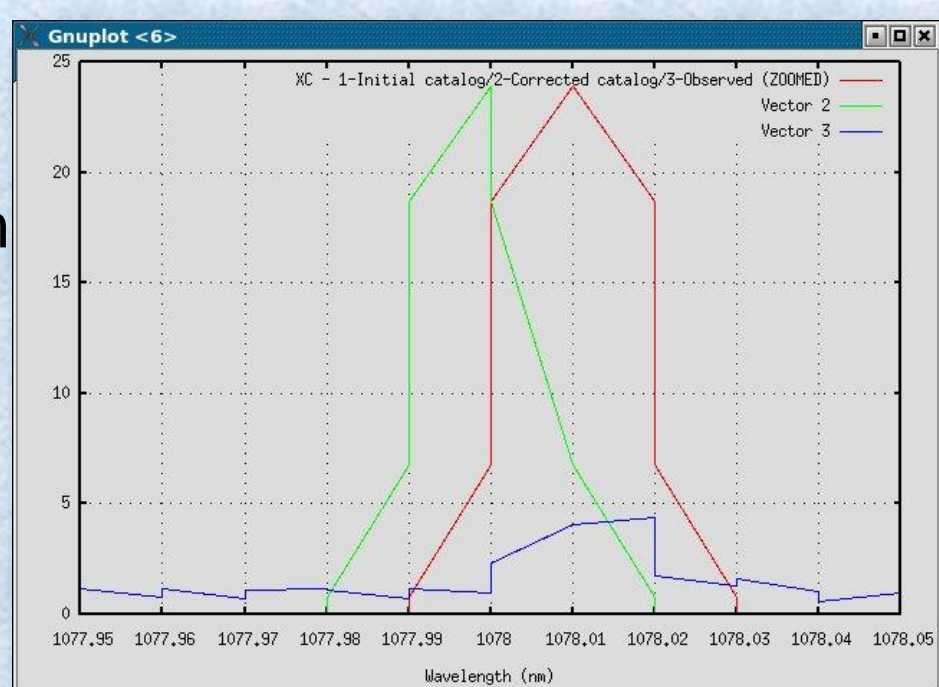


3. Application on CRIRES data through the CRIRES pipeline

<http://www.eso.org/pipelines>

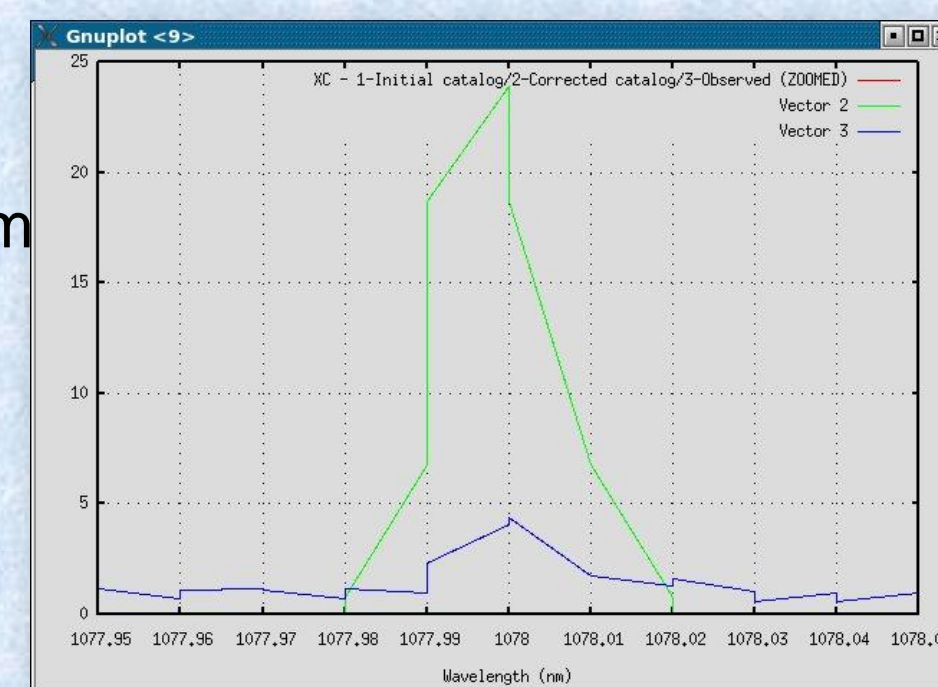
Iteration 1 :

