

# SYNSPEC

# SYNSPEC

- calculation of synthetic spectra of stellar atmospheres and disks
- input model atmospheres TLUSTY or ATLAS

Hubeny, I., Štefl, S., Harmanec, P., 1985, Bull.  
Astron. Inst. Czechosl. 36, 214

<http://www.physics.muni.cz/~krticka/synspec49.tgz>

# Equation to solve

- radiative transfer equation

$$\mu \frac{\partial}{\partial z} I(z, \mu, \nu) = \eta(z, \nu) - \chi(z, \nu) I(z, \mu, \nu)$$

- fixed structure of atmosphere

# Files with the source code

<http://nova.astro.umd.edu/Synspec49/synspec.html>

- synspec45.f code
- PARAMS.FOR some basic settings
- LINDAT.FOR, MODEL.P.FOR
- rotin3.f code for the rotational convolution

compilation:

ifort [-O3 ] synspec45.f

# Input files

<http://nova.astro.umd.edu/>

- file 1: the same as in TLUSTY
- file 5: the same as in TLUSTY
- files of the model ions (as in TLUSTY)
- nonstandard settings (as in TLUSTY)
- file 55: basic parameters of the calculation
- file 56: change of the chemical composition
- file 19: line list
- file 8: input model atmosphere

# File 55 (fort.55)

```
0 30 1      !IMODE, IDSTD, IPRIN
0 0 0 0      !INPUT MODEL, INTRPL, ICHANGE, ICHEM
0           !IOPHLI
0 0 0 0 1    !IFREQ, INLTE, ICNTL, INORIG, IFHE2
0 0 0 0 0 0 0 !ILYCS, IBVCS, IHE1, IHE4471, IHE2UV, I
4000 7000 10 10 1d-6 0.001 ! ALAM0, ALAST, CUTOF0,
                               CUTOFFS, RELOP, SPACE
2.0          ! turbulent velocity (in km/s)
```

# Input parameters

IMODE: mode of calculation

IDSTD: index of the standard depth (unity optical depth)

IPRIN: amount of output

INPUT MODEL: (0=ATLAS, 1=TLUSTY)

IOPHLI: treatment of hydrogen lines

INLTE: treatment of NLTE

ALAM0, ALAST: wavelength limits

RELOP: minimum relative contribution of lines

VTB: turbulent velocity in km/s

# Example tun

```
/home/krticka/prakaf/synspec/synspec  
file syn:  
DOMA=/home/krticka  
JOBID=$1  
MODEL=$2  
LINES=$3  
ln -s $MODEL.7 fort.8  
ln -s $JOBID.55 fort.55  
ln -s $DOMA/GF/$LINES fort.19  
$DOMA/prakaf/synspec/synspec/synspec45 <$MODEL.5  
    >! $JOBID.6  
cp fort.7 $JOBID.7  
cp fort.17 $JOBID.17  
cp fort.12 $JOBID.12  
cp fort.18 $JOBID.18  
cp fort.13 $JOBID.13  
cp fort.10 $JOBID.10  
rm fort.*
```

# Example run

```
/home/krticka/prakaf/synspec/synspec  
file syn:  
run syn slun.s slun gf99.dat  
input files slun.5, slun.7, slun.s.55,  
gf99.dat  
output files slun.s.6, slun.s.7,  
slun.s.17 ...
```

# Output files

- file 6: basic model description
- file 7: emergent flux in lines
- file 17: emergent flux in continuum
- file 12: line identification (except H and He II)

CGS!

# File 6 (fort.6)

- general model output
- tables with input data
- error messages

# File 7 (fort.7)

- wavelength-dependent emergent flux  
 $(\lambda - \text{\AA}, H_\lambda - \text{erg cm}^{-2} \text{s}^{-1} \text{\AA}^{-1})$

# File 17 (fort.17)

- wavelength-dependent emergent flux in continuum  
 $(\lambda - \text{\AA}, H_\lambda - \text{erg cm}^{-2} \text{s}^{-1} \text{\AA}^{-1})$

# File 12 (fort.12)

- line identification table, *estimate* of the equivalent width, line strength (\*)

# Rotational convolution

- code rotin3.f
- input files
  - file 5: parameters of the calculation
  - file 7, file 17 (SYNSPEC output)
- output file
  - file 11: convolved spectrum  
(rotation+instrumental)

# Input file 5

```
'fort.7' 'fort.17' 'fort.11' ! filenames for the  
70.0 0.00 0.001 ! vrot, chard, step  
0.03001 0.001 ! fwhm (instrumental)  
4100 6900 1 ! lam0, lam1, irel
```

