# CASSIS as a plug-in tool in HIPE 10.0 Charlotte Vastel David Teyssier



18-19 April, 2013





Université Paul Sabatier To install, select Plug-ins from the Tools menu. Click Install new and enter the following URL:

http://cassis.irap.omp.eu/download/cassis\_3.2.jar

If you have an old version of this plugin, please remove it (using "File" "Remove completely" on the plug-ins panel) and install the new version.

Open Hipe 10 with enough memory:

java.vm.memory.max=5120m

hcss.ia.pal.pool.lstore.dir=/Users/vastel/.hcss/lstore

Open the oribar\_ALL\_clean.fits file through the navigator or through the Console:

oribar\_ALL\_clean = fitsReader(file = '/Users/vastel/DATA/ HIFI/ORION\_BAR/oribar\_ALL\_clean.fits')

It creates a variable: oribar\_ALL\_clean. Right click on the oribar\_ALL\_clean variable and "open with" "CASSIS Spectrum1dLineAnalysis"

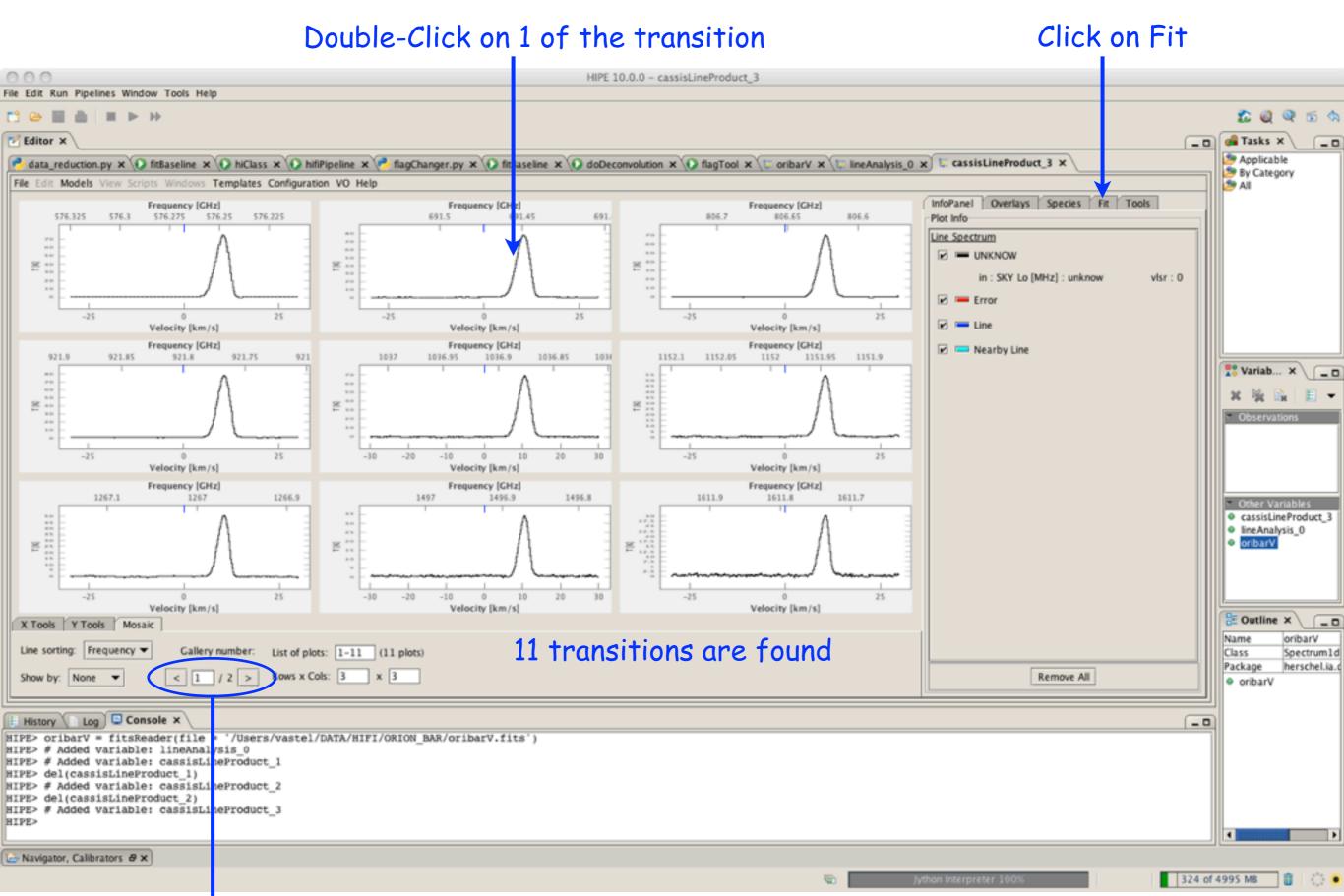
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Select the upper energy: 2000K means that it will look for all the transitions for energy lower than 2000K

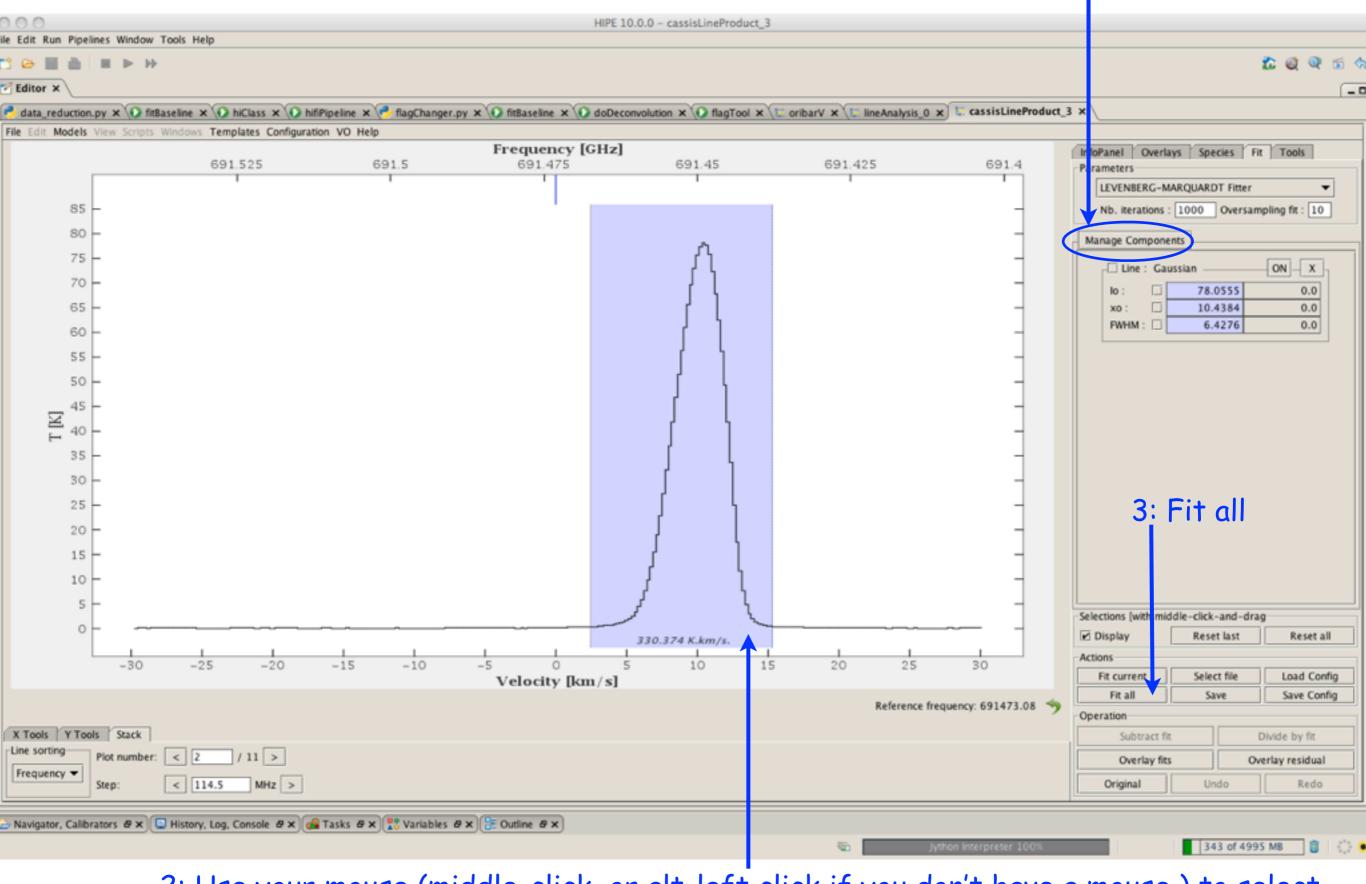
### press DISPLAY

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			Package herschel.ia.c
			<ul> <li>oribarV</li> </ul>
			1
History Log Console ×			
<pre>HIPE&gt; oribarV = fitsReader(file = '/Users/vastel/DATA/HIFI/ORION_BAR/oribarV.fits') HIPE&gt; # Added variable: lineAnalysis_0</pre>			
HIPE> # Added variable: cassisLineProduct_1			
HIPE> del(cassisLineProduct_1) HIPE> # Added variable: cassisLineProduct_2			
HIPE> del(cassisLineProduct_2)			
<pre>HIPE&gt; # Added variable: cassisLineProduct_3 HIPE&gt;</pre>			
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browse the transition from 1 window to another

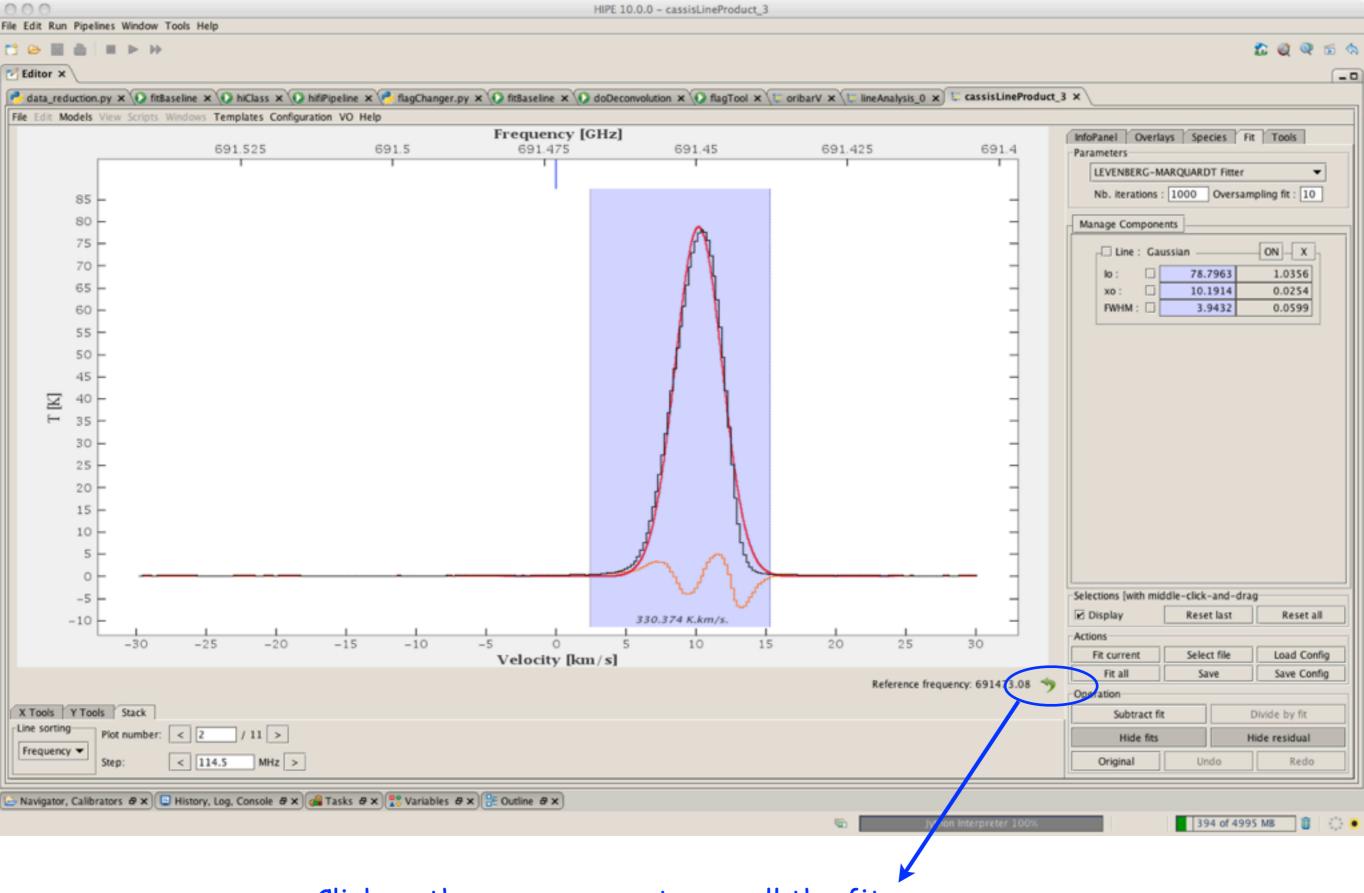
1: Select Gaussian (the baselines are already substracted in these data, but you can also fit the continuum)



