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---

## **agn\_blr\_albedo *measure rayleigh scattering of Lya***

```

title measure rayleigh scattering of Lya
c model from Korista, K., & Ferland, G. 1998, ApJ, 495, 672
c
c commands controlling continuum =====
agn 6.683 -1.20 -1.20 -0.90
ionization parameter 1.0
c
c commands for density & abundances =====
init file = "ism.ini"
abundances old solar 84
hden 11.0
c
c commands controlling geometry =====
stop total column density = 23.75
c
c other commands for details =====
iterate
c
c commands controlling output =====
normalize to "FeKa" 1.78
print line faint -1
print diffuse continua
print lines inward
punch continuum last "agn_blr_albedo.con" units kev
punch dr "agn_blr_albedo.dr"
punch reflected continuum last "agn_blr_albedo.ref"
punch emitted continuum last "agn_blr_albedo.emt"
c
c must assert the luminosity since change in this
c line would appear to change the rest of the spectrum
c agn_blr_albedo.in
c class blr
c =====

```

This model computes the albedo of a fairly standard BLR cloud. This is the type of model that was presented in the BLR albedo paper by Korista & Ferland, 1998, ApJ 495, 672.

The print diffuse continua command enters continuum fluxes into the emission line stack. The asserts then check that these continua have the expected brightness.

---

## **agn\_lex00\_u0 *intermediate-ionization x-ray ionized cloud from Lexington 2000***

```

title intermediate-ionization x-ray ionized cloud from Lexington 2000
stop column density 16
print line sort wavelength
print lines column linear
normalise to "H 1" 1216
print line faint -1

```

```

hden 5
iterate
phi(h) 15.477 range 7.353 to 735.3
interpolate (-8 -3)
continue (-6 1.)
continue (-2 14.5)
continue (0.4771 12.7)
continue (0.8663 10.6)
continue (3.8663 7.6)
continue (6 1)
continue (9 -3)
element abundance helium -1
element abundance carbon -3.432
element abundance nitrogen -3.959
element abundance oxygen -3.097
element abundance neon -3.959
element abundance magnesium -4.4318
element abundance silicon -4.456
element abundance sulphur -4.7959
element abundance argon -5.4318
element abundance iron -4.398
element lithium off
element beryllium off
element boron off
element fluorine off
element sodium off
element aluminium off
element phosphrous off
element chlorine off
element potassium off
element calcium off
element scandium off
element titanium off
element vanadium off
element chromium off
element manganese off
element cobalt off
element nickel off
element copper off
element zinc off
c
punch transmitted continuum "agn_lex00_u0.trn" units keV
c
c agn_lex00_u0.in

```

This is one of the "warm absorber" simulations presented at the Lexington 2000 meeting on nebulae. Pequignot et al. summarized in 2001ASPC..247..533P. It is necessary to also include the command no induced processes to obtain the results presented there. This disables UTA ionization, a process that was not included in the calculations presented in the paper.

---

## **agn\_lex00\_u1 *high-ionization x-ray ionized cloud from Lexington 2000***

```

title high-ionization x-ray ionized cloud from Lexington 2000
stop column density 16
print lines column linear
normalise to "H 1" 1216
print line faint -1
hden 5

```

```

iterate
phi(h) 16.477 range 7.353 to 735.3
interpolate (-8 -3)
continue (-6 1.)
continue (-2 14.5)
continue (0.4771 12.7)
continue (0.8663 10.6)
continue (3.8663 7.6)
continue (6 1)
continue (9 -3)
element abundance helium -1
element abundance carbon -3.432
element abundance nitrogen -3.959
element abundance oxygen -3.097
element abundance neon -3.959
element abundance magnesium -4.4318
element abundance silicon -4.456
element abundance sulphur -4.7959
element abundance argon -5.4318
element abundance iron -4.398
element lithium off
element beryllium off
element boron off
element fluorine off
element sodium off
element aluminium off
element phosphorous off
element chlorine off
element potassium off
element calcium off
element scandium off
element titanium off
element vanadium off
element chromium off
element manganese off
element cobalt off
element nickel off
element copper off
element zinc off
punch transmitted continuum "agn_lex00_u1.trn" units keV
c
c agn_lex00_u1.in

```

This is one of the "warm absorber" simulations presented at the Lexington 2000 meeting on nebulae. Pequignot et al. summarized in 2001ASPC..247..533P. It is necessary to also include the command no induced processes to obtain the results presented there. This disables UTA ionization, a process that was not included in the calculations presented in the paper.

---

## **agn\_lex00\_um1 *low-ionization x-ray ionized cloud from Lexington 2000***

```

title low-ionization x-ray ionized cloud from Lexington 2000
stop column density 16
print lines column linear
print lines sort intensity
normalise to "H 1" 1216
print line faint -2
hden 5
iterate

```

```

phi(h) 14.477 range 7.353 to 735.3
interpolate (-8 -3)
continue (-6 1.)
continue (-2 14.5)
continue (0.4771 12.7)
continue (0.8663 10.6)
continue (3.8663 7.6)
continue (6 1)
continue (9 -3)
element abundance helium -1
element abundance carbon -3.432
element abundance nitrogen -3.959
element abundance oxygen -3.097
element abundance neon -3.959
element abundance magnesium -4.4318
element abundance silicon -4.456
element abundance sulphur -4.7959
element abundance argon -5.4318
element abundance iron -4.398
element lithium off
element beryllium off
element boron off
element fluorine off
element sodium off
element aluminium off
element phosphorous off
element chlorine off
element potassium off
element calcium off
element scandium off
element titanium off
element vanadium off
element chromium off
element manganese off
element cobalt off
element nickel off
element copper off
element zinc off
c
punch transmitted continuum "agn_lex00_um1.trn" units keV
c
c agn_lex00_um1.in

```

This is one of the "warm absorber" simulations presented at the Lexington 2000 meeting on nebulae. Pequignot et al. summarized in 2001ASPC..247..533P. It is necessary to also include the command no induced processes to obtain the results presented there. This disables UTA ionization, a process that was not included in the calculations presented in the paper.

---

## **agn\_reflector *model of Compton reflector***

```

title model of Compton reflector
c
c commands controlling continuum =====
c this is a powerlaw that extends into the x-rays, but by default
c has exponential cutoff in FIR
power law -1. hi=1000000000
c the ionization parameter for this powerlaw
ionization parameter 1.
c add background so that FIR/radio does not have zero flux
background

```

```

iterate
c
c commands for density & abundances =====
hden 11
c
c commands controlling geometry =====
stop total column density 24
c
c other commands for details =====
constant temper 5
atom h-like element hydrogen levels resolved 10
atom h-like element helium levels resolved 10
c
c commands controlling output =====
print line faint 1
print line inward
print diffuse continuum
c
c this is to produce lines with lots of contrast with the continuum
set PunchLWidth 100 km/sec
c a fig in part 3 of hazy uses data from this model. the fig shows lines
c with above and below linewidth
c set PunchLWidth c
punch continuum last "agn_reflector.con" units kev
punch continuum reflected last "agn_reflector.alb" units kev
punch overview last "agn_reflector.ovr"
punch dr last "agn_reflector.dr"
c
table lines "LineList_BLR.dat"
c agn_reflector.in
c class blr
c =====
c

```

This is a model of the Compton reflector in AGN. It is a constant temperature since models of this region often make that assumption. A plot in Part I of Hazy shows the incident and reflected portions of the continuum. The code will complain that the cloud is Compton thick since it is not really designed to simulate this situation.

---

## **agn\_s\_curve\_grid *temperature across Spitzer thermal stability S curve***

```

title temperature across Spitzer thermal stability S curve
c
c commands controlling continuum =====
c this is the generic AGN continuum
table agn
c the ionization parameter will be varied
ionization parameter 0.00 vary
grid from 0 to 3 in 0.25 dex steps
c add background so that FIR/radio does not have zero flux
background
iterate
c
c commands for density & abundances =====
c intended to make sim run as fast as possible
hden 8
init file "ism.ini"
element sulphur off
element chlorine off
element argon off

```

```

c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c commands controlling output
c this produces a very large number of lines in the output mostly
c because Hbeta is very weak in hot gas. Choose Lya as the normalization
c line to make line list smaller
normalize to "totl" 1216
print line faint 10 will be varied
c create file with list of grid parameters
punch grid "agn_S_curve_grid.grd" last
c save average H+ temperature
punch average "agn_S_curve_grid.avr" last
temperature hydrogen 1
end of temperatures
c save heating and cooling
punch heating "agn_S_curve_grid.het" last
punch cooling "agn_S_curve_grid.col" last
punch iron element "agn_S_curve_grid.fe" last
punch xspec atable "agn_S_curve_grid.fit"
c
c
c agn_S_grid.in
c class BLR
c =====
c

```

This computes a series of models that check the temperature through the S curve in the Fields et al. three-phase model of ISM stability.

---

## **agn\_warm\_absorber** *simple warm absorber model*

```

title simple warm absorber model
c
c commands controlling continuum =====
table power law
ionization parameter 0
c
c commands for density & abundances =====
hden 6
abundances old solar 84
c
c commands controlling geometry =====
stop column density 22
c
c other commands for details =====
iterate to convergence
c
c commands controlling output =====
print line faint 10
plot continuum range 1 1000
punch continuum "agn_warm_absorber.con" last units kev
punch continuum "agn_warm_absorber.conA" last units Angstroms
punch fine continuum "agn_warm_absorber.finA" last units Angstroms range 61 120 merge 3
punch dr "agn_warm_absorber.dr" last
c this will be the default units, like the printout
punch line optical depths last "agn_warm_absorber_lab.lin"
c this will be the same, in rygbergs

```

```
punch line optical depths last "agn_warm_absorber_ryg.lin" units rydberg
c
c warrm_absorber
c class nlr
c =====
c
```

this is a simple warm absorber model. It makes a plot of the transmitted continuum, and generates a list of lines with significant optical depths

---

## **aperture\_beam\_int** *test aperture beam command with intensity*

```
title test aperture beam command with intensity
c
c commands for density & abundances =====
hden 0
init file "honly.ini"
c
c other commands for details =====
constant temper 4
c
c commands controlling continuum =====
laser 2
ionization -2
c
c commands controlling geometry =====
stop thickness 10
sphere
aperture beam
c
c commands controlling output =====
c
c aperture_beam_int.in
c class geometry
c =====
```

This is a homogeneous sphere that is especially simple. The model is a test of the aperture command, a command that simulates observing part of an extended object. In this case the aperture is a beam centered on the center of the nebula, with a line of sight extending through the object.

The code carries along a dummy emission line ("Unit 1") with a constant intensity of  $1e-10$  erg cm<sup>-3</sup> s<sup>-1</sup>. The line goes through all of the code's infrastructure, and when the calculation is complete, the program confirms that the "luminosity" of the line is the emitting volume times  $1e-10$ . The aperture command is verified by asserting that the emission line has the correct "luminosity". In this case the inner radius is not specified so the returned value is unity.

---

## **aperture\_beam\_lum** *test aperture beam command with luminosity*

```
title test aperture beam command with luminosity
c
c commands controlling continuum =====
laser 2
q(h) 31.5
c
```

```

c commands for density & abundances =====
hden 0
init file "honly.ini"
c
c commands controlling geometry =====
stop thickness 10
radius 9
sphere
aperture beam
c
c other commands for details =====
constant temper 4
c
c commands controlling output =====
c
c aperture_beam_lum.in
c class geometry
c =====

```

This is a homogeneous sphere that is especially simple. The model is a test of the aperture command, a command that simulates observing part of an extended object. In this case the aperture is a beam centered on the center of the nebula, with a line of sight extending through the object.

