

# SFIT4 – kB matrix calculation

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# KB matrix calculation



# Error calculation

## Recall

Forward model and its derivative

$$y = F(x) + \epsilon$$
$$K_{g,b} = \frac{\partial F}{\partial x_{g,b}}$$

The contribution function matrix is used to calculate the state from the spectrum  $x = Dy$ . The averaging kernel  $A = DK$  provides a mappint from the original state to the retrieved state:  $\hat{x} = x_A + DK(x - x_A)$ . The complete error analysis for the retrieval is (compare Rodgers,2000)

$\hat{x} - x$	$= (A - I)(x - x_A)$	Smoothing error
	$+ D_y \epsilon$	Retrieval noise
	$+ D_y \Delta F(x, b, \hat{b})$	Forward model error
	$+ D_y K_b (b - \hat{b})$	model parameter error



## Model parameter error

The parameter  $x_b$  consist of two groups:

1. Parameters retrieved along with the atmospheric state  $x_g$ , e.g. interfering gases, instrumental parameters like slope or wavenumber shift.
2. Parameters which are not retrieved, but need to be used like FOV, MAX\_OPD, temperature profile

The parameters in group 2 have errors and this are accounted for in the parameter error. In order to calculate them, we need the  $K_b$  matrix for the parameters which are not retrieved:

$$K_b = \frac{F}{\partial b}$$

They are calculated in SFIT4 as the KB matrix. It is switched on in the sfit4.ctl file as

kb = T



# KB matrix calculation

- ▶ The subkeys of kb can be chosen to specify for which parameters error matrices are calculated.
- ▶ Some parameters can also be retrieved, e.g. rt.curvature for the curvature of the transmission function.
- ▶ Parameters which are retrieved but also specified in kb are only retrieved, no *KB* entry is calculated for them. e.g.

...

rt.slope = T

kb.slope = T

...

no kb entry for the slope is calculated.



# KB matrix calculation

For some parameters it does not make sense to retrieve errors, either they are negligible or they are better retrieved.

**kb.dwshift** dwshift shifts all lines of all gases except the retrieval gas. The line position error is taken care of in kb.line

**fw.wshift** this is usually retrieved and thus contributes to the interference error.

**kb.max\_opd** the error for this is usually negligible. Because max\_opd should be known to a very high accuracy. NOTE: Might be different for low resolution spectra.



# KB matrix calculation

Error for apod and phase function

- ▶ If `fw.apodisation_fcn` and/or `fw.phase_fcn` are calculated as polynomial or fourier series (not retrieved!) the parameters are used to calculate the error matrix.
- ▶ if `fw.apodisation_fcn` or `fw.phase_fcn` is set `FALSE`, in the error calculation a three order polynomial is assumed and used to estimate the KB for the ILS