2: Use your mouse (middle-click, or alt-left click if you don't have a mouse ) to select the area to be fitted (here in blue)

HIPE 10.0.0 - cassisLineProduct\_3 dit Run Pipelines Window Tools Help 🏠 🌒 🔍 卮 🔕 e ditor X --data\_reduction.py 🗙 🜘 fitBaseline 🗙 🕼 hiClass 🗙 🜘 hifiPipeline 🗙 🥐 flagChanger.py 🗙 🜘 fitBaseline 🗙 🕼 doDeconvolution 🗙 🚺 flagTool 🗙 🔚 oribarV 🗙 🔚 lineAnalysis\_0 🗙 🖫 cassisLineProduct\_3 🗴 Edit Models ws Templates Configuration VO Help View Scripts Windo Frequency [GHz] InfoPanel Overlays Species Fit Tools 691.425 691.525 691.5 691.475 691.45 691.4 Parameters LEVENBERG-MARQUARDT Fitter Ŧ 85 Nb. iterations : 1000 Oversampling fit : 10 80 Manage Components 75 - Line : Gaussian ON - X 70 78.0555 0.0 lo 10.4384 65 0.0 XO : 6.4276 FWHM : 0.0 60 55 50 45 000 Write the log in a file T [K] 40 What do you want? 35 30 Create a new File Open an existing File Cancel 25 20 15 10 5 Selections (with middle-click-and-drag 0 Display Reset last Reset all 330.374 K.km/s. Actions -25 -20 -100 10 25 -30 -15 -5 15 30 20 Fit current Select file Load Config Velocity [km/s] Fit all Save Save Config Reference frequency: 691473.08 🥎 Operation Tools Y Tools Stack Subtract fit Divide by fit e sorting / 11 > Plot number: < 2 Overlay fits Overlay residual requency 🔻 MHz > < 114.5 Undo Step: Original Redo avigator, Calibrators 🖉 🗙 🔁 History, Log, Console 🦉 🗙 🖓 Tasks 🧬 🗙 👯 Variables 🖉 🗙 🚼 Outline 🦉 🗙 350 of 4995 MB 70 - C - 💌

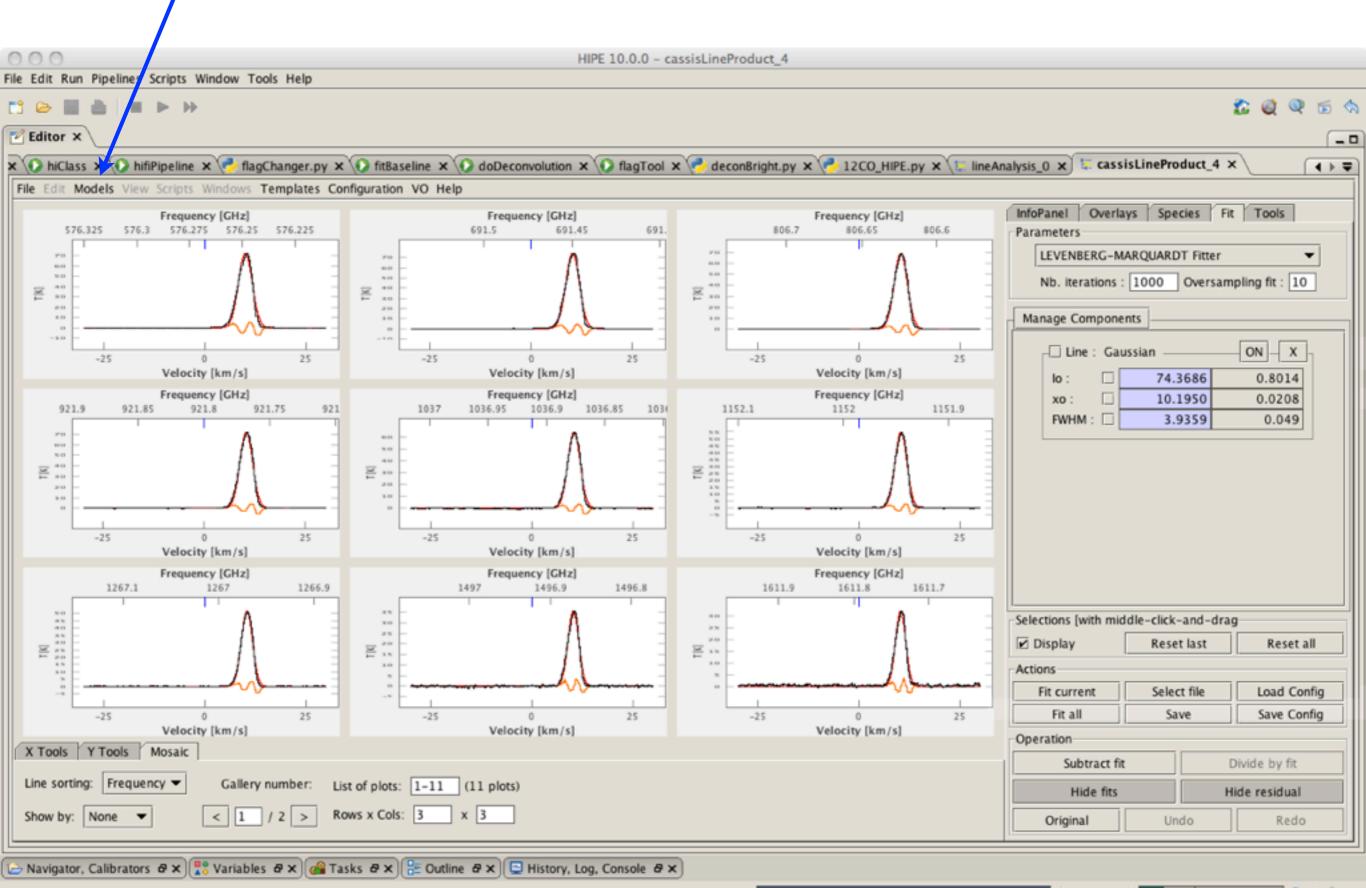
Click on "Create a new file": it will save all the fits (in this case, 1 Gaussian fit per transition, i.e. 11 fits) in the same file



Click on the green arrow to see all the fits

000

### After admiring your beautiful fits, click on "Models", then "Rotational Diagram"

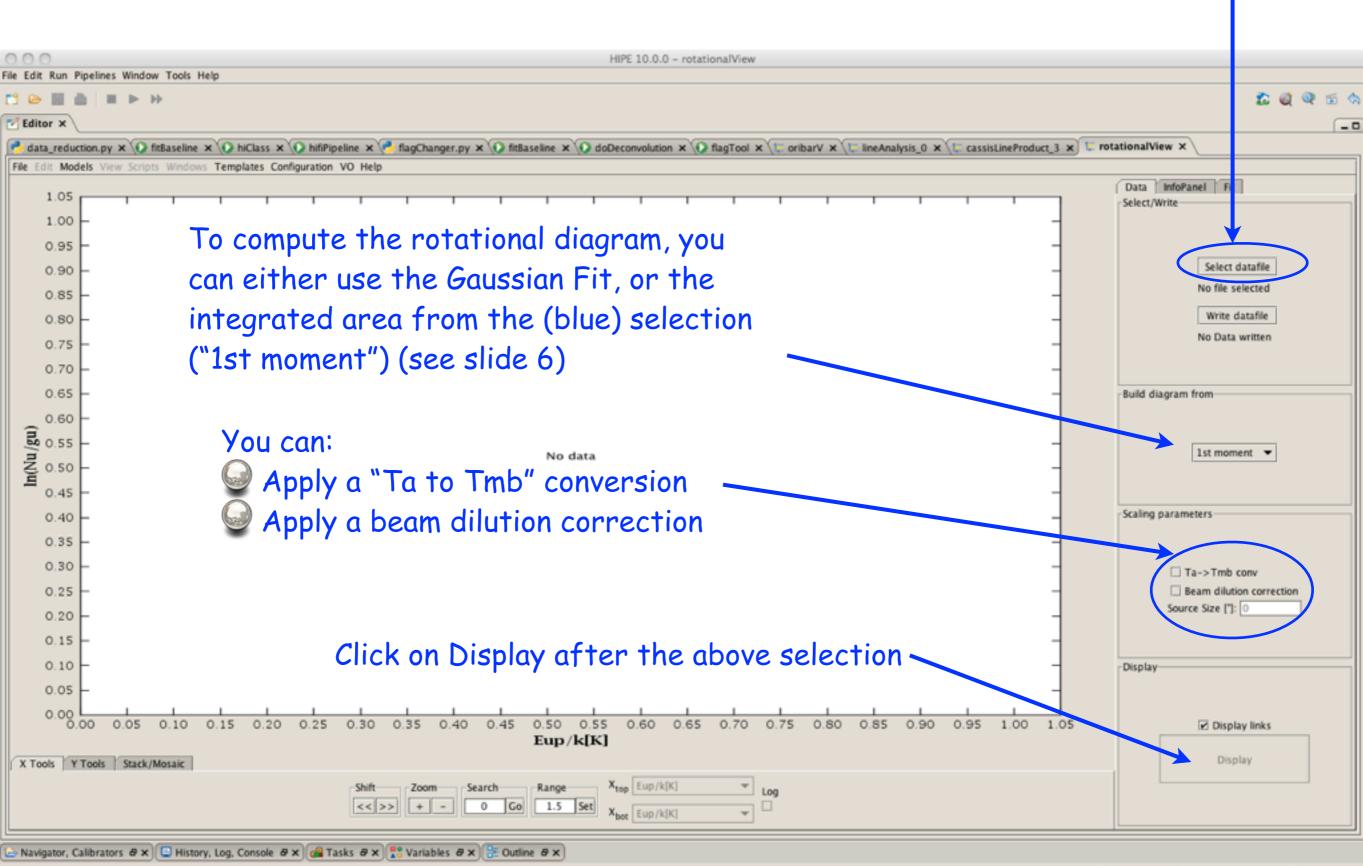


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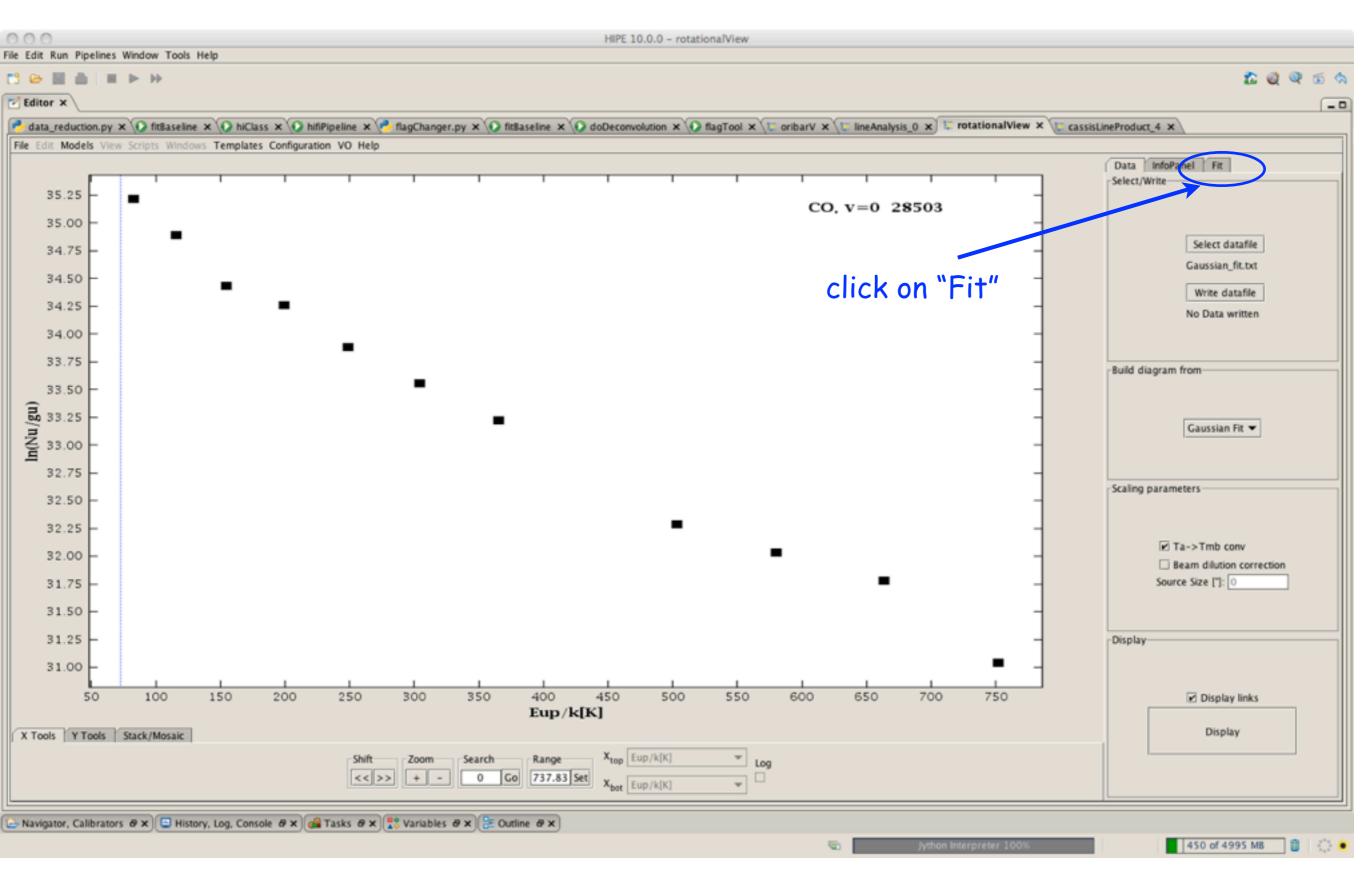
### open the file that you created in Slide 7 (where the fits where saved)