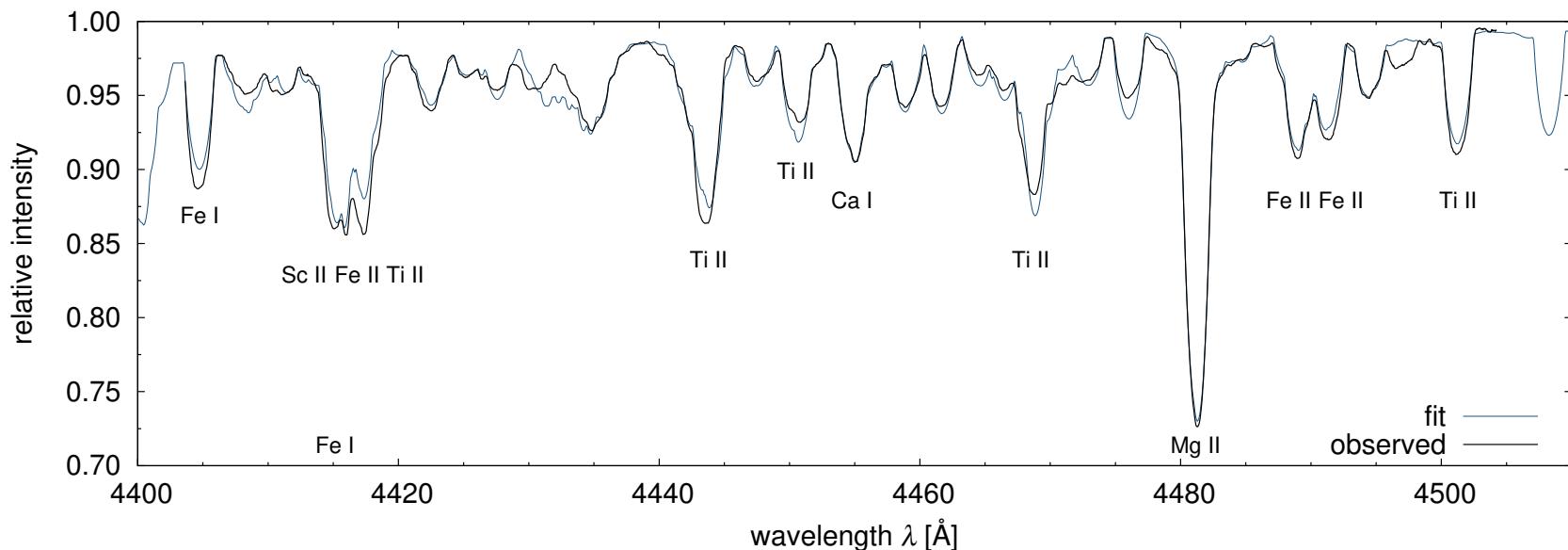
- names of input and output files
- $v_{\text{rot}} \sin i$  in km/s, 0., step
- FWHM of instrumental profile
- wavelength interval, absolute/relative spectrum

# Example run

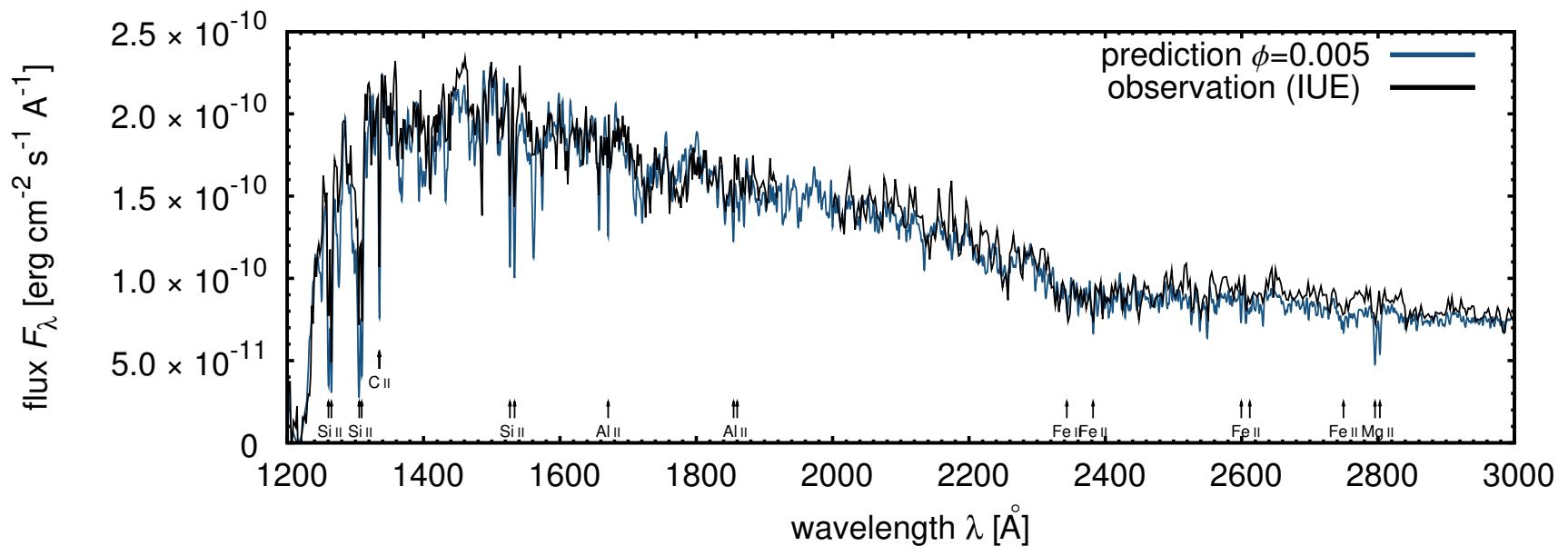
**file rot:**

```
JOBID=$1
SPEK=$2
ln -s ${SPEK}.7 fort.7
ln -s ${SPEK}.17 fort.17
rotin3 <${JOBID}.5
rm fort.7 fort.17
mv fort.11 ${SPEK}.11
execute: rot rvs b.s
```

# Example of visual spectrum



# Example of UV spectrum



# Abundance analysis

minimalization

$$\chi^2 = \sum_{\lambda_i} \frac{(I_i^{\text{obs}} - I_i^{\text{syn}})^2}{\sigma_i^2}$$