

SFIT4 – Retrieval parameters

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Retrieval parameters



Retrieval parameters

Overview

rt The main parameter `rt` can be used to switch off and on the retrieval altogether.

rt.lm Switches on or off the Levenberg-Marquardt iteration scheme.

rt.convergence the iteration is considered converged when $rt.convergence > D_CHI = (CHI_2_MAX_{i-1} - CHI_2_MAX_i)$

rt.max_iteration maximum number of iterations

For all retrieval parameters a priori is given and a standard deviation σ in the form

rt.x.apriori the a priori of a given value. It is actually applied in the forward calculation. Meaning it can also be used in forward calculations.

rt.x.sigma the entry in the S_A matrix corresponding to this parameter.



Retrieval parameters

Overview

`rt` The main parameter `rt` can be used to switch off and on

```
ITER  FIT_RMS      GAMMA    CHI^2_X    CHI^2_Y      CHI^2  CHI^2_OLD      DCHI^2
  1    25.5294    1.00E+05    0.000     886.616
  2    15.7601    1.00E+05    0.000     332.528    332.528    886.616    554.088226
  3     9.8038    1.00E+04    0.004     134.162    134.166    332.528    198.361710
  4     5.1080    1.00E+03    0.038      32.608     32.646    134.166    101.519910
  5     2.6411    1.00E+02    0.105       6.402      6.507     32.646     26.139551
  6     1.9136    1.00E+01    0.217       2.831      3.048      6.507      3.458828
  7     1.8797    1.00E+00    0.281       2.594      2.875      3.048      0.172872
  8     1.8844    1.00E-01    0.300       2.573      2.873      2.875      0.002286

FINAL:      MEAN_SNR= 86.6267  MEAN_FIT_RMS(%)= 1.88443  NVAR= 185  NFIT= 3477

  BAND  SCAN  RMSSNR (CALCULATED)  (EFFECTIVE)  (RETRIEVED)  CHI^2
    1    1      53.98              53.98        49.55         1.19
    2    1      57.52              57.52        56.78         1.03
    3    1     154.48         154.48       110.55         1.95
    4    1     147.91         84.62        50.79         2.92

NEGATIVE MIXING RATIO VALUES FOUND FOR : 03668
NEGATIVE MIXING RATIO VALUES FOUND FOR : 03686
```

forward calculations.

`rt.x.sigma` the entry in the S_A matrix corresponding to this parameter.



Retrieval parameters

Construction of the S_A matrix.

The S_A matrix is constructed from the sigma values given. How this is actually done, depends on the parameters. In principle the S_A is constructed as a diagonal matrix and inverted in the code to yield the S_A^{-1} matrix. Some caveats:

`gas.profile.x.correlation` off diagonals using the sigma values as maxima

`.type = 1` gaussian with FWHM = `.width`

`.type = 2` exponentially with FWHM = `.width`

`.type = 3` not used

`.type = 4` the S_A matrix is read in from `file.sa_matrix`

`.type = 5` the S_A^{-1} matrix is read in from `file.sa_matrix`



Retrieval parameters

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- `gas.profile.x.regmethod` lets you chose between OE optimization and the Thikonov-Phillips regularization with L1 constraint (smoothness constraint)
- `.type = 'OE'` optimal estimation (Rodgers, 2000)
- `.type = 'TP'` Thikonov-Phillips with smoothness constraint
- `.lambda` strength of the regularization in TP. The higher the value the less is the regularization.

The smoothness constraint is calculated from `file.stalayers` in order to adapt for non-unique altitude layering. The matrix is scaled using the `gas.profile.x.sigma` values.



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FIT_RMS mean variance of the residuum

GAMMA the Levenberg Marquardt Parameter

CHI_2_X A measure of the deviation of the retrieved state from the A PRIORI



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CHI_2_Y A measure for the retrieval quality

$$\chi_Y^2 = \frac{(y_M - y_C)^T S_e (y_M - y_C)}{m}$$

$\chi_Y^2 = 1$ if the residuum is reduced to the noise as specified in S_e



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- ▶ In the lower half, the retrieval diagnostics for the last calculation are shown for each MW.
- ▶ A warning if retrieved profiles have negative parts



Retrieval parameters

Wave number scaling and shifting

`rt.wshift` wave number shift.

- ▶ Shift works on the internal grid
band.X.calc_point_space
- ▶ This is only useful for microwindows (small) because the mismatch is a wavenumber dependent polynomial. For small wave number regions, this can be approximated by a shift.
- ▶ This is on top of `rt.wave_factor`, which is a scaling.
- ▶ The artificial grid needs to be more dense than the measured grid

`rt.dwshift` wave number shift for each retrieved gas separately, except the first retrieved one.

- ▶ all lines of each gas are shifted by the same amount (!!!)



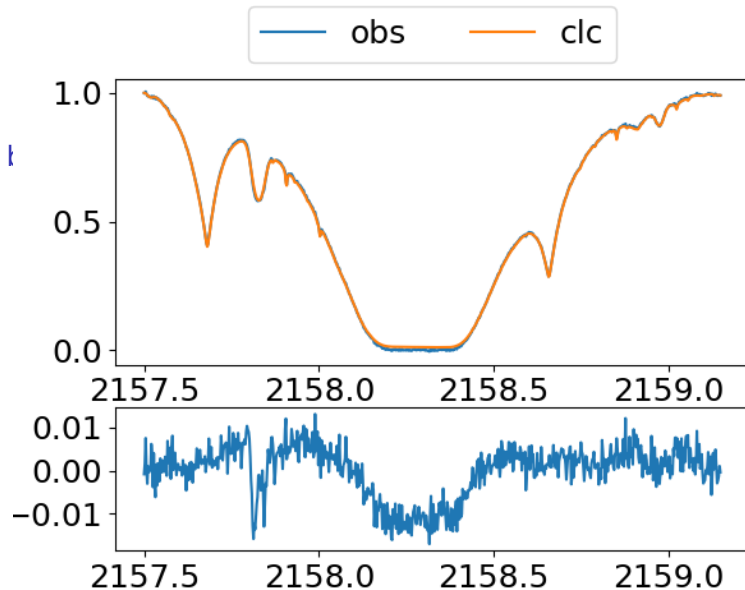
offset in window

`band.zshift` calculates and retrieves an offset in the microwindow. Two types:

`.type = 1` retrieves the offset in this MW **ONLY ONE!!!**



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`.type = 2` uses the offset which is retrieved in another microwindow. THIS MUST BE LATER THAN THE MW USED FOR **ZSHIFT.TYPE=1**

`band.zshift.apriori` is also an **FW parameter**



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`band.zshift.apriori` is also an **FW parameter RETRIEVAL ONLY POSSIBLE IF THERE IS AN SATURATED PART IN THE MW.**



channeling in window

see discussion led by Minqiang Zhou