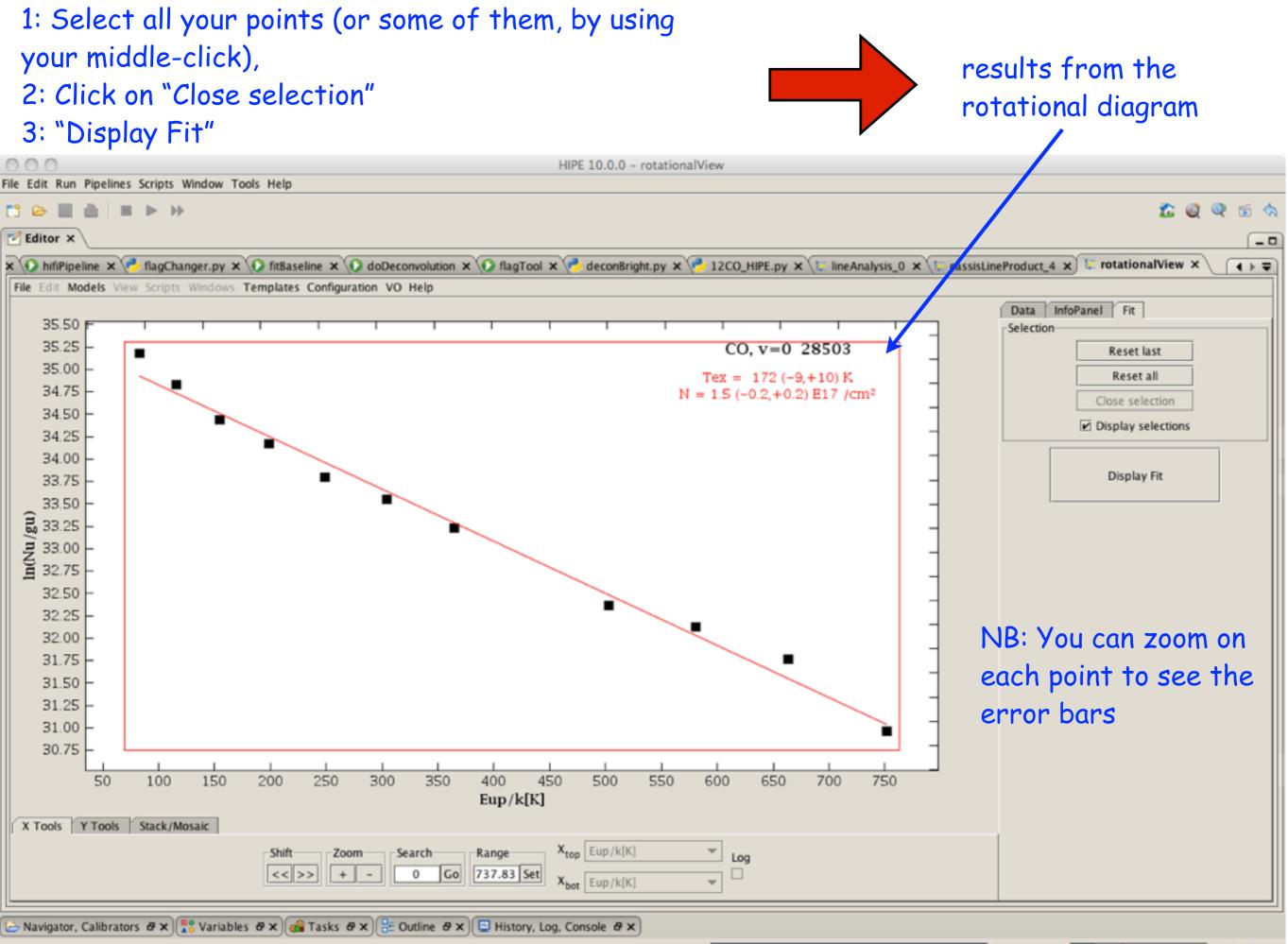


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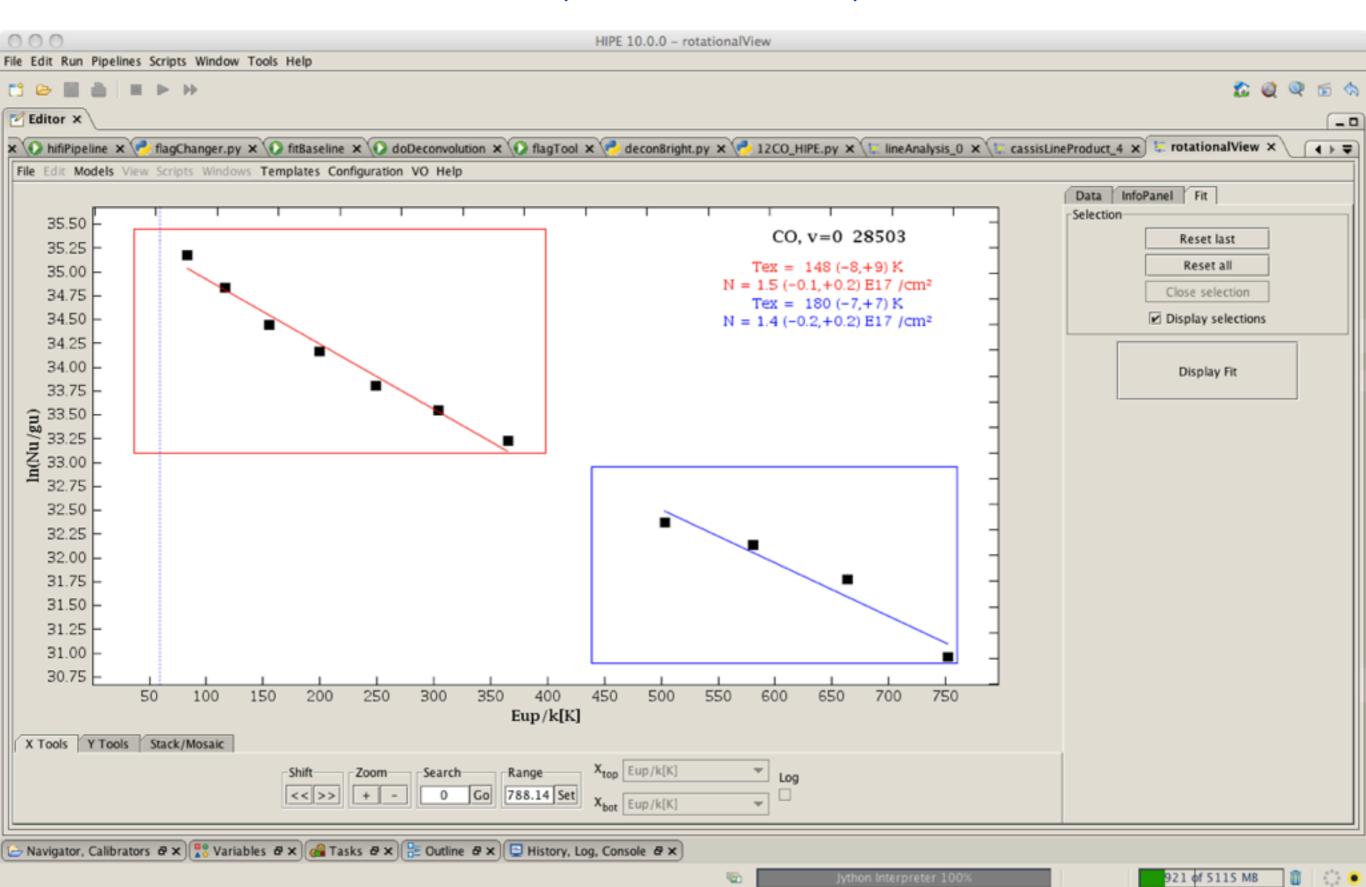
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### Here are the 11 transitions. You can double-click on each transition to see the informations