The code carries along a dummy emission line ("Unit 1") with a constant intensity of  $1e-10$  erg cm<sup>-3</sup> s<sup>-1</sup>. The line goes through all of the code's infrastructure, and when the calculation is complete, the program confirms that the "luminosity" of the line is the emitting volume times  $1e-10$ . The aperture command is verified by asserting that this emission line has the correct "luminosity".

## **aperture\_slit** *test aperture slit command with luminosity*

```

title test aperture slit command with luminosity
c
c commands controlling continuum =====
laser 2
q(h) 31.5
c
c commands for density & abundances =====
hden 0
init file "honly.ini"
c
c commands controlling geometry =====
radius 9 10
sphere
aperture slit
c
c other commands for details =====
constant temper 4
c
c commands controlling output =====
c
c aperture_slit.in
c class geometry
c =====
c

```

This is a homogeneous sphere that is especially simple. The model is a test of the aperture command, a command that simulates observing part of an extended object. In this case the aperture is a long slit centered on the center of the nebula, extending beyond the outer reaches of the matter.



The code carries along a dummy emission line ("Unit 1") with a constant intensity of  $1e-10$  erg cm<sup>-3</sup> s<sup>-1</sup>. The line goes through all of the code's infrastructure, and when the calculation is complete, the program confirms that the "luminosity" of the line is the emitting volume times  $1e-10$ . The aperture command is verified by asserting that the emission line has the correct "luminosity".

---

## **blr\_f92 *standard blr cloud in Ferland et al. 1992***

```

title standard blr cloud in Ferland et al. 1992
c
c commands controlling continuum =====
table agn
ionization parameter -1
c
c commands for density & abundances =====
hden 11
abundances old solar 84
c
c commands controlling geometry =====
stop column density 25
c
c other commands for details =====
atom h-like element hydrogen levels resolved 10
atom h-like element helium levels resolved 10
iterate to convergence
c
c commands controlling output =====
print h-like departure hydrogen
print h-like populations hydrogen
print h-like departure helium
print h-like populations helium
plot continuum
normalize to "H 1" 1216 = 105.90
print line faint -2
punch overview last "blr_f92.ovr"
punch convergence reason "blr_f92.cvr"
punch element nitrogen last "blr_f92.nit"
punch element calcium last "blr_f92.ca"
punch dr "blr_f92.dr"
punch transmitted continuum "blr_f92.trn" last
c
c blr_f92.ini
c class blr
c =====

```

This is similar to one of the BLR models presented in Ferland et al. (1992) for the well-studied Seyfert galaxy NGC 5548. It has a very large column density and is marginally optically thick to electron scattering. The spectrum is given relative to Lya, and the intensity of this line is reset to produce a spectrum that is on the same intensity scale as that paper.

---

## **blr\_fp89 *final F+P 1989 BLR model table 3***

```

title final F+P 1989 BLR model table 3
c
c commands controlling continuum =====
table agn

```

```

ionization parameter -0.5
c
c commands for density & abundances =====
hden 9.5
init file="c84.ini"
abundances old solar 84
c
c commands controlling geometry =====
stop column density 25.5
c
c other commands for details =====
iterate to convergence
failures 2
c
c commands controlling output =====
c set a relatively high faint level since lines relative to Lya=100
print line faint -1
print h-like departure coefficients
plot continuum
normalize to "H 1" 1216 = 85.60
punch heating "blr_fp89.het"
punch convergence reason "blr_fp89.cvr"
punch cooling "blr_fp89.col"
punch overview last "blr_fp89.ovr"
punch element nitrogen last "blr_fp89.nit"
punch results last "blr_fp89.rlt"
punch dr "blr_fp89.dr"
c
c blr_fp89.in
c class blr
c =====

```

Ferland and Persson (1989) presented this calculation of a BLR cloud. The differences between the present predictions and those given by FP are largely due to improved treatment of Balmer line escape and destruction. The spectrum is given relative to a Lya intensity of 100. The column density is VERY large, to reproduce intensities of low-ionization lines, especially the Ca II lines.

---

## **blr\_hizqso *high Z quasar cloud***

```

title high Z quasar cloud
c
c commands controlling continuum =====
table agn
ionization parameter 1
c
c commands for density & abundances =====
c actual elec den will be about 100x larger
hden 8
abundances starburst 35
c
c commands controlling geometry =====
c this is more gramage than it appears because of high Z
stop column density 21
c
c other commands for details =====
c once elec den fail occurs due to tripping over He+ - He I-front
c code does recover so not a problem. The default for this is 10,
c and is set low because (hopefully) this would never occur
c hydrogen is a minor constituent in this high metallicity gas
c the electron density is well above the hydrogen density

```

```

failures 3
iterate 3
c
c commands controlling output =====
normalize "totl" 1909
punch dr "blr_hizqso.dr"
punch convergence reason "blr_hizqso.cvr"
punch overview last "blr_hizqso.ovr"
punch element nitrogen last "blr_hizqso.nit"
punch element chlorine last "blr_hizqso.cl"
punch element iron last "blr_hizqso.fe"
punch element neon last "blr_hizqso.ne"
punch element calcium last "blr_hizqso.ca"
punch element zinc last "blr_hizqso.zn"
punch heating last "blr_hizqso.het"
punch cooling last "blr_hizqso.col"
c
c
c blr_hizqso.in
c class blr
c =====

```

This is a model of a very high metallicity BLR cloud. It checks the intensities of some of the brighter lines, and is a check that the code can converge a cloud with this high Z.

Secondary ionization is very important when H is highly ionized, due to very high He abundance. Sec ionization becomes important at the He<sup>+</sup> - He ionization front, where H<sup>+</sup>/H is 1e-5.

---

## **blr\_kk81** *old blr*

```

title old blr
c
c commands controlling continuum =====
interpolate (0 -5) (.05 -5) (.1 0) (1 -0.5) (7.353 -2.233)
continue (735 -3.233) (800 -15) (8.e6 -15)
f(nu) -7.32148
constant gas pressure
c
c commands for density & abundances =====
hden 9.60206
init file="c84.ini"
abundances he-1 c-3.699 n-4 o-3.1549 ne-4 na=-8 mg-4.5229
continue al-10 si-4.4229 s-10 ar-10 ca-10 fe-4.5229 ni=-8
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
failures 2
iterate to convergence
c
c commands controlling output =====
print line faint -1
print ages
normalise "H 1" 1216 100
plot continuum range -3
punch overview last "blr_kk81.ovr"
punch dr last "blr_kk81.dr"
punch continuum last "blr_kk81.con"
punch ionizing continuum "blr_kk81.ion"

```

```

c
c
c blr_kk81.in
c class blr
c =====
c

```

This is the "standard" BLR model presented by Kwan and Krolik (1981).

>>refer blr cloud Kwan, J., & Krolik, J. 1981, ApJ, 250, 478

Compare line intensities to previous versions of CLOUDY by entering into table on page Error! Bookmark not defined..

The code caution that the resulting total pressure was not constant is to be expected. The KK calculation assumed constant gas pressure, but internally generature line radiation pressure is significant. Because of this the sum of gas plus radiation pressure was not constant although the gas pressure was.

## **blr\_level2 *test dominant level2 lines***

```

title test dominant level2 lines
c
c commands controlling continuum =====
table agn
ionization -2
c
c commands controlling geometry =====
stop column density 19
c
c commands for density & abundances =====
hden 10
element phosphorus abundance 0
c
c other commands for details =====
iterate
c
c commands controlling output =====
punch overview "blr_level2.ovr"
punch dr "blr_level2.dr"
c
c blr_level2.in
c class blr
c =====

```

This model checks predictions for the "level2" lines. These are lines that are normally very weak, have Opacity Project wavelengths, and g-bar collision strengths. Phosphorus is given a large abundance so that its level2 lines are significant.

## **blr\_n09\_p18 *BLR model, density 1e09 cm-3, flux of H-ion phots 1e18 cm2 s-1***

```

title BLR model, density 1e09 cm-3, flux of H-ion phots 1e18 cm2 s-1
c

```

```

c commands controlling continuum =====
table agn
phi(h) 18
c
c commands for density & abundances =====
hden 9
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
init "c84.ini"
iterate convergence
c these are to try to speed things up
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch overview "blr_n09_p18.ovr" last
punch dr "blr_n09_p18.dr" last
punch convergence reason "blr_n09_p18.cvr"
c
c blr_n09_p18.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n09\_p18\_z20 *BLR model, density 1e09 cm-3, flux of H-ion photos 1e18 cm2 s-1, Z=20***

```

title BLR model, density 1e09 cm-3, flux of H-ion photos 1e18 cm2 s-1, Z=20
c
c commands controlling continuum =====
table agn
phi(h) 18
c
c commands for density & abundances =====
hden 9
abundances starburst 20
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
iterate to convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
print h-like hydrogen departure coefficients
punch overview "blr_n09_p18_Z20.ovr"
punch dr "blr_n09_p18_Z20.dr"
punch convergence reason "blr_n09_p18_Z20.cvr"
punch convergence error "blr_n09_p18_Z20.cve"
c
c blr_n09_p18_Z20.in
c class blr

```

c =====

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n09\_p20 *BLR model, density 1e09 cm-3, flux of H-ion photos 1e20 cm2 s-1***

```

title BLR model, density 1e09 cm-3, flux of H-ion photos 1e20 cm2 s-1
c
c commands controlling continuum =====
table agn
phi(h) 20
c
c commands for density & abundances =====
hden 9
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
init "c84.ini"
iterate convergence
c these are to try to speed things up
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch overview "blr_n09_p18.ovr" last
punch dr "blr_n09_p18.dr" last
punch convergence reason "blr_n09_p18.cvr"
c
c blr_n09_p18.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n09\_p20\_z20 *BLR model, density 1e09 cm-3, flux of H-ion photos 1e20 cm2 s-1, Z=20***

```

title BLR model, density 1e09 cm-3, flux of H-ion photos 1e20 cm2 s-1, Z=20
c
c commands controlling continuum =====
table agn
phi(h) 20
c
c commands for density & abundances =====
abundances starburst 20
hden 9
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
init "c84.ini"

```

```

iterate convergence
c these are to try to speed things up
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch overview "blr_n09_p20_Z20.ovr" last
punch dr "blr_n09_p20_Z20.dr" last
punch convergence reason "blr_n09_p20_Z20.cvr"
c
c blr_n09_p20_Z20.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

This simulation is optically thin in the Lyman continuum - no H ionization front is present. As a result it can be difficult to converge.

---

### **blr\_n09\_p22 *BLR model, density 1e09 cm-3, flux of H-ion phots 1e20 cm2 s-1***

```

title BLR model, density 1e09 cm-3, flux of H-ion phots 1e20 cm2 s-1
c
c commands controlling continuum =====
table agn
phi(h) 22
c
c commands for density & abundances =====
hden 9
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
init "c84.ini"
iterate convergence
c these are to try to speed things up
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch overview "blr_n09_p22.ovr" last
punch dr "blr_n09_p22.dr" last
punch convergence reason "blr_n09_p22.cvr"
c
c blr_n09_p22.in
c class blr
c =====

```

This is one of the models that sample the LOC plane.

---

### **blr\_n09\_p22\_z20 *BLR model, density 1e09 cm-3, flux of H-ion phots 1e22 cm2 s-1, Z=20***

```

title BLR model, density 1e09 cm-3, flux of H-ion phots 1e22 cm2 s-1, Z=20
c
c commands controlling continuum =====
table agn
phi(h) 22
c
c commands for density & abundances =====
abundances starburst 20
hden 9
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
init "c84.ini"
iterate convergence
c these are to try to speed things up
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch overview "blr_n09_p22_Z20.ovr" last
punch dr "blr_n09_p22_Z20.dr" last
punch convergence reason "blr_n09_p22_Z20.cvr"
c
c blr_n09_p22_Z20.in
c class blr
c =====

```

This is one of the models that sample the LOC plane.

---

## **blr\_n11\_p20 *BLR model, density 1e11 cm-3, flux of H-ion phots 1e20 cm2 s-1***