degree = 1
WL_ERR = 0.3 nm
nsamples = 100



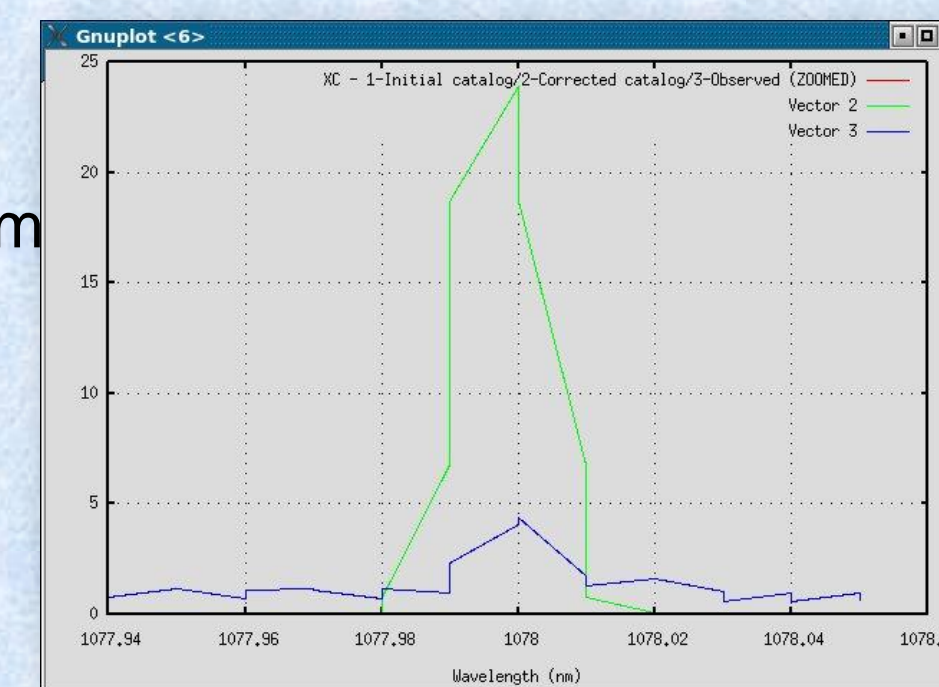
Iteration 2 :

degree = 2
WL_ERR = 0.06 nm
nsamples = 10



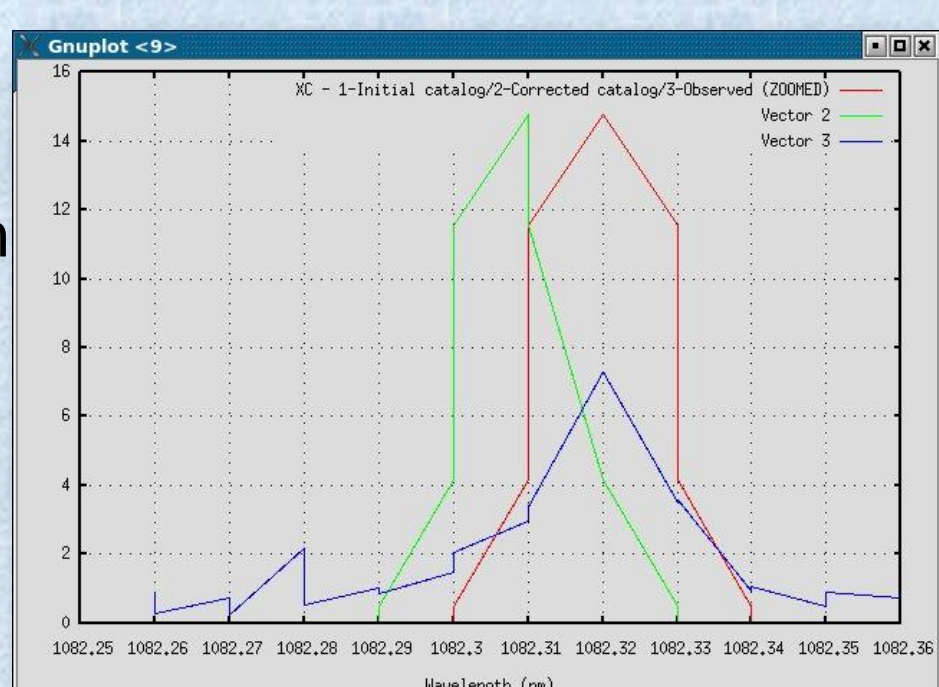
Iteration 3 :