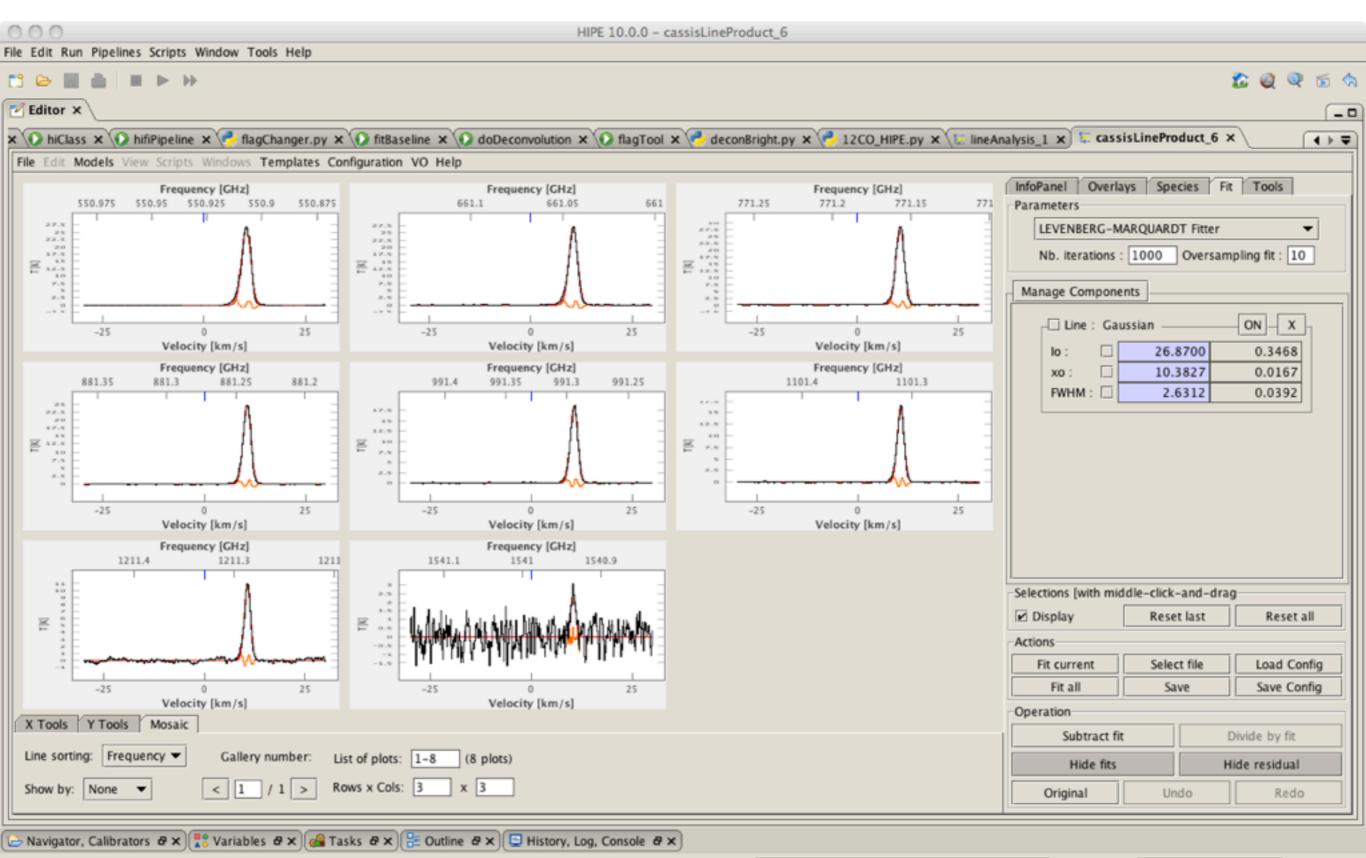




### You can also decide for different components. Below 2 components with different Tex and N.



### Same thing for <sup>13</sup>CO (select Eup lower than 630 K)

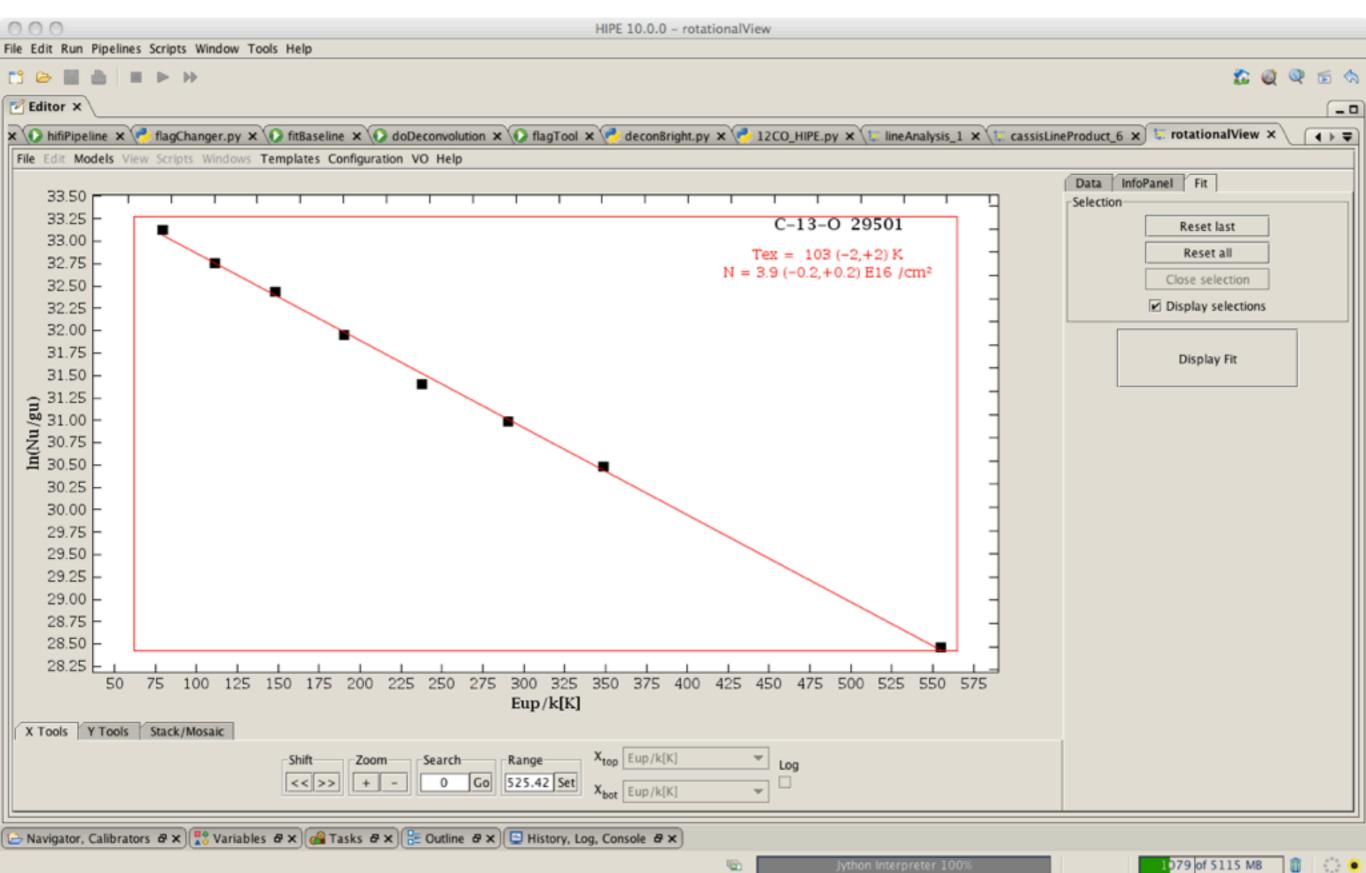


(a)

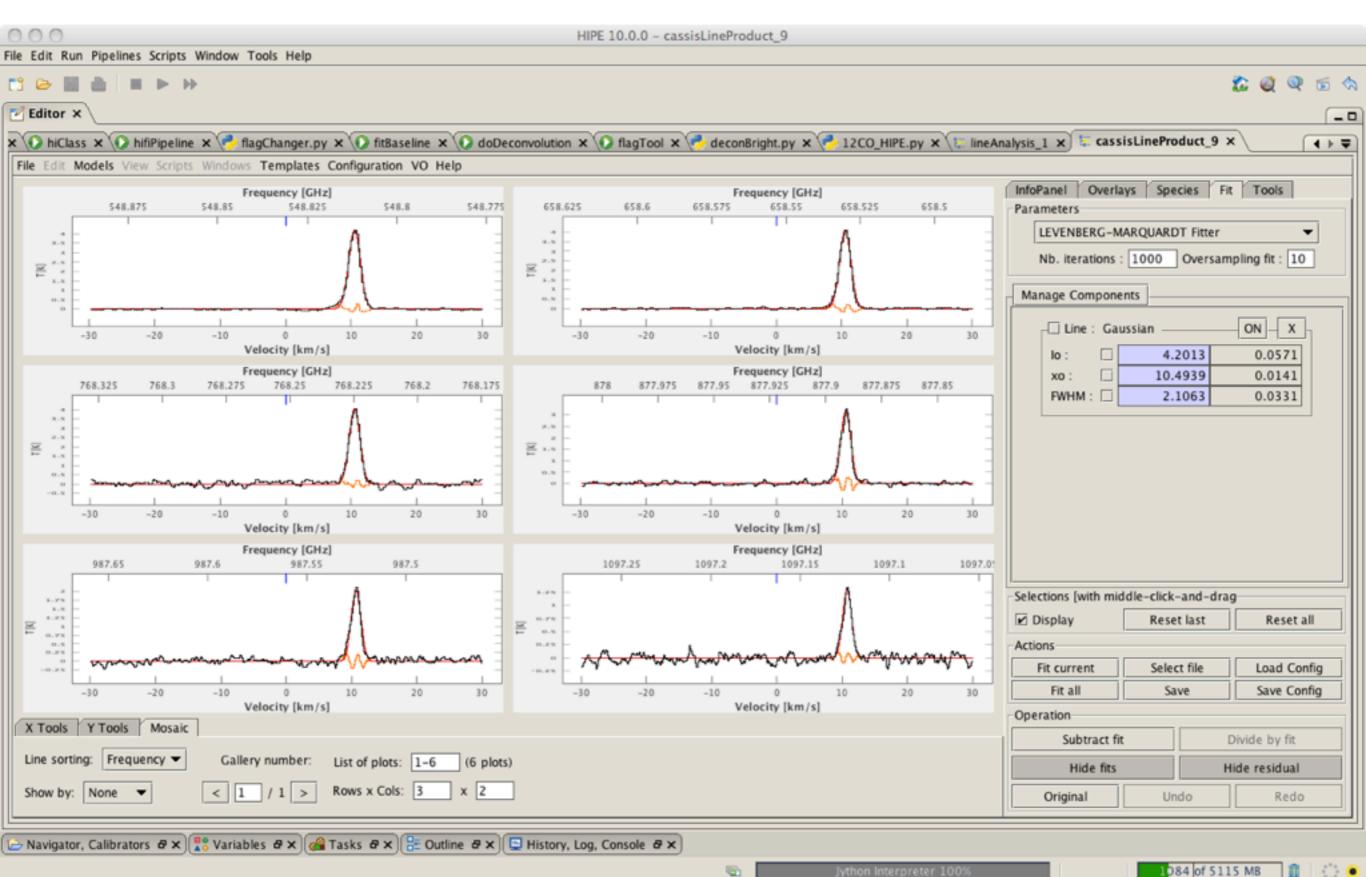
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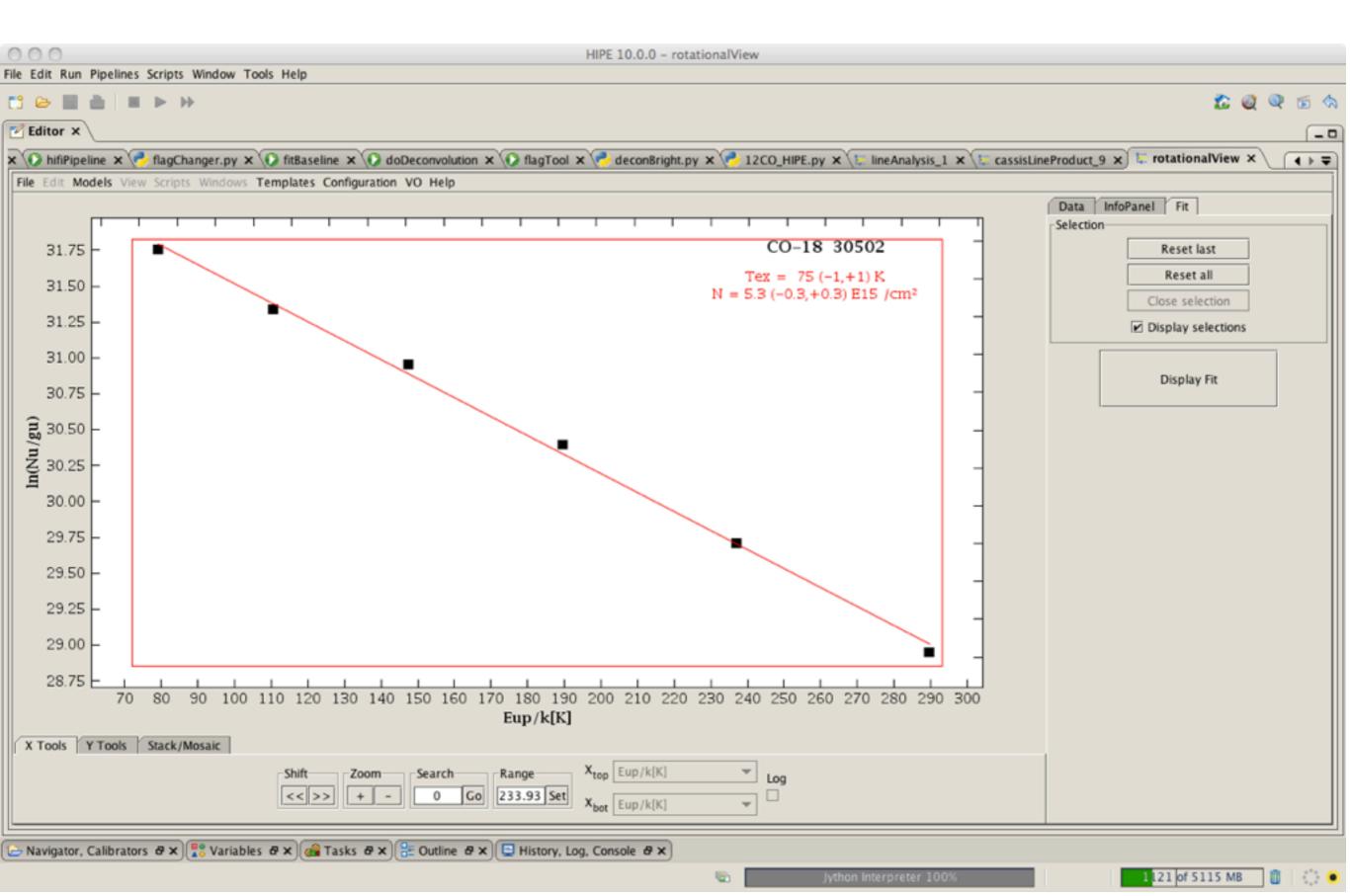
### results for <sup>13</sup>CO