```

title BLR model, density 1e11 cm-3, flux of H-ion phots 1e20 cm2 s-1
c
c commands controlling continuum =====
table agn
phi(h) 20
c
c commands for density & abundances =====
hden 11
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
iterate convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch convergence reason "blr_n11_p20.cvr"
punch overview "blr_n11_p20.ovr" last
punch dr "blr_n11_p20.dr" last
c
c blr_n11_p20.in

```



```
c class blr
c =====
c
```

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n11\_p20\_z20 *BLR model, density 1e11 cm-3, flux of H-ion photos 1e20 cm2 s-1, Z=20***

```
title BLR model, density 1e11 cm-3, flux of H-ion photos 1e20 cm2 s-1, Z=20
c
c commands controlling continuum =====
table agn
phi(h) 20
c
c commands for density & abundances =====
hden 11
abundances starburst 20
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
iterate convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch cooling "blr_n11_p20_Z20.col"
punch heating "blr_n11_p20_Z20.het"
punch element calcium "blr_n11_p20_Z20.cal"
punch overview "blr_n11_p20_Z20.ovr"
punch dr "blr_n11_p20_Z20.dr"
punch temperature history "blr_n11_p20_Z20.tem"
c
c blr_n11_p20_Z20.in
c class blr
c =====
```

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n12\_p19 *BLR model, density 1e12 cm-3, flux of H-ion photos 1e19 cm2 s-1***

```
title BLR model, density 1e12 cm-3, flux of H-ion photos 1e19 cm2 s-1
c
c commands for density & abundances =====
c
c commands controlling continuum =====
table agn
phi(h) 19
c
c commands controlling geometry =====
```

```

hden 12
stop column density 23
c
c other commands for details =====
iterate convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch overview "blr_n12_p19.ovr" last
punch dr "blr_n12_p19.dr" last
c
c
c blr_n12_p19.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

### **blr\_n12\_p19\_z20 *BLR model, density 1e12 cm-3, flux of H-ion photos 1e19 cm2 s-1, Z=20***

```

title BLR model, density 1e12 cm-3, flux of H-ion photos 1e19 cm2 s-1, Z=20
c
c commands controlling continuum =====
table agn
phi(h) 19
c
c commands for density & abundances =====
hden 12
abundances starburst 20
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
failures 5
iterate convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c
c commands controlling output =====
normalize to "totl" 1216 100
punch convergence reason "blr_n12_p19_Z20.cvr"
punch overview "blr_n12_p19_Z20.ovr" last
punch dr "blr_n12_p19_Z20.dr" last
c
c blr_n12_p19_Z20.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

### **blr\_n13\_p18 *BLR model, density 1e13 cm-3, flux of H-ion photos 1e18***

## *cm2 s-1*

```

title BLR model, density 1e13 cm-3, flux of H-ion phots 1e18 cm2 s-1
c
c this is a very low ionization cloud
c the conditions, and some lines, are surprisingly sensitive
c to the treatment of hydrogen molecules
c
c commands controlling continuum =====
table agn
phi(h) 18
c
c commands for density & abundances =====
hden 13
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
iterate to convergence
c
c commands controlling output =====
normalize to "tot1" 1216 100
init "c84.ini"
no level2 lines
punch overview "blr_n13_p18.ovr" last
punch dr "blr_n13_p18.dr" last
punch convergence reason "blr_n13_p18.cvr"
c
c blr_n13_p18.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n13\_p18\_z20** *BLR model, density 1e13 cm-3, flux of H-ion phots 1e18 cm2 s-1, Z=20*

```

title BLR model, density 1e13 cm-3, flux of H-ion phots 1e18 cm2 s-1, Z=20
c
c this is a very low ionization cloud
c the conditions, and some lines, are surprisingly sensitive
c to the treatment of hydrogen molecules
c
c commands controlling continuum =====
phi(h) 18
table agn
c
c commands for density & abundances =====
hden 13
abundances starburst 20
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
c this sim needs roughly 10 iterations to converge

```

```

iterate convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c abort if any failures occur - we want to pass this test
failures 1
c
c commands controlling output =====
normalize to "totl" 1216 100
punch overview "blr_n13_p18_Z20.ovr" last
punch overview "blr_n13_p18_Z20.ovr1"
punch dr "blr_n13_p18_Z20.dr" last
punch convergence reason "blr_n13_p18_Z20.cvr"
c
c
c blr_n13_p18_Z20.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n13\_p22 *BLR model, density 1e13 cm-3, flux of H-ion phots 1e22 cm2 s-1***

```

title BLR model, density 1e13 cm-3, flux of H-ion phots 1e22 cm2 s-1
c
c commands controlling continuum =====
table agn
phi(h) 22
c
c commands for density & abundances =====
hden 13
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
iterate convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c
c commands controlling output =====
print h-like departure hydrogen
print h-like departure helium
normalize to "totl" 1216 100
punch overview "blr_n13_p22.ovr" last
punch dr "blr_n13_p22.dr" last
c
c blr_n13_p22.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

## **blr\_n13\_p22\_z20 *BLR model, density 1e13 cm-3, flux of H-ion phots***

## ***1e18 cm2 s-1, Z=20***

```

title BLR model, density 1e13 cm-3, flux of H-ion phots 1e18 cm2 s-1, Z=20
c
c commands controlling continuum =====
table agn
phi(h) 22
c
c commands for density & abundances =====
hden 13
abundances starburst 20
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
iterate convergence
c these are to try to speed things up
init "c84.ini"
no level2 lines
c
c commands controlling output =====
normalize to "tot1" 1216 100
punch overview "blr_n13_p22_Z20.ovr"
punch dr "blr_n13_p22_Z20.dr"
punch heating "blr_n13_p22_Z20.het"
punch cooling "blr_n13_p22_Z20.col"
c
c blr_n13_p22_Z20.in
c class blr
c =====

```

This is one of the 5 models that sample the LOC plane.

---

## ***blr\_nf84 early model of blr***

```

title early model of blr
c
c commands controlling continuum =====
ionization parameter -1.92
power law -1 100 0.01
c
c commands for density & abundances =====
hden 9.5
init, file="ism.ini"
abundances cameron
stop lyman optical 6
c
c commands controlling geometry =====
c
c other commands for details =====
constant gas pressure
iterate to convergence
c
c commands controlling output =====
print diffuse continuum
plot continuum
punch overview last "blr_nf84.ovr"

```

```

punch dr last "blr_nf84.dr"
punch results last "blr_nf84.rlt"
c
table lines "LineList_strong.dat"
c blr_nf84.in
c class blr
c =====
c

```

This is an example of a "conventional" BLR calculation. The parameters are similar to those of Table 1 of Netzer and Ferland (1984). Notice that the ratio of Lyalpha to Hbeta ratio is much larger than observed.

>>refer blr model Netzer, H., & Ferland, G. J. 1984, PASP, 96, 593

---

## **blr\_nf84\_45deg *early model of BLR, with illumination at 45 degree angle***

```

title early model of BLR, with illumination at 45 degree angle
c
c commands controlling continuum =====
ionization parameter -1.92
power law -1 100 0.01
c this continuum is a beam coming in at 45 degrees
illuminate 45 degrees
c
c commands for density & abundances =====
hden 9.5
abundances cameron
stop lyman optical 6
c
c commands controlling geometry =====
c
c other commands for details =====
init file "ism.ini"
constant gas pressure
iterate to convergence
c
c commands controlling output =====
print diffuse continuum
plot continuum
punch overview "blr_nf84_45deg.ovr"
punch dr "blr_nf84_45deg.dr"
punch results last "blr_nf84_45deg.rlt"
c
c blr_nf84_45deg.in
c class blr
c =====
c

```

This is an example of a "conventional" BLR calculation. The parameters are similar to those of Table 1 of Netzer and Ferland (1984). Notice that the ratio of Lyalpha to Hbeta ratio is much larger than observed.

>>refer blr model Netzer, H., & Ferland, G. J. 1984, PASP, 96, 593

---

## **blr\_rnfa *table 1 of Rees et al. ApJ 347, 648***

```

title table 1 of Rees et al. ApJ 347, 648
c
c commands controlling continuum =====
table agn
ionization parameter -2
c
c commands for density & abundances =====
hden 10
init file="c84.ini"
abundances he -1 c -3.328 n -4.0088 o -3.0809 ne -4 na -20
continue mg -4.3768 al -5.5686 si -4.36653 s -4.76955
continue a -5.4202 ca -5.6383 fe -4.4815 ni -20
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
iterate to convergence
c
c commands controlling output =====
normalise "H 1" 1216 = 100
print line faint -1
punch overview last "blr_rnfa.ovr"
punch continuum last "blr_rnfa.con" units microns
punch transmitted continuum last "blr_rnfa.trn" units rydbergs
punch element nitrogen last "blr_rnfa.nit"
punch dr last "blr_rnfa.dr"
punch results last "blr_rnfa.rlt"
c
c blr_rnfa.in
c class blr
c =====
c

```

This is the lower density cloud computed in Rees et al. (1989). Table 1 of that paper lists the predictions, which were a mean of those of Hagai Netzer's ION and roughly version 76 of CLOUDY. The lines are generally still in good agreement with the predictions of that paper. In particular the changes in the line fluxes shown in Figure 1 of that paper are reproduced quite well.

---

## **blr\_rnfb *table 1 of Rees et al. ApJ 347, 648***

```

title table 1 of Rees et al. ApJ 347, 648
c
c blr model from
c
c commands controlling continuum =====
table agn
ionization parameter -2
c
c commands for density & abundances =====
hden 12
init file="c84.ini"
abundances he -1 c -3.328 n -4.0088 o -3.0809 ne -4 na -20
continue mg -4.3768 al -5.5686 si -4.36653 s -4.76955
continue a -5.4202 ca -5.6383 fe -4.4815 ni -20
c
c commands controlling geometry =====
stop column density 23
c

```

```

c other commands for details =====
iterate to convergence
c
c commands controlling output =====
print line faint -1
punch overview last "blr_rnfb.ovr"
punch element nitrogen last "blr_rnfb.nit"
punch dr last "blr_rnfb.dr"
punch results last "blr_rnfb.rlt"
normalise "H 1" 1216 = 91.8
c
c blr_rnfb.in
c class blr
c =====
c

```

This is a very dense cloud, and was computed in Rees et al. (1989). Table 1 of that paper lists the predictions, which were a mean of those of Hagai Netzer's ION and roughly version 76 of CLOUDY. The lines are generally still in good agreement with the predictions of that paper. In particular the changes in the line fluxes shown in Figure 1 of that paper are reproduced quite well. The fluxes of Ly $\alpha$  and H $\beta$  are not reproduced with great precision by this model because of changes in collision rates for hydrogen and especially the form of the escape probability function for subordinate lines. As Figure 1 of RNF showed the line intensities are very sensitive to density for these parameters.

## **coll\_coronal** *model of active region of solar corona*

```

title model of active region of solar corona
c
c commands controlling continuum =====
c temperature of corona
coronal equilibrium, 2.5e6 K
c assume the sun is a blackbody at 5770K
blackbody 5770 K
luminosity solar linear 1
c
c commands for density & abundances =====
hden log=10
c
c commands controlling geometry =====
c its height above photosphere and thickness
radius 11 thickness = 10
c the gas fully covers the center of symmetry
sphere
c
c other commands for details =====
iterate
c
c commands controlling output =====
normalize to "o 8" 18.97A
punch dr "coll_coronal.dr" last
punch continuum "coll_coronal.con" last units angstroms
c
c
c coll_coronal.in
c class coronal
c =====
c

```

This is a rough model of the solar corona. The test checks that the coronal equilibrium commands work. The



gas is predominantly collisionally ionized.