degree = 3
WL_ERR = 0.01 nm
nsamples = 10



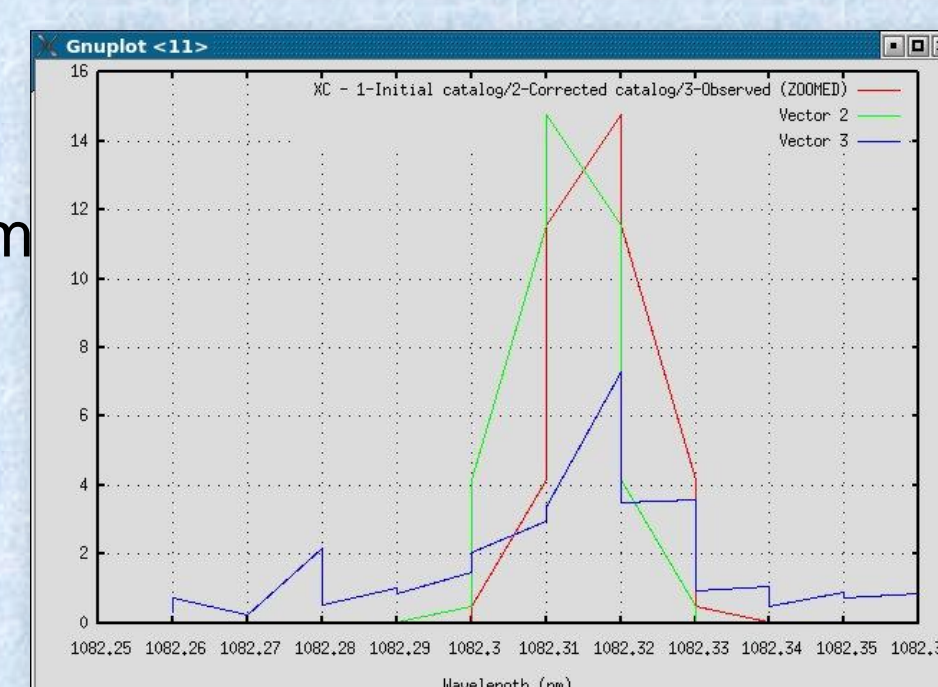
Iteration 1 :

degree = 1
WL_ERR = 0.3 nm
nsamples = 100



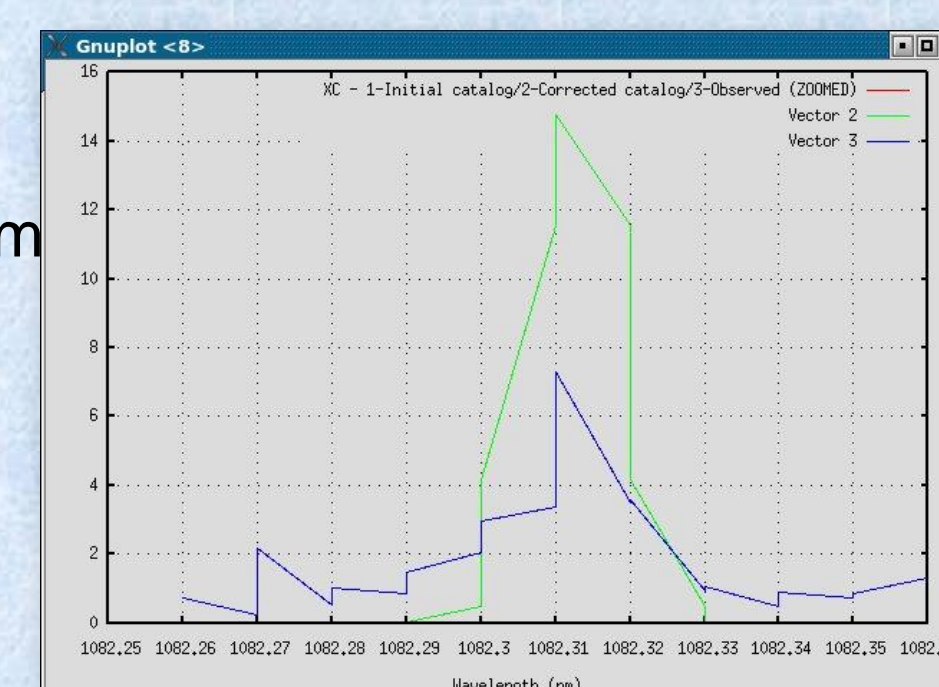
Iteration 2 :

degree = 2
WL_ERR = 0.06 nm
nsamples = 10



Iteration 3 :

degree = 3
WL_ERR = 0.01 nm
nsamples = 10



A calibration lamp is used here to calibrate the instrument in wavelength. The plot shows the **calibrated extracted spectrum** and the **catalog emission lines**. The **red** plot shows the catalog position regarding the spectrum using the guess polynomial solution. A good solution is obtained when **green** and **blue** are well aligned.