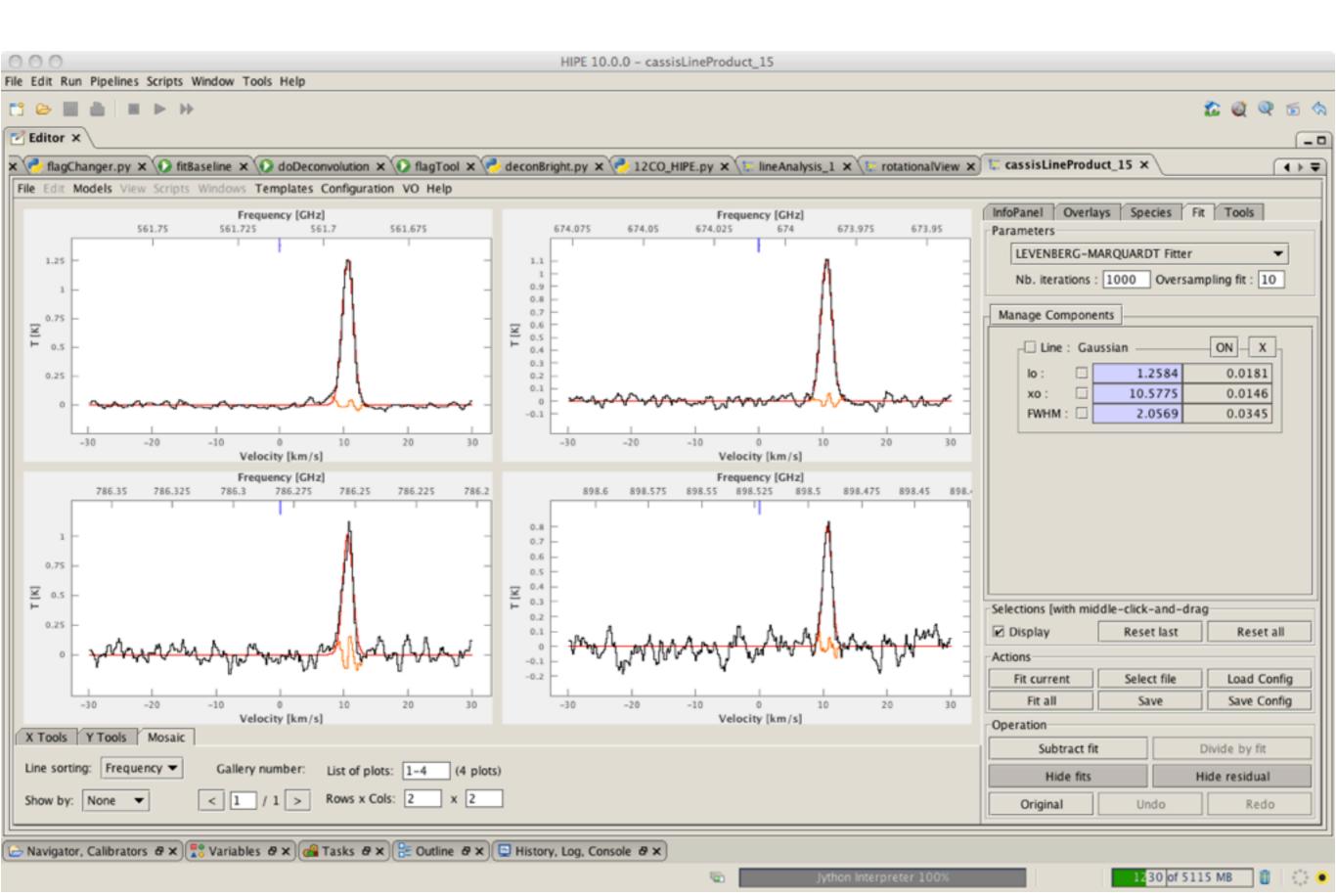
### Same thing for C<sup>18</sup>O (select Eup lower than 300 K)



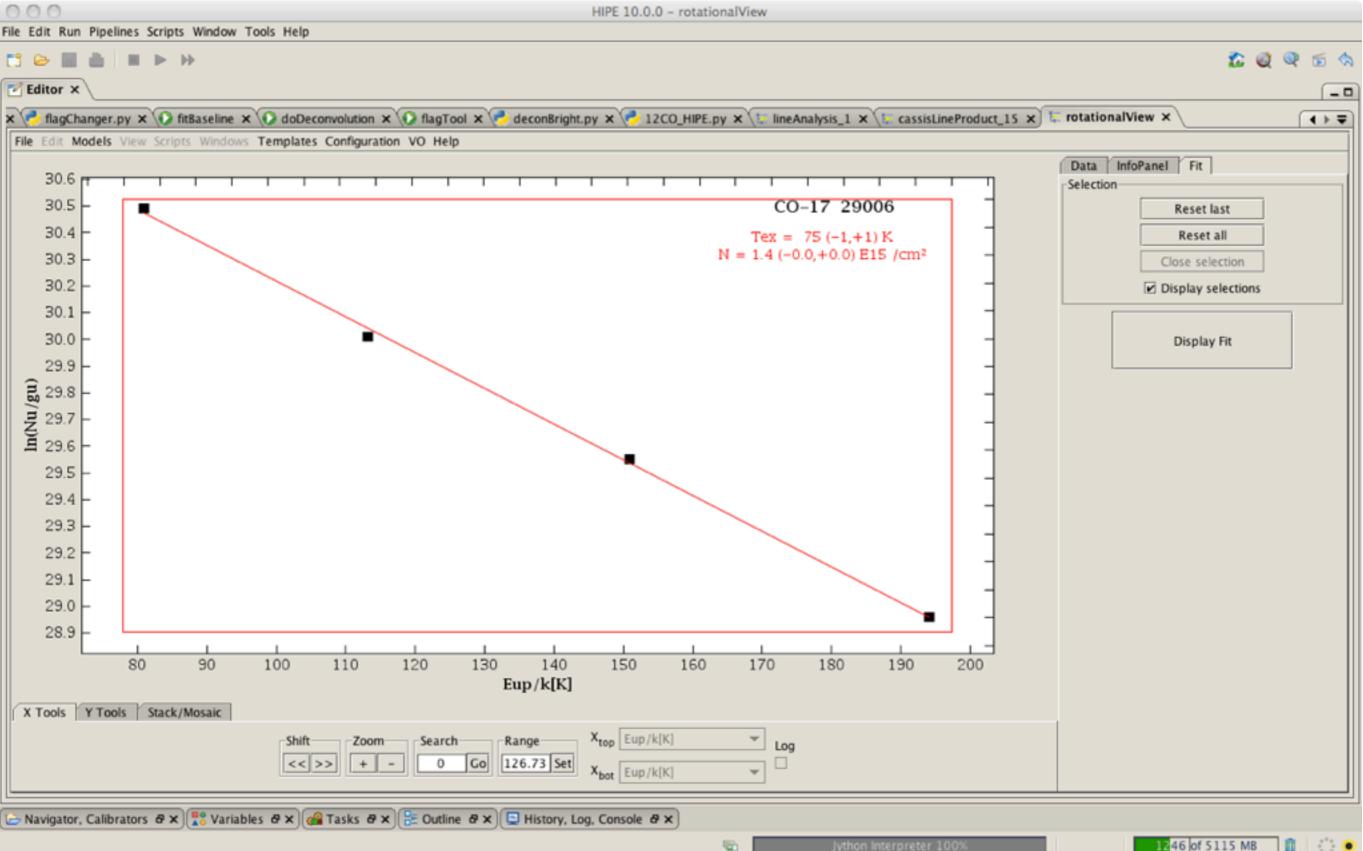
### results for C<sup>18</sup>O



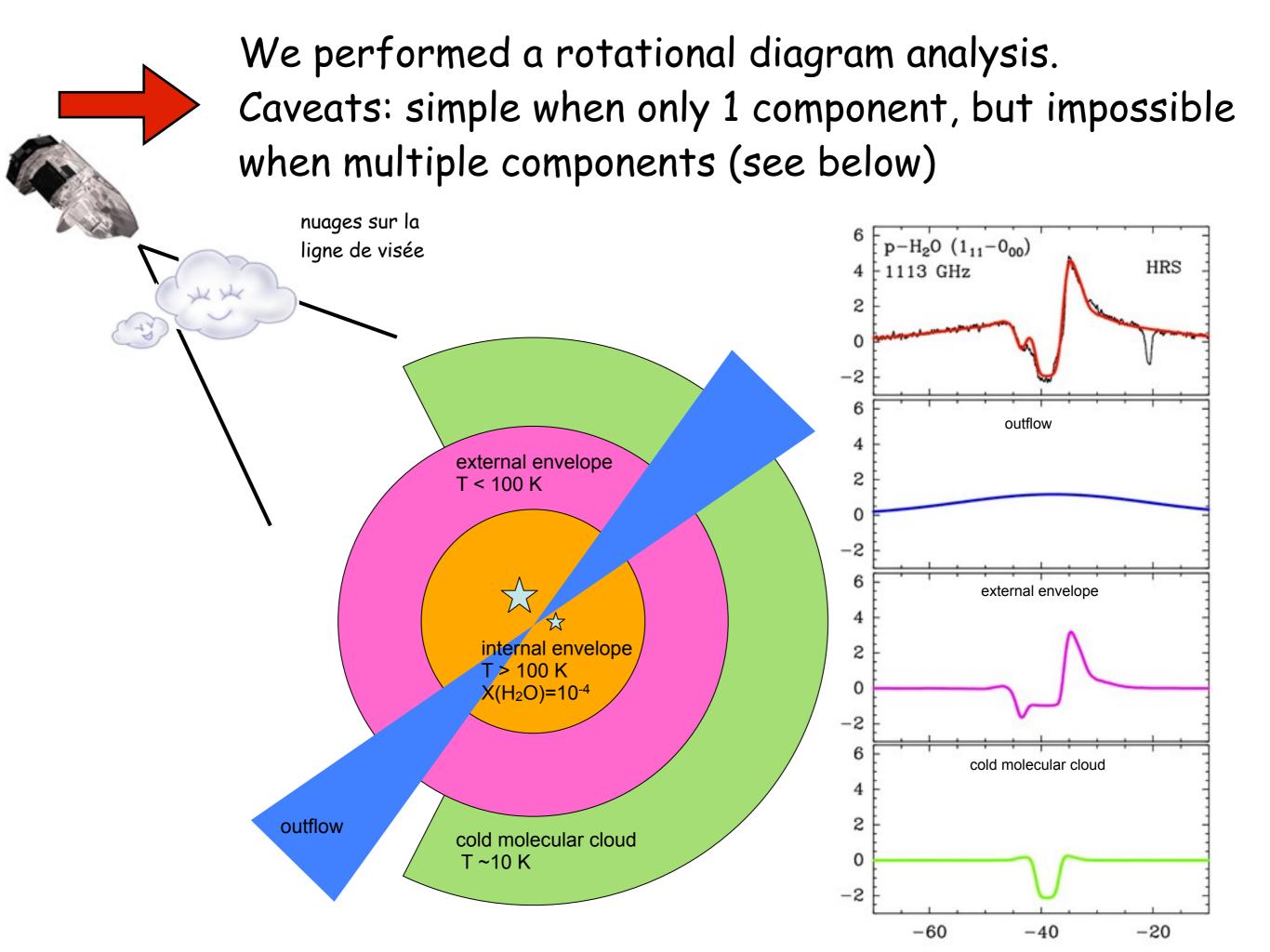
### Same thing for C<sup>17</sup>O (select Eup lower than 300 K), with JPL database (no hyperfine structure)



### results for $C^{17}O$



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Then we might need to disentangle the many components, before doing the rotational diagram fitting for each of the components.

Also, we might need to explore the many parameters needed for complete radiative transfer modeling: excitation temperature, column density, velocity in the local standard of rest, linewidth, beam dilution, interaction with the continuum, (and H<sub>2</sub> density, kinetic temperature for non-LTE modeling). Then you must explore all those parameters (beware that it can be time consuming) within grids of parameters spacing. That can be done within CASSIS with the  $\chi^2$  minimization. The best fit can be found when the  $\chi^2_{red}$  is

 $\chi_i^2 = \sum_{i=1}^N \frac{(I_{obs,j} - I_{model,j})^2}{rms_i^2 + cal_i^2(I_{obs,j} - I_{cont,j})^2}$ 

minimum.