---

## ***coll\_heat\_only test code in limit where ONLY mechanical heating is present***

```

title test code in limit where ONLY mechanical heating is present
c
c commands controlling continuum =====
no photoionization
blackbody 5000 STE
c
c commands controlling geometry =====
stop column density 15
c
c commands for density & abundances =====
hden 10
c
c other commands for details =====
hextra -1.5
iterate
c
c commands controlling output =====
c this command must be tested somewhere
print cooling, zone 5
c output files
punch overview "coll_heat_only.ovr"
punch heating "coll_heat_only.het"
punch cooling "coll_heat_only.col"
punch convergence reason "coll_heat_only.cvr"
punch dr "coll_heat_only.dr"
punch continuum "coll_heat_only.con" units angstroms
c
c
c coll_heat_only.in
c class coronal
c =====

```

This test is an optically thin collisionally ionized gas with no photoionization at all.

---

## ***coll\_t3 coronal equilibrium at 10<sup>4</sup> K***

```

title coronal equilibrium at 104 K
c
c commands controlling continuum =====
coronal 3
c add component of cosmic rays to drive chemistry
cosmic ray background
c
c other commands for details =====
iterate
c
c commands controlling geometry =====
stop column density 15
c
c commands for density & abundances =====
hden 10

```

```

c
c commands controlling output =====
punch overview "coll_t3.ovr"
punch continuum "coll_t3.con" units angstroms
c
c
c coll_t3.in
c class coronal
c =====

```

This tests conditions of collisional equilibrium at low densities. This is one of a series of sims coll\_t?.in which test ionization over a range of temperatures. This one, unlike the others, includes cosmic rays. Chemistry is important at this low temperature and the chemical network will collapse without a source of ionization. The cosmic rays provide this source of ionization.

---

### **coll\_t4 coronal equilibrium at 10<sup>4</sup> K**

```

title coronal equilibrium at 10^4 K
c
c commands controlling continuum =====
coronal 4
c
c other commands for details =====
iterate
c
c commands controlling geometry =====
stop column density 15
c
c commands for density & abundances =====
hden 10
c
c commands controlling output =====
punch overview "coll_t4.ovr"
punch continuum "coll_t4.con" units angstroms
c
c
c coll_t4.in
c class coronal
c =====

```

this tests conditions of collisional equilibrium at low densities

---

### **coll\_t5 coronal equilibrium at 10<sup>5</sup> K**

```

title coronal equilibrium at 10^5 K
c
c commands controlling continuum =====
coronal 5
c
c commands controlling geometry =====
stop column density 15
c
c commands for density & abundances =====
hden 10
c
c other commands for details =====

```

```

iterate
c
c
c commands controlling output =====
c normalize to strong HeII line
normalize to "He 2" 303.8A
punch overview "coll_t5.ovr"
punch continuum "coll_t5.con" units angstroms
c
table lines "LineList_He_like.dat"
c coll_t5.in
c class coronal
c =====

```

This is a test collisional ionization equilibrium at  $1e5$  K.

---

## **coll\_t6 coronal equilibrium at $10^6$ K**

```

title coronal equilibrium at  $10^6$  K
c
c commands controlling continuum =====
coronal 6
c
c commands controlling geometry =====
stop column density 15
c
c commands for density & abundances =====
hden 10
c
c other commands for details =====
iterate
c
c commands controlling output =====
normalize to "c 5" 40.27A
c output files
punch overview "coll_t6.ovr"
punch continuum "coll_t6.con" units angstroms
c
c coll_t6.in
c class coronal
c =====

```

This test is an optically thin collisionally ionized gas.

---

## **coll\_t7 coronal equilibrium at $10^7$ K**

```

title coronal equilibrium at  $10^7$  K
c
c commands controlling continuum =====
coronal 7
c
c commands controlling geometry =====
stop column density 15
c
c commands for density & abundances =====
hden 10

```

```

c
c other commands for details =====
c
normalize to "o 8" 18.97A
c commands controlling output =====
punch overview "coll_t7.ovr"
punch continuum "coll_t7.con" units angstroms
punch cooling "coll_t7.col"
c
c coll_t7.in
c class coronal
c =====

```

Test with only collisional ionization at a high temperature.

---

## **dynamics\_orion\_flow** *Orion nebula blister with wind*

```

title Orion nebula blister with wind
c
c commands controlling continuum =====
blackbody 40000
phi(h) 13.0
brems 6
phi(h) 10
c
c commands for density & abundances =====
hden 4
abundances hii region no grains
grains orion no qheat single
c
c commands controlling geometry =====
wind -7 km/s advection
c this accounts for matter beyond end of the model
double optical depths
sphere
stop AV 5
stop temperature linear 5
c
c other commands for details =====
iterate 4
magnetic field -4
cosmic rays background
c this has no effect on dynamics, but is to desaturate the lines as per flow
turbulence 8 km/sec no pressure
c try to speed up simulation without hurting dynamics
init "fast.ini"
c use Bakes & Tielens heating
set PAH Bakes
* init "c84.ini"
* failures 5
c
c commands controlling output =====
c want to print line intensities are surface brightness, per arcsec^2
print lines surface brightness arcsec
print diffuse continuum
print line faint -1.5
punch overview "dynamics_orion_flow.ovr" no hash
punch hydrogen 21 cm "dynamics_orion_flow.21cm" no hash
punch dr "dynamics_orion_flow.dr" no hash
punch continuum last "dynamics_orion_flow.con" units microns no hash
punch pressure "dynamics_orion_flow.pre" no hash

```

```

punch heating "dynamics_orion_flow.het" no hash
punch cooling "dynamics_orion_flow.col" no hash
punch molecules "dynamics_orion_flow.mol" no hash
punch wind "dynamics_orion_flow.wnd" last
c
c dynamics_orion_flow.in
c class dynamics hii pdr
c =====
c

```

This is a model similar in spirit to the blister geometry H+ region model computed by Baldwin et al. (1991), but with a D-critical flow. Many physical processes have been disabled to make this simulation faster. Grain physics is not done so the gas temperature is incorrect. The main purpose is to do a quick test of the dynamical flow with grain opacities included. The slow directory contains a full simulation of a flow like Orion.

---

## **dynamics\_veryfast** *very fast wind model*

```

title very fast wind model
c
c commands controlling continuum =====
table agn
ionization parameter -2
c
c commands for density & abundances =====
hden 9
c this is a set of commands to speed up calc
init file "fast.ini"
c
c commands controlling geometry =====
stop thickness 8.4
c this is a subsonic wind with advection
wind velo -5 advection
c
c other commands for details =====
c absolutely no convergence problems should occur
failures 1
c this degrades the continuum resolution by factor of 3, to
c speed up calc
set continuum resolution 3
iterate 4
c
c commands controlling output =====
punch pressure last "dynamics_veryfast.pre"
punch total opacity last "dynamics_veryfast.opc"
punch ionizing continuum last "dynamics_veryfast.ion"
punch continuum last "dynamics_veryfast.con"
punch dr "dynamics_veryfast.dr"
c
c
c dynamics_veryfast.in
c class dynamics
c =====
c

```

This is meant to be a very fast calculation to use when running extensive debug-enabled runtimes.

---

## **dynamics\_veryfast\_rec** *very fast wind model*

```

title very fast wind model
c
c commands controlling continuum =====
table agn
ionization parameter -2
c
c commands for density & abundances =====
hden 9
c this is a set of commands to speed up calc
init file "fast.ini"
c
c commands controlling geometry =====
stop thickness 8.4
c this is a subsonic wind with advection
wind velo 5 advection
c
c other commands for details =====
c absolutely no convergence problems should occur
failures 1
c this degrades the continuum resolution by factor of 3, to
c speed up calc
set continuum resolution 3
iterate 4
c
c commands controlling output =====
punch pressure last "dynamics_veryfast_rec.pre"
punch total opacity last "dynamics_veryfast_rec.opc"
punch ionizing continuum last "dynamics_veryfast_rec.ion"
punch continuum last "dynamics_veryfast_rec.con"
punch dr "dynamics_veryfast_rec.dr"
c
c
c dynamics_veryfast_rec.in
c class dynamics
c =====
c

```

This is meant to be a very fast calculation to use when running extensive debug-enabled runtimes.

---

## **dynamics\_wind** *test of equations of motion in a very highly ionized wind*

```

title test of equations of motion in a very highly ionized wind
c
c test motion in electron scattering limit
c radiative acceleration (e- only) is 9.54E-7 cm s^-2
c terminal velocity (e- only) is 7.6 km s^-1
c
c commands controlling continuum =====
table agn
luminosiy (total) 45
c
c commands for density & abundances =====
hden 4
c this only includes H and He

```

```

init file="hheonly.ini"
c
c commands controlling geometry =====
radius (parsecs) 1
stop thickness (parsecs) -1
c a slow wind velocity, will assert velocity at outer radius
wind velo 0.1 ball
c
c other commands for details =====
c this will speed things up a bit
no level2
no radiation pressure
constant temperature 8
c
c commands controlling output =====
punch pressure "dynamics_wind.pre"
punch wind "dynamics_wind.wnd"
punch dr "dynamics_wind.dr"
c
c dynamics_wind.in
c class limit dynamics
c =====
c

```

This tests the radiative acceleration and terminal velocity of a wind in which only electron scattering is important. The parameters were chosen so that electron scattering is the dominant opacity source, so that the equations can be solved both numerically (in the example) and analytically (the expected solution given above). In a realistic wind the gas would be more neutral and line driving would dominate. The force multiplier, given in the punch wind output, is nearly unity as a result.

#### Checks:

- The radiative acceleration is correct (e- 9.543910-7 cm s-2).
- The terminal velocity should be 7.57 km s-1.
- Force multiplier near unity (no line driving since so highly ionized).
- Thickness of cloud correct ( $R-R_0 + dr/2$  should be 3.086391017 cm).

### ***feii\_hin test feii in high density limit***

```

title test feii in high density limit
c
c commands controlling continuum =====
black 25000 K
ionization -10
c
c commands for density & abundances =====
hden 18
c force a high electron density to insure collisions dominate
eden 18
c
c commands controlling geometry =====
constant tempera 4
stop zone 1
c
c other commands for details =====
atom feii
iterate
c want to test collision only limit

```

```

no induced processes
c
c commands controlling output =====
c
c
c feii_hin.in
c class limit
c =====
c

```

This checks that, at high particle densities, in which the gas should be in collisional equilibrium, the level populations of the large model Fe<sup>+</sup> ion go to the proper values, where the departure coefficients are all equal to unity.

---

### ***feii\_hirad*** *feii in case of high radiation density limit*

```

title feii in case of high radiation density limit
c
c commands controlling continuum =====
black 8000 K lte
c
c commands for density & abundances =====
hden 9
element iron abundance 2
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
atom feii
constant tempera 8000
iterate
c
c commands controlling output =====
punch feii relative populations range 0 200 "feii_hirad.lv1" last
punch feii relative populations range 201 371 "feii_hirad.lv2" last
c
c
c feii_hirad.in
c class limit
c =====
c

```

This checks that, at high radiation densities, in which the gas is irradiated by a blackbody in strict thermodynamic equilibrium, the level populations of the large model Fe<sup>+</sup> ion go to the proper values, where the departure coefficients are all equal to unity.

---

### ***feii\_pump*** *test feii in continuum pumped limit*

```

title test feii in continuum pumped limit
c
c commands controlling continuum =====
black 25000 K
ionization -5

```



```

c
c commands for density & abundances =====
hden 10
abundances old solar 84
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
Case B
iterate
constant tempera 4
atom feii
c
c commands controlling output =====
punch feii continuum "feii_pump.con"
c
c feii_pump.in

```

This is a constant temperature low ionization cloud, with BLR-like densities, which includes the large FeII atom. The tests check on the emission predicted in the Fe II bands.