I<sub>obs</sub>: observed intensity within a channel I<sub>model</sub>: modeled intensity within a channel

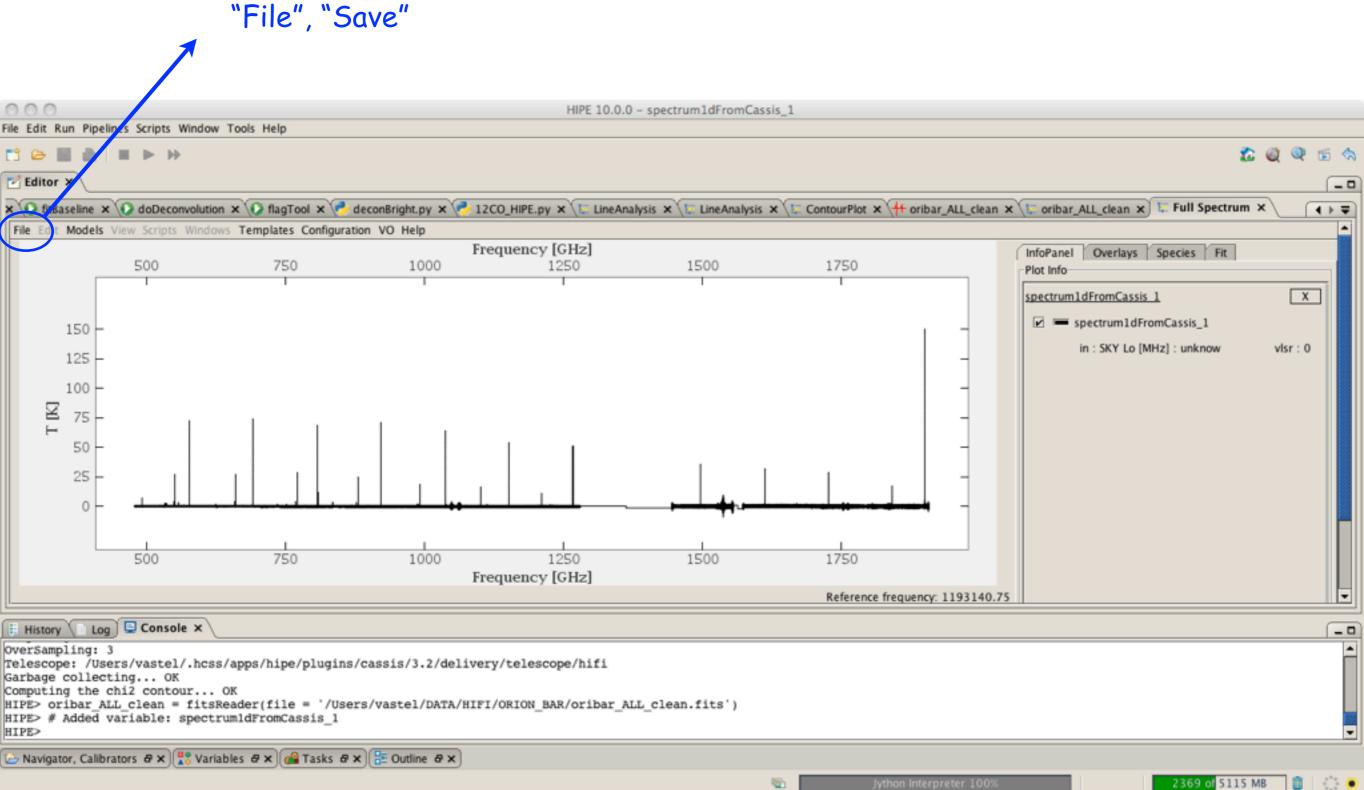
taking into account all spectra

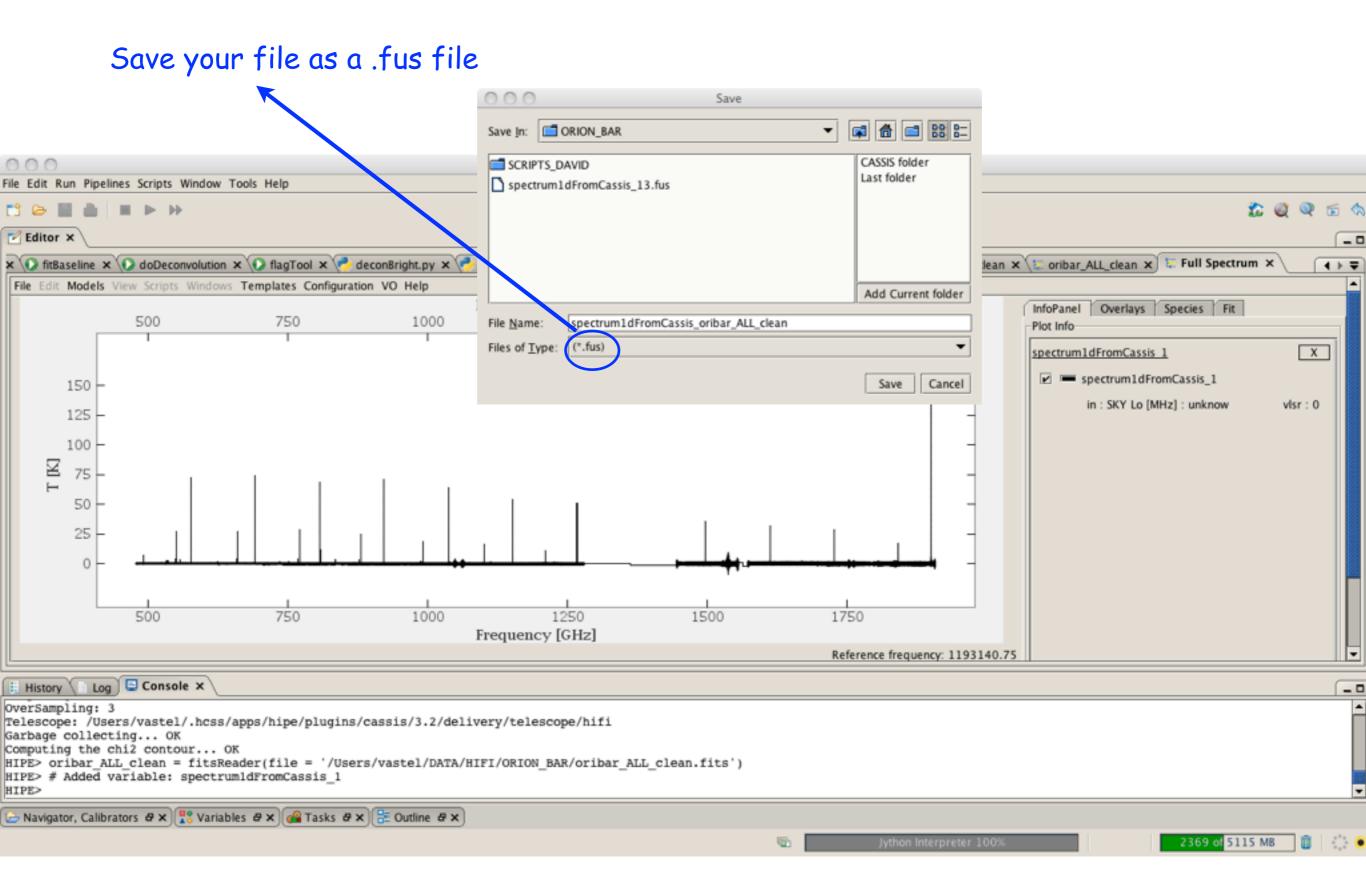
$$\chi^2_{reduit} = \frac{1}{N_{spectre} \sum_{i=1}^{N_{spectre}} w_i} \sum_{i=1}^{N_{spectre}} \frac{\chi^2_i w_i}{\frac{N}{N_{ind}} - dof}$$

N<sub>spectre</sub>: total number of spectra N<sub>ind</sub>: total number of independent points dof: degree of liberty wi: weigth for each spectrum

### Right click on the oribar\_ALL\_clean variable and "open with" "CASSIS Spectrum1dAnalysis"

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🗙 🕢 fitBaseline 🗙 🕡 doDeconvolution 🗙 🕡 flagTool 🗙 🥐 deconBright.py 🗙 🍊 12CO_HIPE.py 🗙 🚺	🗄 LineAnalysis 🗙 🕼 LineAnalysis 🗙 🕼 ContourPlot 🗙 🕂 oribar_ALL_clean 🗙 😳 oribar_ALL_clean 🗙 🌾 Full Spectrum 🗙 🕢 📢 🛡
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Computing the chi2 contour OK HIPE> oribar_ALL_clean = fitsReader(file = '/Users/vastel/DATA/HIFI/ORION_BAR/or	ribar_ALL_clean.fits')
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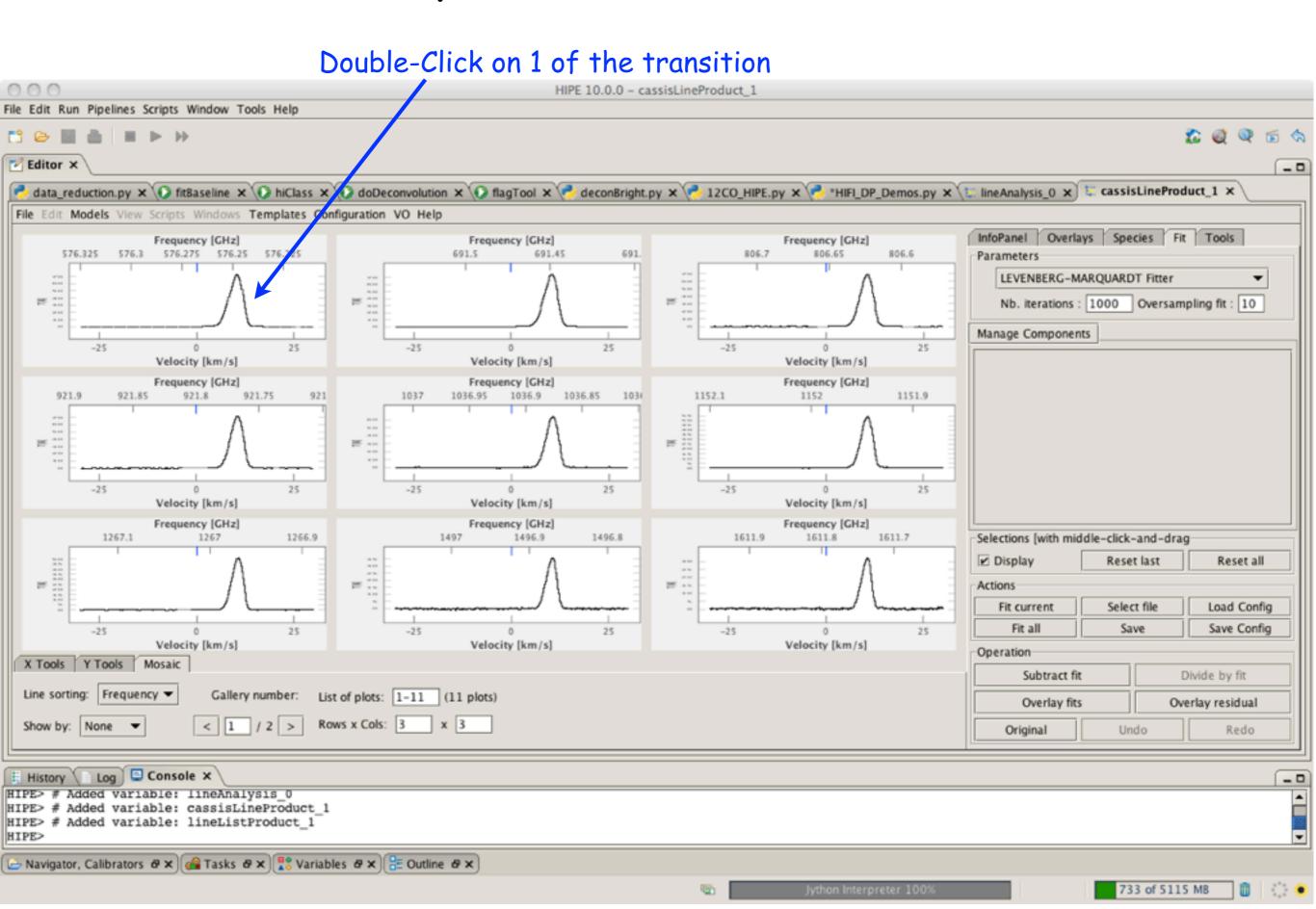