This model tests the large FeII model in the optically thin, continuum pumped limit. The zone thickness is set to a small value (1 cm) so that full continuum hits atom.

### ***feii\_ste thermal equilibrium of FeII in STE limit***

```

title thermal equilibrium of FeII in STE limit
c
c commands controlling continuum =====
black 8000 K lte
hden 9
c
c commands for density & abundances =====
element iron abundance 2
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
atom feii
iterate
c
c commands controlling output =====
punch feii continuum last "feii_ste.con"
punch feii column density last "feii_ste.col"
punch feii populations all last "feii_ste.pop"
c
c feii_ste.in
c class function
c =====
c

```

This model has a very high iron abundance, 100x H, and most Fe is in the form of Fe+. It is irradiated by a blackbody in strict thermodynamic equilibrium. We check that the temperature of the gas is equal to the radiation temperature, to confirm that the thermal properties of the model FeII atom obey thermodynamics.

## **feii\_t4 *FeII emission in typical intermediate density photoionized cloud***

```

title FeII emission in typical intermediate density photoionized cloud
c
c commands controlling continuum =====
black 5e5K
ionization parameter -2
c
c commands for density & abundances =====
hden 8
element iron abundance 2
element iron ionization -2 0 -2
c
c commands controlling geometry =====
stop thickness 5
c
c other commands for details =====
constant temperature 4
atom feii
iterate 3
c
c commands controlling output =====
set line precision 6
print lines column
print line inward
print line sort wavelength
print line faint 0
punch feii continuum last "feii_t4.con" units angstroms
punch feii inward continuum last "feii_t4.cin" units angstroms
punch feii outward continuum last "feii_t4.cout" units angstroms
punch feii column density last "feii_t4.col"
punch feii populations all last "feii_t4.pop"
punch line emissivity "feii_t4.ems"
Fe2b 2797
Inwd 2797
end of lines
c
c feii_t4.in
c class function
c =====
c

```

This model has a very high iron abundance, 100x H, and most Fe is in the form of Fe<sup>+</sup>. It has intermediate density and should produce an FeII spectrum something like an AGN. The set line precision 6 increases the number of significant figures in the wavelengths for each line. This is needed to get the right FeII inward band.

---

## **func\_abund\_fluc *check fluctuating heavy-element abundances***

```

title check fluctuating heavy-element abundances
c commands controlling continuum =====
blackbody 5
ionization -2
c
c commands for density & abundances =====
hden 3

```

```

c this varies the abundances between default and 1e-4 times default,
c over a 1e10 cm period - we will compute half a cycle
fluctuation abundances 10.3 0 -4
c
c commands controlling geometry =====
c stop at half a cycle so pick off lowest abundances
stop thickness 10
c
c other commands for details =====
c speed things up a bit
init "c84.ini"
c
c commands controlling output =====
print column density
punch overview "func_abund_fluc.ovr"
punch abundances "func_abund_fluc.abn"
c
c commands giving asserts
c
c func_abund_fluc.in
c class function
c =====
c

```

This checks that the variable abundances option works

---

## ***func\_distance check that distance and print flux earth commands work***

```

title check that distance and "print flux earth" commands work
c
c commands controlling continuum =====
c luminosity in ionizing radiation, this and distance
c were chosen so that flux at earth associated with total
c luminosity of object is 1 erg cm-2 s-1
luminosity 40
laser 2
c
c commands for density & abundances =====
hden 4
init file "honly.ini"
c
c commands controlling geometry =====
stop zone 1
radius 17
c
c other commands for details =====
distance 19.450395
constant temperature 4
c
c commands controlling output =====
c give flux at Earth
print lines, flux at Earth
c
c func_distance.in
c class function
c =====
c

```

Normally the code predicts the intensity or luminosity of the emission lines. This test confirms that it can predict the flux recieved at the Earth instead. The model is the simplest and fastest that can be computed - a H-only constant temperature single zone. The total luminosity is set to  $1e40$  erg/s, and the ionization source is a laser at 2 ryd. With these set, the total luminosity in ionizing radiation, the total luminosity in the incident continuum (the emission line labeled "Inci 0") will be  $1e40$ .

The code will predict the flux at the Earth if both the distance to the object is specified with the distance command, and this is requested with the print flux earth command. The distance was chosen so that the total flux at the Earth will be 1 erg/s. This is asserted at the end of the calculation.

---

### **func\_dlaw** *test model with dlaw table*

```

title test model with dlaw table
c
c commands controlling continuum =====
phi(H) 15
table agn
c
c commands for density & abundances =====
dlaw table radius
continue 16 9
continue 17 7
continue 18 5
continue 19 3
continue 20 1
continue 21 -1
end of dlaw
init "honly.ini"
c
c commands controlling geometry =====
sphere
filling factor -5
radius 17 20
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
punch dr "func_dlaw.dr"
punch overveiw "func_dlaw.ovr"
c
c
c func_dlaw.in
c class function
c =====
c

```

this model tests the dlaw density table command

---

### **func\_fulltrace** *test full trace output*

```

title test full trace output
c
c commands controlling continuum =====
trace h-like hydrogen full no print

```

```

stop zone 1
table agn
ionization parameter -0.5
c
c commands for density & abundances =====
hden 9.5
init file="honly.ini"
c
c other commands for details =====
atom h-like element hydrogen collapsed levels 1
c
c commands controlling output =====
c func_fulltrace.in
c class function
c =====

```

this checks that trace output functions correctly

---

## **func\_globule** *test of globule command*

```

title test of globule command
c
c commands controlling continuum =====
black 40000 K
ionization parameter -1.5
c
c commands for density & abundances =====
globule initial density=5 scale depth=14
init file "ism.ini"
abundances hii region no qheat
c
c commands controlling geometry =====
c
c other commands for details =====
failures 2
no diffuse line pumping
c
c commands controlling output =====
print line faint -1.5
punch overview last "func_globule.ovr"
punch results last "func_globule.rlt"
punch dr "func_globule.dr" last
punch heating "func_globule.het" last
punch cooling "func_globule.col" last
c
c func_globule.in
c class function
c =====
c

```

This model uses the globule command, tests that the zoning logic works for this extreme case, and that the code is able to converge the globule model.

---

## **func\_grid\_line\_ratios** *test generating line ratios in a grid run*

```

title test generating line ratios in a grid run

```

```

c
c commands controlling continuum =====
blackbody 40,000K
ionization parameter -2
c
c commands for density & abundances =====
c these are to speed up the calculation, only do H, O, and Ne
init "honly.ini"
element oxygen on
element neon on
element sulphur on
element oxygen ionization 1 1 1 0.01
element neon ionization 1 1 1 0.01
element sulphur ionization 1 1 1 0.01
c vary the hydrogen density
hden 4 vary
grid 2 6.1 1
c
c other commands for details =====
c these are constant temperature models, vary T
constant temperature 4 vary
grid 4000 17000 3000
stop zone 1
c
c commands controlling output =====
punch line list "func_grid_line_ratios.pun" "func_grid_line_ratios.dat" ratio no hash
punch grid "func_grid_line_ratios.grd"
c
c commands giving (lack of) assert =====
c
c func_grid_line_line_ratios.in
c class function
c =====

```

This uses the grid command to compute line ratios for a wide range of density and temperature. The ionization is set to a uniform value and only a few elements are included. this makes the calculation faster and prevents recombination [O III] 4363 from becoming important (there is no O+3).

These are the line ratios mentioned as limits in the Johnstone et al. Spitzer cooling flow filament paper (2007).

---

## **func\_hotgas\_coolstar *test very soft continuum, very hot gas***

```

title test very soft continuum, very hot gas
c
c commands controlling continuum =====
c put in the cosmic background as the only continuum source
CMB
c gas has constant temperature of 1e6 K
constant temperature, t=6
c
c commands for density & abundances =====
hden 10.
c
c commands controlling geometry =====
stop zone 1
c specify a thin cell of gas - 1 cm thick
set dr 0
c
c other commands for details =====
c iterate since gas is optically thin

```

```

iterate
c
c commands controlling output =====
c
c func_hotgas_coolstar.in
c class limit
c =====
c

```

This is a test where the CMB is the only continuum source. It does not extend to energies where the code needs to work. There are special cases used in this situation, for continuum addressing, so this checks whether those still function.

---

## **func\_ion\_increase** *test model where ionization increases with depth*

```

title test model where ionization increases with depth
c
c commands controlling continuum =====
ionization parameter -4
table agn
c
c commands for density & abundances =====
hden 7 -4
init "hheonly.ini"
element oxygen on
element iron on
c
c commands controlling geometry =====
sphere
filling factor -5
radius 17 20
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
punch dr "func_ion_increase.dr"
punch overveiw "func_ion_increase.ovr"
punch element oxygen "func_ion_increase.oxy"
punch element iron "func_ion_increase.fe"
c
c
c func_ion_increase.in
c class function
c =====
c

```

This density falls off faster than  $1/r^2$  so the ionization increases with depth. Most sims have decreasing rather than increasing ionization.

---

## **func\_lines** *create output file with list of func\_lines*

```

title create output file with list of func_lines
atom h2 levels large
atom feii
test

```

```

c
c commands controlling continuum =====
c commands for density & abundances =====
c commands controlling geometry =====
c other commands for details =====
c
c commands controlling output =====
c this creates the file "func_lines.lab" that contains a list
c of all emission labels
punch line labels "func_lines.lab"
c this creates the file "func_lines.lis" that contains a list
c of the emission func_lines that are transported.
punch line data "func_lines.lis"
c
c There is nothing extra to assert, since test does include many asserts
c
c func_lines.in
c class function
c =====
c

```

This runs the standard "test" case, and then creates the line data and labels files. Test by itself includes many asserts, so no further asserts are needed here.

The file func\_lines.lab is a useful list of all lines predicted by the code. Cut and paste this into other places when you need to find a particular emission line.

The func\_lines lines.dat gives atomic data for all the lines, and their critical density at 10000 K. The large H2 and Fe II model atoms are turned on to include their lines.

## **func\_map** *map of heating vs cooling*

```

title map of heating vs cooling
c
c commands controlling continuum =====
table agn
ionization parameter -2.5
c
c commands for density & abundances =====
hden 0
abundances old solar 84
c
c commands controlling geometry =====
set dr 0
stop lyman optical depth -7
stop zone 1
c
c other commands for details =====
c
c commands controlling output =====
c save map information to generate plot for hazy
punch map file="func_map.dta"
set nmaps 100
c
c map.in
c class function
c =====
c

```



This is a test of the continuity of the code over a very large range of temperature. It was used to produce one of the thermal maps shown in Hazy.

Checks:

- No breaks in the heating and cooling curves where various approximations change.

---

### **func\_sdrmin** *test set drmin command*

```

title test set drmin command
c
c commands controlling continuum =====
blackbody 8e4
luminosity 37.5
c
c commands for density & abundances =====
hden 4
c
c commands controlling geometry =====
radius 17
sphere
stop thickness 17.03
stop temperature off
c
c other commands for details =====
set drmin 15.5
cosmic rays background
c
c commands controlling output =====
punch dr "func_sdrmin.dr" last
c

```

This simulation tests the SET DRMIN command. It is a toy model of a planetary nebula that is designed to extend into the PDR. The minimum stepsize is deliberately set much too large so that we would immediately notice if the command was broken. It also tests if we still hit the correct outer radius when SET DRMIN is used. Without the SET DRMIN command this sim would need 410 zones (trunk@2760).