# Open the 12CO\_HIPE.py script

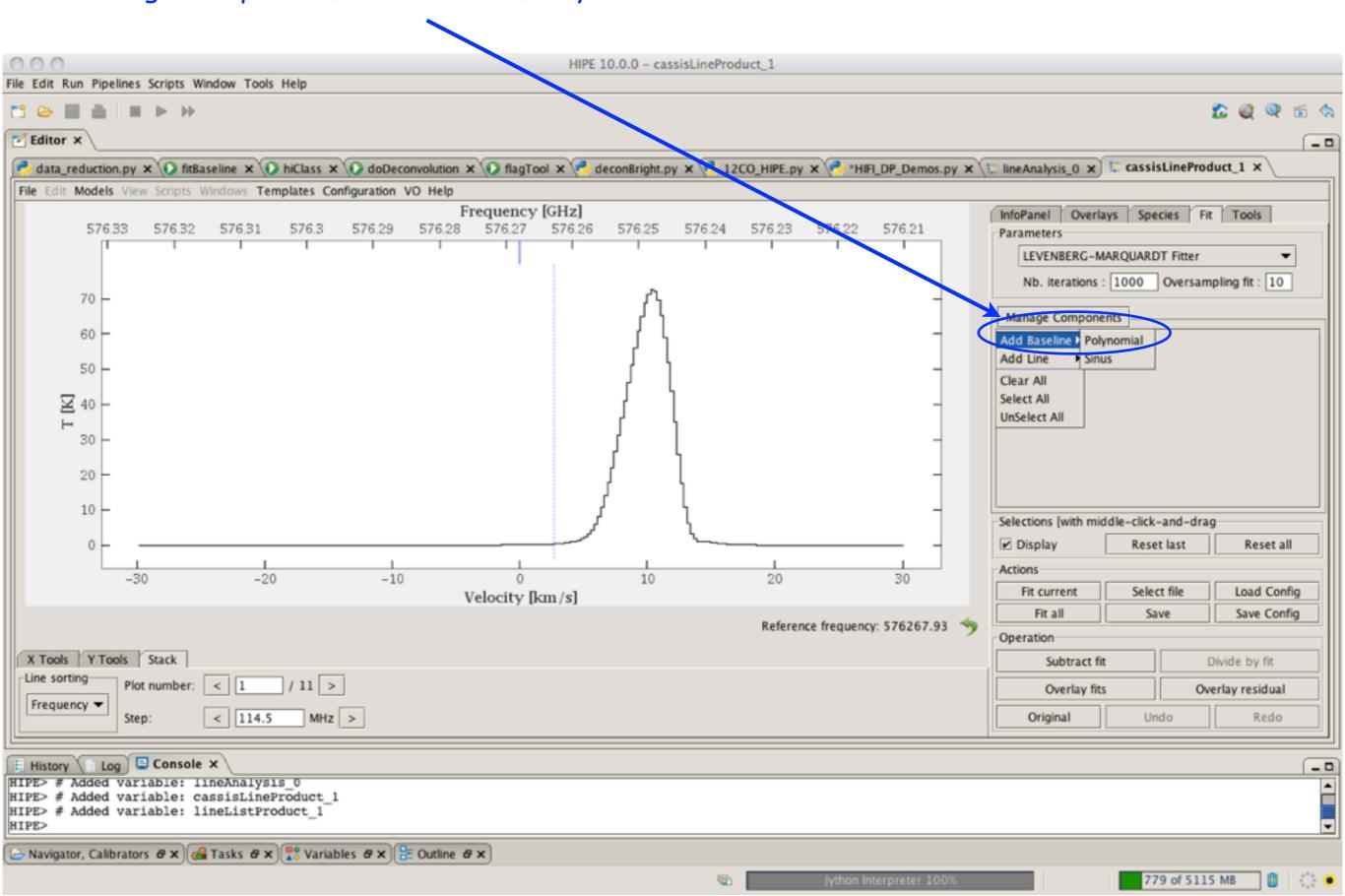
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23456789	<pre>import ScriptEnvironment import time from Component import Component from Range import Range from ScriptLineAnalysisRG import UserInputs # INPUTS # Range.unit = "km/s" # Possible units are GHz and km/s, MHz, cm-</pre>	1 and micrometer to be implemented	
11 12 13	<pre>rl = Range(0.0, 20.0) sourceName = "Orion_Bar" speciesName = "12CO" myDir = "/Users/vastel/DATA/HIFI/ORION_BAR/" myName = sourceName+speciesName</pre>		
17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32	<pre>USER INPUTS USER INPUT USER IN</pre>	telescope/hifi",	
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Total Comput. OverSar Telesco Garbago	<pre>y Log Console x rameters 1: PhysicalModel(VISr = 10.25, tex = 155.556, nmol = 2.22605e+17, Size b of points used in the computation = 1753 / 5244 </pre>	e = 300, fwhm = 3.5, tbg = 2.73, model = Lte, continuum = continuum-0)	
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Interactive part with the user to modify the parameters of the python script

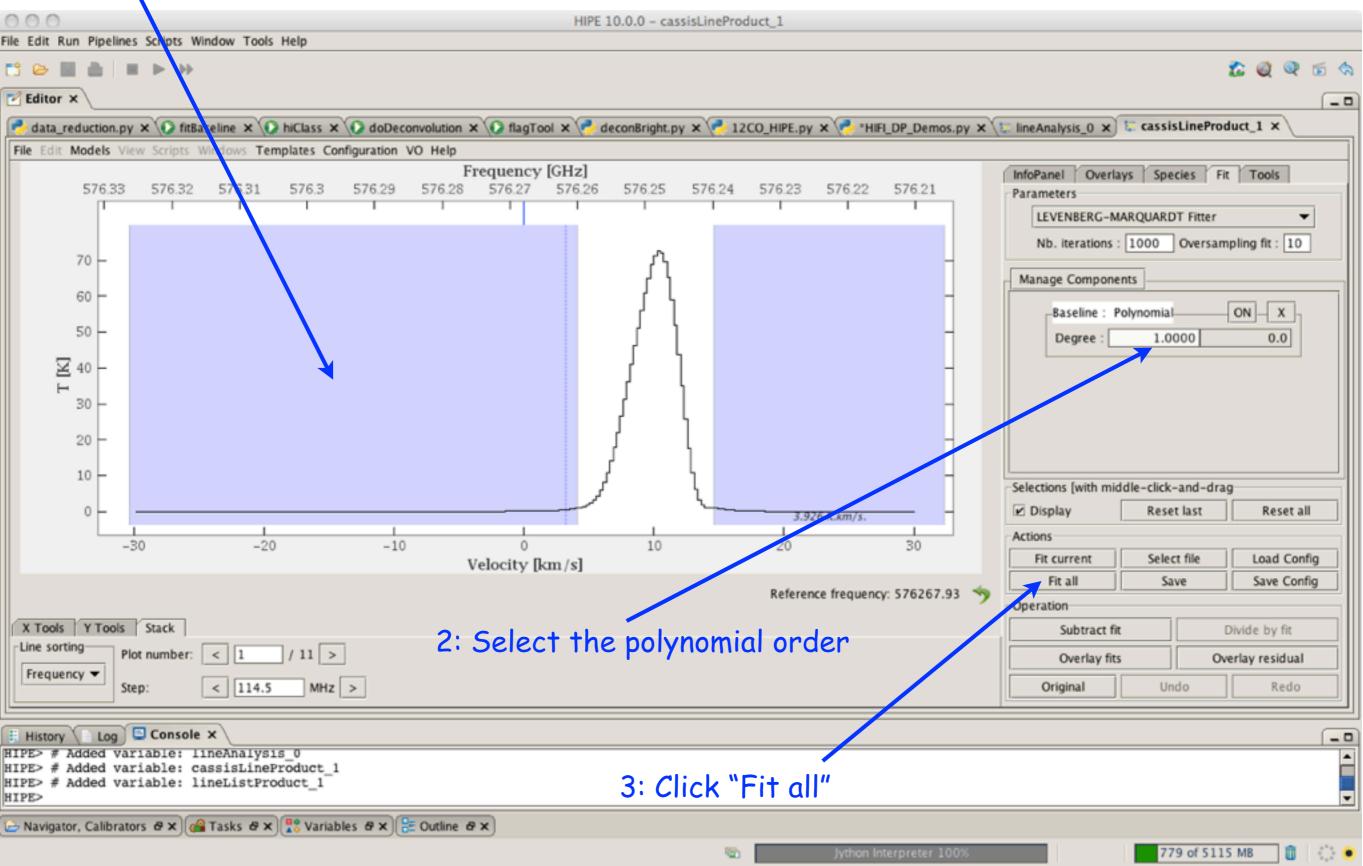
## How to compute the rms with CASSIS?



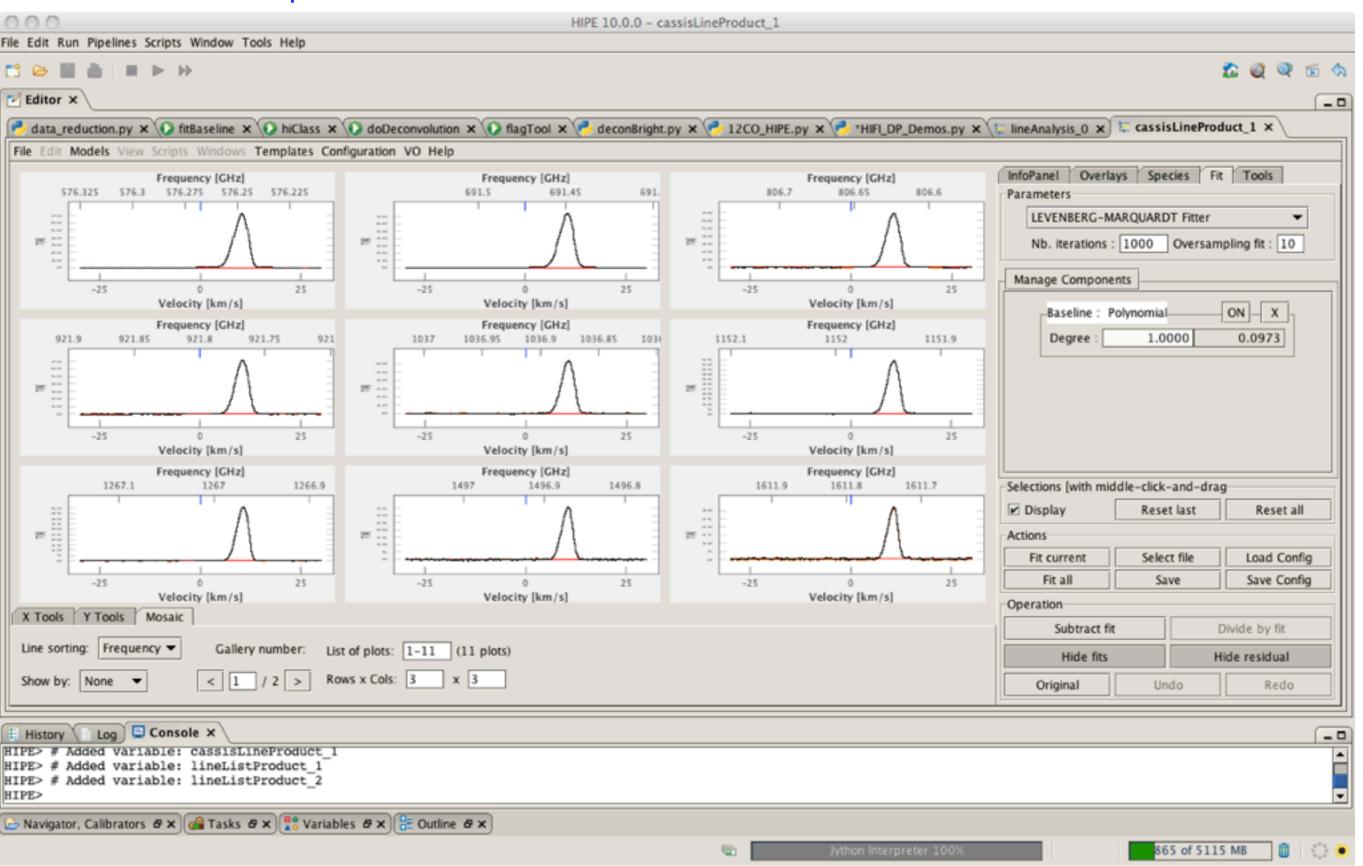
### Manage Component, Add Baseline, Polynomial