---

### **func\_set\_ion** *test impact of setting ionization*

```

title test impact of setting ionization
c
c commands for density & abundances =====
init "ism.ini"
hden 0
element hydrogen ionization 0 1
element helium ionization 0 1 0
element carbon ionization 0 1
element oxygen ionization 0 0.1 0.2 0.3 0.2 0.1 0.1
c
c commands controlling continuum =====
blackbody 5
ionization parameter -2
c
c commands controlling geometry =====
c force model to be 1 cm thick

```

```

set dr 0
c
c other commands for details =====
constant temperature 4
stop zone 1
c
c commands controlling output =====
normalize to "O 3" 5007
punch dr "func_set_ion.dr"
c
c
c func_set_ion.in

c class function c ===== c this script exercises the option
to specify the ionization of a species

```

---

### **func\_stoplevel1 *test stop line command***

```

title test stop line command
c
c commands controlling continuum =====
ionization parameter -2
blackbody 50000
c
c commands for density & abundances =====
hden 2
abundances old solar 84
init file "ism.ini"
c test option to turn off elements with small abundances
element limit off -5
c
c commands controlling geometry =====
c stop calc at line ratio, the point of this test
stop line "C 2" 157.6m reaches 0.0358
c
c other commands for details =====
constant temperature 10000 K
c
c commands controlling output =====
punch dr "func_stoplevel1.dr" last
c
c func_stoplevel1.in
c class function
c =====
c

```

This is an example of a simple calculation that stops when a line reaches a specified intensity. The option to turn off elements with trivial abundances is used.

---

### **func\_stoplevel2 *test stop line command***

```

title test stop line command
c
c commands controlling continuum =====
ionization parameter -2
blackbody 50000

```

```

c
c commands for density & abundances =====
hden 2
abundances old solar 84
init file "ism.ini"
c
c commands controlling geometry =====
c stop calc at line ratio, the point of this test
stop line "C 2" 157.6m reaches 0.002 relative to "o 3" 5007 A
c
c other commands for details =====
constant temperature 10000 K
c
c commands controlling output =====
normalise to "o 3" 5007
punch dr "func_stopleveline2.dr" last
c
c func_stopleveline2.in
c class function
c =====
c

```

this is an example of a simple calculation that stops when a certain emission line ratio is reached

---

### ***func\_t10 test very soft continuum, very hot gas***

```

title test very soft continuum, very hot gas
c
c commands controlling continuum =====
table agn
ionization parameters -2
c
c commands for density & abundances =====
hden 1.
c
c commands controlling geometry =====
stop zone 1
c specify a thin cell of gas - 1 cm thick
set dr 0
c
c other commands for details =====
c set constant temperature to high-T limit
constant temperature, t=10 log
c iterate since gas is optically thin
iterate
c
c commands controlling output =====
c
c
c func_t10.in
c class limit
c =====
c

```

This is a test of the highest temperature the code can do.

---

### ***func\_t3 test low temperature limit of code, 3K***

```

title test low temperature limit of code, 3K
c
c commands controlling continuum =====
table agn
ionization parameters -2
c
c commands for density & abundances =====
hden -1
c
c commands controlling geometry =====
stop zone 1
c specify a thin cell of gas 1 cm thick
set dr 0
c
c other commands for details =====
c set constant temperature to low-T limit
constant temperature, t=3.0 linear
c iterate since gas is optically thin
iterate
c
c commands controlling output =====
c
c
c
c func_t3.in
c class limit
c =====
c

```

This is a test of the lowest temperature the code can do. It runs a constant temperature of 3K

---

## **func\_test** *run smoke test*

```

title run smoke test
c commands controlling continuum =====
test
c commands for density & abundances =====
c commands controlling geometry =====
c other commands for details =====
c include this in one test to make sure it works
set continuum shield ferland
c
c commands controlling output =====
c increase precision of all printed wavelengths
set line precision 6
c following will print physical constants used by the code
print constants
c print the path to the data directory
print path
c the citation for the current version of the code
print citation
c following prints column densities
print column density
punch overview "func_test.ovr"
c func_test.in
c class function
c =====
c

```

This runs the smoke test command, which include several asserts. The tests the behavior of increasing the number of significant figures in printed wavelengths.

---

## **func\_testmole** *this runs the standard, one command, test, which contains many asserts*

```

title this runs the standard, one command, test, which contains many asserts
c commands controlling continuum =====
test mole
c
c commands for density & abundances =====
c commands controlling geometry =====
c other commands for details =====
c
c commands controlling output =====
punch overview "func_testmole.ovr"
c func_testmole.in
c class function
c =====
c

```

This runs the "test mole" command, which include several asserts.

---

## **func\_trans\_punch** *first of func\_trans\_punch/transread pair, punch continuum*

```

title first of func_trans_punch/transread pair, punch continuum
c
c commands controlling continuum =====
ionization parameter -1
table agn
c
c commands for density & abundances =====
hden 10
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
iterate
c
c commands controlling output =====
print incident continuum
punch ionization last "func_trans_punch.ion"
punch continuum last "func_trans_punch.con"
c this continuum will be used by transread.in
punch last transmitted continuum "func_trans_punch.trn"
c
c this contains a large number of asserts that both tests use
init file "trans.dat"
c
c func_trans_punch.in
c class function

```

```
c =====
c
```

func\_trans\_punch.in and transread.in are a pair of tests that check that the code can punch a transmitted continuum then read it.

This sim must come before func\_trans\_read since it generates the punch file needed by func\_trans\_read. Alphabetical order insures this.

## **func\_trans\_read *second of transpunch/transread pair, used transmitted continuum***

```
title second of transpunch/transread pair, used transmitted continuum
c the transmitted continuum produced by func_trans_punch
c
c commands controlling continuum =====
ionization parameter -1
table read "func_trans_punch.trn"
brems 6
ionization parameter -6
c
c commands for density & abundances =====
hden 10
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
iterate
c
c commands controlling output =====
c this contains a large number of asserts used by both commands
init file "trans.dat"
print incident continuum
punch ionization last "func_trans_read.ion"
c
c
c func_trans_read.in
c class function
c =====
c
```

func\_trans\_punch.in and func\_trans\_read.in are a pair of tests that check that the code can punch a transmitted continuum then read it.

## **grains\_conserve\_pp *test energy conservation with vastly optically thick dust***

```
title test energy conservation with vastly optically thick dust
c
c commands controlling continuum =====
black 500 STE
```

```

c
c commands for density & abundances =====
init "honly.ini"
hden 0
grains ism single graphite
c
c commands controlling geometry =====
stop Av 1000
sphere
stop temperature 100K
c
c other commands for details =====
iterate
constant temper 500 K
no grain gas collisional energy exchange
no molecules
cosmic rays background
c
c commands controlling output =====
punch grain temperature "grains_conserve_pp.grntem"
punch continuum units microns "grains_conserve_pp.con"
c
c
c grains_conserve_pp.in
c class limit
c =====
c

```

This tests that multiple absorption / reemission by dust conserves energy Energy density and grain temperatures should be exactly 500 K

---

## ***grains\_conserve\_sp test energy conservation with vastly optically thick dust, sphere geometry***

```

title test energy conservation with vastly optically thick dust, sphere geometry
c
c commands controlling continuum =====
black 500 STE
c
c commands for density & abundances =====
init "honly.ini"
hden 3.25 -2
grains ism single graphite
c
c commands controlling geometry =====
radius from 21 to 26
* stop Av 1000
sphere
stop temperature 100K
c
c other commands for details =====
iterate
constant temper 500 K
no grain gas collisional energy exchange
no grain electrons
no molecules
cosmic rays background -10
c
c commands controlling output =====

```

```

punch grain temperature "grains_conserve_sp.grntem"
punch continuum units microns "grains_conserve_sp.con"
c
c
c grains_conserve_sp.in
c class limit
c =====
c

```

This tests that multiple absorption / reemission by dust conserves energy in a spherical geometry. Radiation density temperature in last zone should be  $500 * \sqrt{1e21/1e26} = 1.5811 \text{ K}$

---

## **grains\_hot** *test temperature of gas and dust in high energy density environment*

```

title test temperature of gas and dust in high energy density environment
c
c commands controlling continuum =====
table power law
c intensity of incident radiation field set by its energy density
energy density temp 550 K
c
c commands for density & abundances =====
hden 5.0
init "ism.ini"
c Orion grains and abundances
abundances orion
c
c commands controlling geometry =====
stop zone 1
set drmax 10
c
c other commands for details =====
c
c commands controlling output =====
normalize to "O 8" 18.97A
punch overview "grains_hot.ovr" last
c
c
c grains_hot.in
c class limit
c =====
c

```

This tests the grains in an extreme condition - irradiation by an AGN near the illuminated face of the molecular torus. The gas is predominantly heated by the grain electron photo-ejection.

---

## **grains\_hot\_wd01** *test temperature of gas and dust in high energy density environment*

```

title test temperature of gas and dust in high energy density environment
c
c commands controlling continuum =====
table power law

```



```

energy density temp 550
c
c commands for density & abundances =====
hden 5.0
init "ism.ini"
abundances orion no grains
grains orion no qheat
c
c revert to WD01 treatment
no grain x-ray treatment
c
c commands controlling geometry =====
stop zone 1
set drmax 10
c
c other commands for details =====
c
c commands controlling output =====
normalize to "O 8" 18.97A
punch overview "grains_hot_wd01.ovr" last
c
c
c grains_hot_wd01.in
c class limit
c =====
c

```

This tests the grains in an extreme condition - irradiation by an AGN near the illuminated face of the molecular torus. The gas is predominantly heated by the grains. The grain treatment has been reverted to Weingartner & Draine, 2001, which is NOT appropriate for these conditions. It is however a good test whether the old treatment is not broken....

---

## **grains\_lte** *check that grains equilibriate at correct temp in ste limit*

```

title check that grains equilibriate at correct temp in ste limit
c this also tests that the code works when H He ionization is ZERO
c
c commands controlling continuum =====
black 3 lte
c
c commands for density & abundances =====
hden 5
eden 5
init file "hheonly.ini"
c need to normalize to some line, oxygen will provide some
element oxygen on
c ism grains with high dust to gas ratio
grain abund 1
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
iterate
c the equilibrium temp would not be correct since small H2 molecule
c does not go to ste, being an equivalent two-level system
constant temperature 3
c
c commands controlling output =====

```

```

normalize to "O 1" 63.17m
punch heating last "grains_lte.het"
punch cooling last "grains_lte.col"
c
c grains_lte.in
c class limit
c =====
c

```

This test irradiates a set of grains with a true blackbody in strict thermodynamic equilibrium. We expect the grains (and everything else) to equilibrate at the blackbody temperature. The gas temperature is forced to the radiation temperature because the current molecule network (based on ISM approximations) does not go to LTE in the high-radiation density limit. The calculation asserts that all grain temperatures are very close to the radiation temperature.

---

## **grains\_qheat** *cool atomic ISM with Si grain quantum heating*

```

title cool atomic ISM with Si grain quantum heating
c
c commands controlling continuum =====
table draine
extinguish 20
c
c commands for density & abundances =====
hden 0
c turns on grains with default abundance (log abundance ratio to ism = 0)
c and with quantum heating enabled by default
c (disabled with "no grains qheat" option)
grains 0. "silicate_ism_10.opc"
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
cosmic rays background
iterate
c
c commands controlling output =====
normalize to "C 2" 157.6m
print diffuse continuum
print line faint -2
print line column
punch grains temperature last "grains_qheat.qht"
punch continuum "grains_qheat.con" units microns last
punch two photon continuum "grains_qheat.2nu" units microns
c
c
c grains_qheat.in
c class limit
c =====
c

```

This sim produces dust emission with a Wein trail that is dominated by quantum heating emission.

---

## **grains\_temp** *test all grain species temperature*

```

title test all grain species temperature
c
c commands controlling continuum =====
black 50000
ionization parameter -2
c
c commands for density & abundances =====
hden 3
grains_temp "graphite_ism_01.opc"
grains_temp "silicate_ism_01.opc"
grains_temp "graphite_orion_01.opc"
grains_temp "silicate_orion_01.opc"
grains_temp "silicate_0m010.opc"
grains_temp "silicate_0m100.opc"
grains_temp "grey_ism_01.opc"
grains_temp "silicate_1m000.opc"
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
iterate
constant temper 4
c
c commands controlling output =====
c
c
c grains_temp.in
c class limit
c =====
c

```

This turns on all the grain species that are included in the distribution. A model of an ionized layer is done and the monitors confirm the resulting grain temperatures.