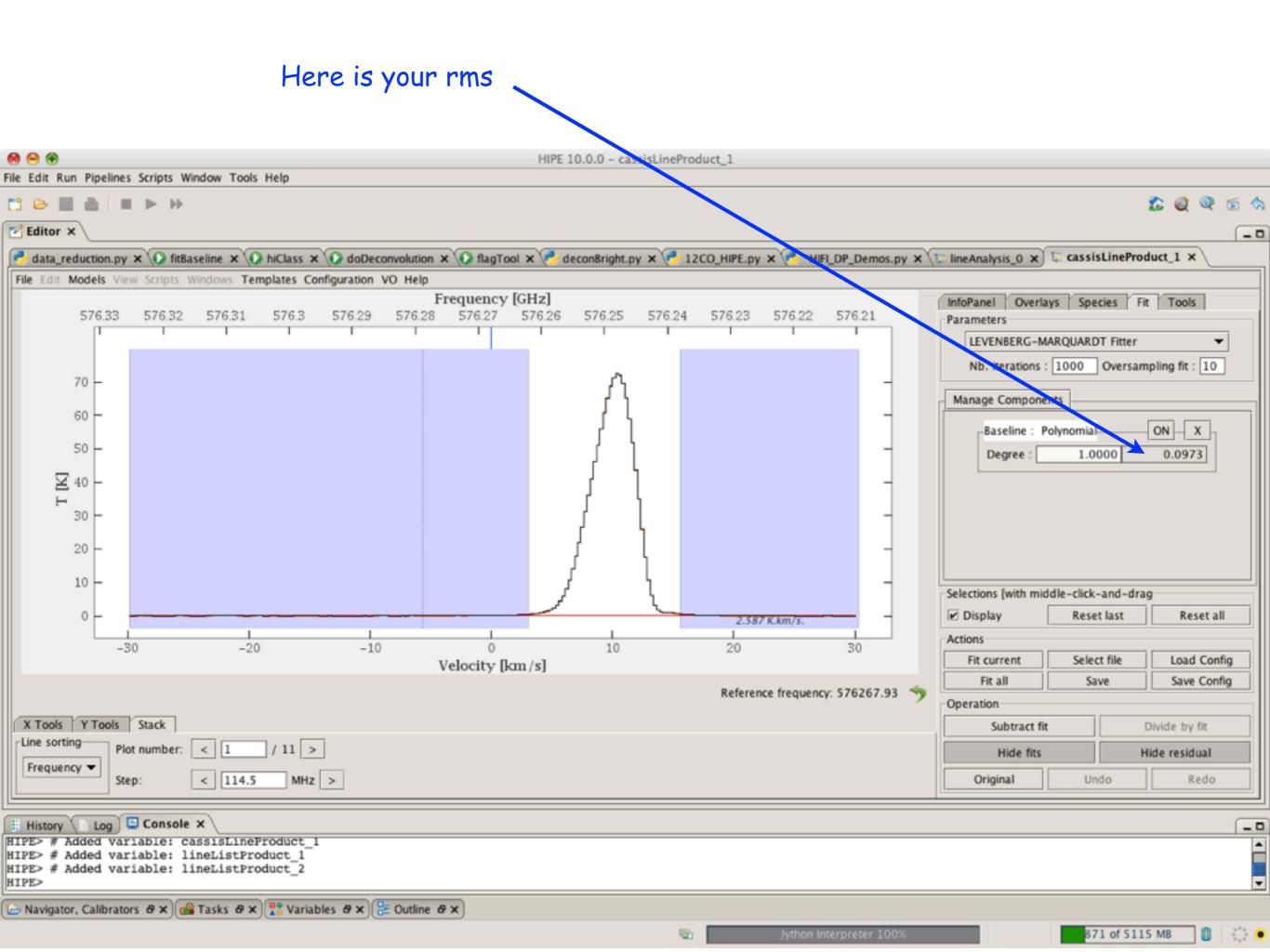
### 1: Use your mouse (middle-click, or alt-left click if you don't have a mouse ) to select the area to be fitted (here in blue). You must avoid all the lines.

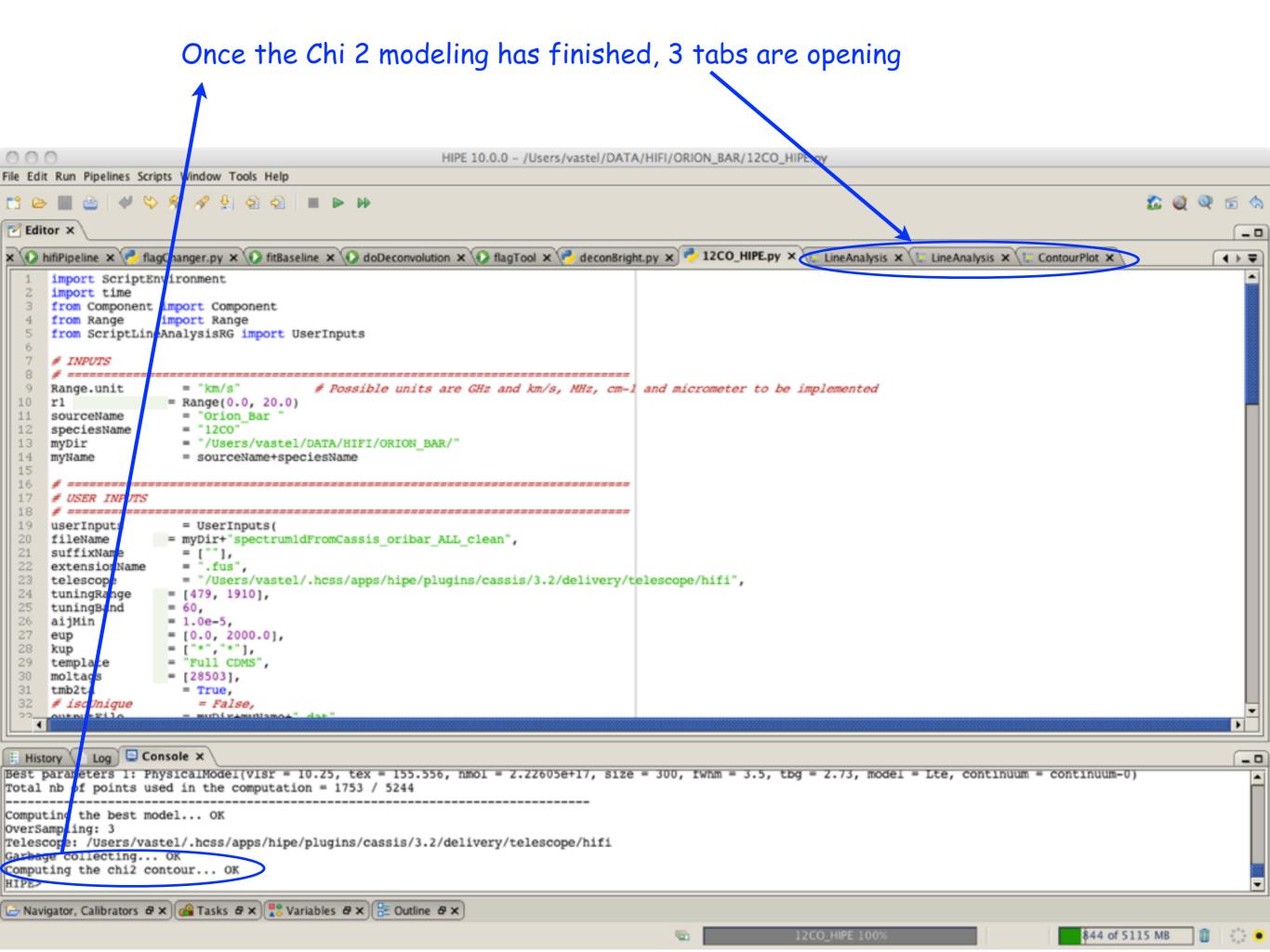


#### Follow the same steps as in Slides 7 and 8.

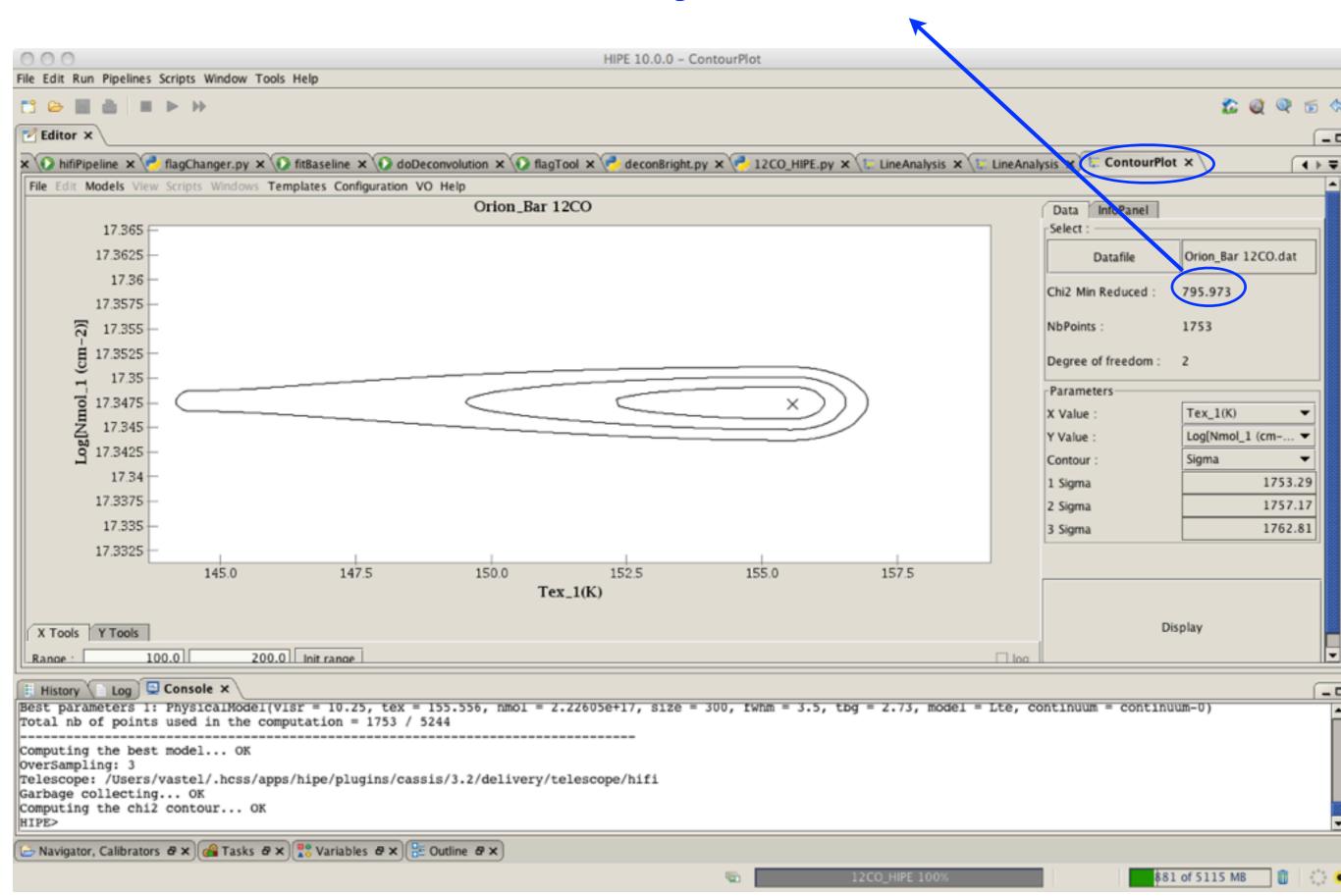


Then, double-click on each spectra and get your rms (see an example in the next slide).





### Not so good but we fixed fwhm and VIsr!!!



# These are the parameters found for the best $\chi^2$ fit, as well as the fixed parameters, that are used for the LTE modeling of the transition.

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	CH 13502
Jup min: * max: * Kup min: * max: * Lup min: * max: * Mup min: * max: *	CH+         13503         Save config           CH+, v=1-0         13504
LTE-RADEX 🗹	
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Component 1 🔟 🔶	
Mode: Full LTE	Tbg [K]: 2.73 N(H <sub>2</sub> ) [cm <sup>-2</sup> ]: 7.5E22 Continuum
Molecules: Operations  Geometry: Slab	V <sub>lsr</sub> : 10.25 km/s   Continuum-0
Species Tag Database Compute N(S	Sp) (cm <sup>-2</sup> ) Abundance (/H2) Tex (K) FWHM (km/s) Size (")
CO, v=0 28503 CDMS	2.23E17 0.001 155.556 3.50 300.00
History Log Console ×	
Best parameters 1: PhysicalModel(VIsr = 10.25, tex = 155.556, nmol = 2.22605e+17, size =	
Total nb of points used in the computation = 1753 / 5244	= 300, fwhm = 3.5, tbg = 2.73, model = Lte, continuum = continuum-0)
Computing the best model OK	
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# These are the modeled transitions, for the best $\chi^2$ fit. Note that we fixed the beam dilution, the FWHM as well as the VIsr.