---

## **grains\_temp\_all** *test all grain species temperature*

```

title test all grain species temperature
c
c commands controlling continuum =====
black 50000
ionization parameter -2
c
c commands for density & abundances =====
hden 3
c this turns on orion-style distributed grains
grains orion
c this turns on ism distributed grains
grains ism
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
constant temper 4
iterate
c
c commands controlling output =====
punch dr "grains_temp_all.dr" last

```

```
punch grain charge "grains_temp_all.grnchr" last
punch grain temperature "grains_temp_all.grntem" last
punch grain drift velocity "grains_temp_all.grndft" last
c
```

```
c grains_temp_all.in c class limit c ===== c
```

This is a relatively quick test of grains. The Orion and ISM silicate and graphitic grains are turned on and their equilibrium temperature checked.

---

## **h2\_cr *H2 with background cosmic ray ionization***

```
title H2 with background cosmic ray ionization
c
c commands controlling continuum =====
c this is hot star SED
black 30000
intensity -5. range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c background CR ionization rate of 2.5e-17 s^-1,
cosmic rays, background
c
c commands for density & abundances =====
hden 5
c
c commands controlling geometry =====
set dr 0
stop zone 5
stop temperature 400
c
c other commands for details =====
c failures should not occur
failures 1
atom h2
iterate
c want clean test of CR physics so set constant kinetic temperature
constant temperature 500K
c must turn off line photoexcitation and Solomon process
c since we want to test cr heating and ionization
no induced processes
c
c commands controlling output =====
print column density
print ages
punch overview "h2_cr.ovr" last
punch h2 lines "h2_cr.lin" last all
punch h2 rates "h2_cr.rat" last
punch h2 column density "h2_cr.col" last
c
c
c h2_cr.in
c class limit
c =====
c
```

This test conditions of cosmic ray ionization. Solar abundances with no dust are assumed so this involves gas-phase chemistry alone. The Solomon process is disabled with the "no induced processes" command so H2 is mainly dissociated by cosmic rays. This forms a pair with h2\_cr\_grains, which does include grains.

---

## **h2\_cr\_grains** *background cosmic ray ionization by suprathermal electrons only*

```

title background cosmic ray ionization by suprathermal electrons only
c
c commands controlling continuum =====
c this is hot star continuum
black 30000
intensity -5. range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c background CR ionization rate should be 2.5e-17 s^-1,
c Tielens & Hollenbach 1985; McKee AstroPH 9901370
cosmic rays, background
c
c commands for density & abundances =====
abundances ism
hden 5
c
c commands controlling geometry =====
set dr 0
stop zone 5
stop temperature 400
c
c other commands for details =====
set h2 grain formation thermal pump
atom h2
iterate
constant temperature 500K
c must turn off line photoexcitation
c since we want to test cr heating and ionization
no induced processes
c
c commands controlling output =====
normalize to "H2 " 2.121m
print column density
print ages
punch overview "h2_cr_grains.ovr" last
punch h2 lines "h2_cr_grains.lin" last all
punch h2 rates "h2_cr_grains.rat" last
punch h2 column density "h2_cr_grains.col" last
punch grains H2rates "h2_cr_grains.h2r" last
c
c h2_cr_grains.in
c class limit
c =====
c

```

This tests conditions of cosmic ray ionization and grain formation pumping. Solomon process is turned off with the "no induced processes" command so cosmic rays are the main dissociation process. This forms a pair with h2\_cr which does not include grains, so relies only on gas-phase chemistry.

---

## **h2\_hminus** *H2 populations in H- dominated limit*

```

title H2 populations in H- dominated limit

```

```

c this is one of a pair with h2_solomon.in
c
c commands controlling continuum =====
c this is hot star continuum
black 30000
intensity -5. range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c background CR ionization rate should be 2.5e-17 s^-1,
c Tielens & Hollenbach 1985; McKee AstroPH 9901370
cosmic rays, background
c
c commands for density & abundances =====
c no grains in this sim so H2 forms by H- route
abundances ism no grains
hden 5
c
c commands controlling geometry =====
set dr 0
stop zone 5
stop temperature 400
c
c other commands for details =====
c default is Takahashi 2001 - test thermal distribution here
set h2 grain formation thermal pump
atom h2
iterate
constant temperature 500K
c
c commands controlling output =====
print column density
print ages
punch overview "h2_hminus.ovr" last
punch h2 lines "h2_hminus.lin" last all
punch h2 rates "h2_hminus.rat" last
punch h2 column density "h2_hminus.col" last
c
c h2_hminus.in
c class limit
c =====
c

```

This tests large H2 model in limit of H- formation and Solomon destruction.

---

## ***h2\_pdr\_leiden\_f1 low density and flux model 1***

```

title low density and flux model 1
atom H2 He collisions new
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 5
c this is to insure that no H-ionizing radiation strikes the cloud
extinguish 24
c
c commands for density & abundances =====
c hydrogen density
hden 3.
grains ism 1.16 no qheat

```

```

c
c commands controlling geometry =====
c
c other commands for details =====
c turn on the large H2 model
atom h2
c turn down the number of convergence failures - there should be none
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the gas kinetic temperature to a constant 50 Kelvin
constant temperature 50 linear
c
c commands controlling output =====
c default normalization line is Hbeta, which is not produced by this gas
normalize to "C 2" 157.6m
print line faint -4
punch overview "h2_pdr_leiden_f1.ovr"
punch line emissivity "h2_pdr_leiden_f1.ems"
H2 2.121m
12CO 647.2m
C 2 157.6m
O 1 63.17m
Si 2 34.81m
end of lines
punch leiden lines "h2_pdr_leiden_f1.lin"
punch leiden "h2_pdr_leiden_f1.lei"
punch dr "h2_pdr_leiden_f1.dr"
punch molecules "h2_pdr_leiden_f1.mol"
punch grain physical conditions "h2_pdr_leiden_f1.grn"
punch overview "h2_pdr_leiden_f1.ovr"
punch H2 lines "h2_pdr_leiden_f1.h2lin" all
punch H2 column density "h2_pdr_leiden_f1.h2col"
punch H2 populations matrix zone "h2_pdr_leiden_f1.pop"
punch H2 rates "h2_pdr_leiden_f1.rat"
c
c
c h2_pdr_leiden_f1.in
c class pdr
c =====
c

```

This is one of the tests in Rollig et al. 2007, A&A, 467, 187

---

## **h2\_solomon** *H2 populations in solomon dominated limit*

```

title H2 populations in solomon dominated limit
c this is a pair with h2_hminus
c
c commands controlling continuum =====
c this is hot star continuum
black 30000
intensity -5. range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c background CR ionization rate should be 2.5e-17 s^-1,
c Tielens & Hollenbach 1985; McKee AstroPH 9901370
cosmic rays, background
c

```

```

c commands for density & abundances =====
abundances ism
hden 5
c
c commands controlling geometry =====
set dr 0
stop zone 5
stop temperature 400
c
c other commands for details =====
c default is Takahashi 2001 - test thermal distribution here
set h2 grain formation thermal pump
atom h2
iterate
constant temperature 500K
c
c commands controlling output =====
c don't print so many lines
print line faint 10
print line column
print line sort wavelength
print column density
print ages
punch overview "h2_solomon.ovr" last
punch h2 lines "h2_solomon.lin" last all
punch h2 rates "h2_solomon.rat" last
punch h2 column density "h2_solomon.col" last
c
c h2_solomon.in
c class limit
c =====
c

```

This test H2 in case of grain formation and solomon destruction

---

## ***h2\_t2000 test large H2 molecule in shock-like conditions***

```

title test large H2 molecule in shock-like conditions
c
c commands controlling continuum =====
c there is a continuum but it is unimportant
c first continuum is FIR hot grain continuum produced in
c unmodeled HII Region
black 30000
intensity -5 range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c must provide source of ionization for chemistry
cosmic rays, background
c
c commands for density & abundances =====
hden 6
grains ism, abundance log 0.16 single
init file="ism.ini"
abundances he -1.01 c -3.52 n-8 o-3.30 ne-8 mg-5.89
continue si -6.10 s -5.10 cl=-7 ar-8 fe -6.60
c
c commands controlling geometry =====
stop zone 10
c force zone thickness to 1 cm
set dr 0

```



```

c need to lower stop temperature since gas is cooler than 4000K
stop temperature 3
c
c other commands for details =====
failures 2
atom h2
constant temperature 2000
iterate
c this stop Lyman line excitation of H, which is then photoionized
c by Balmer continuum
case b
c
c commands controlling output =====
normalize to "H2 " 2.121m
c
punch pdr "h2_t2000.pdr"
punch h2 rates "h2_t2000.rat"
punch h2 column density "h2_t2000.col"
punch h2 lines "h2_t2000.lin" all last
punch h2 populations "h2_t2000.pop" all last
punch raw continuum "h2_t2000.raw"
punch continuum "h2_t2000.con"
punch overview "h2_t2000.ovr"
c
c h2_t2000.in
c class limit
c =====

```

This is a collisionally dominated H<sub>2</sub> simulation. The temperature has been fixed at 2000K and the large molecule turned on. The calculation checks the returned value of the ortho to para densities. Cosmic rays and the incident continuum have little effect, the density is high, so the populations should be close to LTE.

---

## ***h2\_t500 test large H2 molecule in PDR-like conditions***

```

title test large H2 molecule in PDR-like conditions
c
c commands controlling continuum =====
c first continuum is FIR hot grain continuum produced in
c unmodeled HII Region
blackbody, t = 75 K
intensity -2.7 (total)
c this is hot star continuum
black 30000
intensity -5 range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c
c commands for density & abundances =====
hden 6
grains ism, abundance log 0.16 single, no qheat
init file="ism.ini"
abundances he -1.01 c -3.52 n-8 o-3.30 ne-8 mg-5.89
continue si -6.10 s -5.10 cl=-7 ar-8 fe -6.60
turbulence 2.7 km/sec
c
c commands controlling geometry =====
stop zone 10
c force zone thickness to 1 cm
sphere
set dr 0

```

```

c stop when gas is fully neutral
stop efrac -10
c stop when gas is cold
stop temperature 10 linear
c
c other commands for details =====
failures 2
atom h2
constant temperature 500
iterate
case b
cosmic rays background
c
c commands controlling output =====
normalize to "12CO" 863.0m
c uv lines are strongly pumped by stellar continuum, break out contribution
print line pump
print line optical depths
print line faint -2
print ages
c
punch pdr "h2_t500.pdr"
punch H2 rates "h2_t500.rat"
punch H2 lines "h2_t500.lin" all
punch H2 column density "h2_t500.col"
punch H2 populations "h2_t500.pop"
punch H2 levels "h2_t500.lev"
punch raw continuum "h2_t500.raw"
punch continuum "h2_t500.con"
punch overview "h2_t500.ovr"
c
table lines "LineList_PDR_H2.dat"
c h2_t500.in
c class limit
c =====

```

This is a dense molecular gas with background cosmic rays and the incident radiation field set to a small value. The lower levels are in LTE.

---

## ***h\_otsopen test ots, inward fractions for pure hydrogen, open geo, filling factor***

```

title test ots, inward fractions for pure hydrogen, open geo, filling factor
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
hden 5
init file "honly.ini"
c
c commands controlling geometry =====
stop efrac 0.05 %stop when 5 percent ionized
constant temper 7500
filling factor 0.01
c
c other commands for details =====
c must iterate to get inward line fractions, that are asserted below

```

```

iterate
diffuse ots
c not iterating does not affect the solution
no level2
no induced processes (OK)
c
c commands controlling output =====
print diffuse continuum
print line inward
punch overview "h_otsopen.ovr" last
punch dr "h_otsopen.dr" last
punch continuum "h_otsopen.con" last units angstrom
normalize to "Ca B" 4861 = 1 % normalize to exact result
c
c this is an open geometry, so do not expect H atom to exactly be case B
c due to Lyman line leakage. we iterated, so have a good estimate of
c inward and outward fractions, which we test
c h_otsopen.in
c class limit
c =====
c

```

This tests the total emission from a hydrogen Stromgren sphere using the OTS approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

## ***h\_otspp plane parallel conservation and hydrogenic emission for pure hydrogen***

```

title plane parallel conservation and hydrogenic emission for pure hydrogen
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
hden 5
init file "honly.ini"
c
c commands controlling geometry =====
sphere static (OK)
stop efrac 0.05 %stop when 5 percent ionized
c
c other commands for details =====
constant temper 7500
no level2
no induced processes (OK)
c needed to prevent 2s / Bal from becoming optically thick
turbulence 20 km/s
diffuse ots
atom h-like levels resolved 10
c
c commands controlling output =====
print diffuse continuum
normalize to "Ca B" 4861 = 1 % normalize to exact result for comp ion struc
punch overview "h_otspp.ovr"
punch dr "h_otspp.dr"
c
c h_otspp.in
c class limit

```

```
c =====
c
```

This tests the total emission from a plane parallel pure hydrogen Stromgren sphere using the OTS approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

## ***h\_otssp spherical conservation and hydrogenic emission for pure hydrogen***

```
title spherical conservation and hydrogenic emission for pure hydrogen
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
hden 5
init file "honly.ini"
c
c commands controlling geometry =====
radius 13
c not iterating does not affect the solution
sphere static (OK)
stop efrac 0.05 %stop when 5 percent ionized
c
c other commands for details =====
constant temper 7500
diffuse ots
no level2
no induced processes (OK)
c
c commands controlling output =====
print diffuse continuum
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "h_otssp.ovr"
punch dr "h_otssp.dr"
c
c h_otssp.in
c class limit
c =====
c
```

This tests the total emission from a spherical pure hydrogen Stromgren sphere using the OTS approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

## ***h\_outopen test open geometry***

```
title test open geometry
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
```

```

hden 0
init file "honly.ini"
c
c commands controlling geometry =====
stop efrac 0.05 %stop when 5 percent ionized
filling 0.01
c
c other commands for details =====
diffuse outward
iterate
atom h-like levels resolved 15
no level2
no induced processes (OK)
constant temper 4
c
c commands controlling output =====
print diffuse continuum
c normally will not print out inward parts of lines
print line inward
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "h_outopen.ovr" last
punch dr "h_outopen.dr" last
c
c this is an open geometry, so do not expect H atom to exactly be case B
c due to Lyman line leakage. we iterated, so have a good estimate of
c inward and outward fractions, which we test
c h_outopen.in
c class limit
c =====
c

```

This tests the total emission from an open geometry, hydrogen Stromgren sphere, using the outward only approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

## ***h\_outpp plane parallel H-only, close, test hydrogenic emission***

```

title plane parallel H-only, close, test hydrogenic emission
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
c the hydrogen density
hden 5
c an ini file that will include H-only
init file "honly.ini"
c
c commands controlling geometry =====
stop efrac 0.05 %stop when 5 percent ionized
c need this to prevent lyman lines from escaping
sphere static (OK)
c
c other commands for details =====
c prevent continuum pumping of lyman lines
no induced processes (OK)
c must include this to keep Ha optically thin
turbulence 20 km/s
c this is the approximation we are testing

```

```

diffuse outward
c force a constant temperature, then give stopping criteria
c if stop were not present model would go to limiting number of zones
constant temper 7500
c increase number of levels to get full emission
atom h-like levels resolved 20
c
c commands controlling output =====
c need to turn this on so that continua at various energies will
c be included in the emission line list
print diffuse continuum
c normalize results to perfect case b - our predictions should
c be close to this
normalize to line "Ca B" 4861
c the output and asserts
punch overview "h_outpp.ovr"
punch dr "h_outpp.dr"
c
c lots of continuum emission points turned on with print diffuse continuum
c energies dominated by two photon emission not included since very
c sensitive to Ly $\alpha$  transport
c h_outpp.in
c class limit
c =====
c

```

This tests the total emission from a plane parallel pure hydrogen Stromgren sphere using the outward only approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

## ***h\_outsp spherical conservation and hydrogenic emission for pure hydrogen***

```

title spherical conservation and hydrogenic emission for pure hydrogen
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
hden 5
init file "honly.ini"
c
c commands controlling geometry =====
radius 13
c not iterating does not affect the solution
sphere static (OK)
stop efrac 0.05 %stop when 5 percent ionized
c
c other commands for details =====
constant temper 7500
no level2
atom h-like levels resolved 20
no induced processes (OK)
diffuse outward
c
c commands controlling output =====
print diffuse continuum
normalize to "Ca B" 4861 = 1 % normalize to exact result

```

```

punch overview "h_outsp.ovr"
punch dr "h_outsp.dr"
c
c lots of continuum emission points turned on with print diffuse continuum
c energies dominated by two photon emission not included since very
c sensitive to Ly $\alpha$  transport
c h_outsp.in
c class limit
c =====
c

```

This tests the total emission from a spherical pure hydrogen Stromgren sphere using the outward only approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

## **h\_t4\_conemis** *continuous emission from H atom*

```

title continuous emission from H atom
c
c commands controlling continuum =====
c use laser so that stellar continuum is not in predicted continuum
laser 2
ionization -1
c
c commands for density & abundances =====
c high density to suppress two-photon emission
hden 8
init "honly.ini"
c
c commands controlling geometry =====
c this thickness is the inverse square of the density, so the
c total emission from the computed slab will in effect have the
c square of the density removed, so predictions will be  $j_{pi j} / n_e n_p$ 
stop thickness -16
c
c other commands for details =====
c include a large number of levels to capture IR emission
c increase default resolution stored in continuum_mesh.ini
set continuum resolution 0.1
iterate
constant temperature 4
c
c commands controlling output =====
print diffuse continuum
print lines column punch continuum h_t4_conemis.con last no units microns
punch continuum "h_t4_conemis.con" last no title units microns punch spectrum
h_t4_conemis.ncon last no units microns
punch spectrum "h_t4_conemis.ncon" last no title units microns
c
c h_t4_conemis.in
c class limit
c =====
c

```

This tests the continuous emission from the model H atom. The gas temperature is 10,000 K and the continuous emissivity is asserted for a range of wavelengths.

This was used to generate the plot in Hazy 2 comparing the emission from a pure hydrogen plasma with those

of Ferland 1980.

//>>refer HI emission Ferland, G. J. 1980, PASP, 92, 596

---

## **h\_t4\_conemis\_lon *low-den continuous HI emission with 2-nu important***

```

title low-den continuous HI emission with 2-nu important
c
c commands controlling continuum =====
laser 2
ionization -1
c
c commands for density & abundances =====
hden 0
init "honly.ini"
c
c commands controlling geometry =====
stop thickness 0
c
c other commands for details =====
c increase default resolution stored in continuum_mesh.ini
set continuum resolution 0.1
case b
iterate
constant temperature 4
c
c commands controlling output =====
print diffuse continuum
print lines column
punch continuum "h_t4_conemis_lon.con" last units microns
punch two photon continuum "h_t4_conemis_lon.2nu" last rydbergs
punch spectrum "h_t4_conemis_lon.ncon" last units microns
c
c h_t4_conemis_lon.in
c class limit
c =====
c

```

This is a mate to hatomt10.in except that the density is low enough for two-photon emission to be very imporant in the optical and uv.

---

## **h\_t4\_conemis\_thick *HI continuous emissivity, used for plot in hazy***

```

title H I continuous emissivity, used for plot in hazy
c
c commands controlling continuum =====
black 5
ionization parameter -2
c
c commands for density & abundances =====
c high density to suppress two-photon emission
hden 7
c turn off all elements, except for hydrogen, so that
c we really test hydrogen emission

```



```

init file= "honly.ini"
c
c commands controlling geometry =====
c this is necessary to stop model at H ionization front
c constant temperature models will go on to infinity without this
stop efrac -0.8
c
c other commands for details =====
turbulence 200 km/s ; needed to keep balmer lines thin
c increase continuum resolution by factor of ten
set continuum resolution 0.1
c line fluorescence would be very important with above turbulence,
c turn it off
no induced processes
c need large atom to get the many Rydgerg levels in the infrared
c atom h-like levels resolved 10
c what full outward flux so use spherical geometry
sphere
constant temperature 4
c
c commands controlling output =====
c results will be relative to interplated Hummer&Storey Hbeta
c this tests hydrogen atom too
normalize "Ca B" 4861
c by default, the continuum flux at many wavelengths is not printed,
c this turns that information on
print diffuse continuum
c output the predicted continuum
punch emitted continuum last iteration "h_t4_conemis_thick.con" units micron
punch dr last "h_t4_conemis_thick.dr"
c
c h_t4_conemis_thick.in
c class limit
c =====
c

```

This checks that the predicted hydrogen continuum is in good agreement with exact results in the optically thin nebular limit.

Checks:

- This output was used to generate figure h\_t4\_conemis\_thick in Part I of HAZY.
- Continuum relative to Hbeta should agree with Ferland (1980) filter averaged results.
- Hbeta should agree with Case B predictions, and Q(H) 4861.

## **heatomt10 *continuous emission from HeI***

```

title continuous emission from HeI
c
c commands controlling continuum =====
laser 1.9 ryd
ionization -1
c
c commands for density & abundances =====
c need high density to supress two-photon emission
hden 10
c force high density for electrons, He+ density will be the same
set eden 12
c set He/H ratio to 100, He and elec will have same density

```

```

element abundance helium 2
init "hheonly.ini"
c
c commands controlling geometry =====
c this thickness will get unit emission when combined with above densities
stop thickness -24
c
c other commands for details =====
iterate
constant temperature 4
c
c commands controlling output =====
print continuum
c
set continuum resolution 0.1
set nFnu diffuse outward, diffuse inward
print continuum
print line faint 1
print line column punch spectrum heatomt10.spc last no units microns
punch spectrum "heatomt10.spc" last no title units microns punch continuum heatomt10.con
last no units microns
punch continuum "heatomt10.con" last no title units microns punch diffuse continuum
heatomt10.dif last no units microns
punch diffuse continuum "heatomt10.dif" last no title units microns
c
c heatomt10.in
c class limit
c =====
c

```

This tests continuous emission from the He I atom. The laser is used so that the incident continuum is not included in the total emission.

---

## **heatomt10lon *test low-den continuous emission from H atom, 2-nu is important***

```

title test low-den continuous emission from H atom, 2-nu is important
c
c commands controlling continuum =====
laser 2
ionization -1
c
c commands for density & abundances =====
hden -8
element helium abundance log 3
set eden -5
init "hheonly.ini"
c
c commands controlling geometry =====
stop thickness log 10
c
c other commands for details =====
no scattering opacity
c force all helium to be in He+
element helium ionization -4 0 -4
case b
c increase default resolution stored in continuum_mesh.ini
set continuum resolution 0.1

```

```

iterate
constant temperature 4
c
c commands controlling output =====
normalize to "he 1" 4471
print lines column
print diffuse continuum
print lines column
punch continuum "heatomt10lon.con" last units microns
punch diffuse continuum "heatomt10lon.dif" last units microns
punch two photon continuum "heatomt10lon.2nu" last
punch spectrum "heatomt10lon.ncon" last units microns
c
c heatomt10lon.in
c class limit
c =====
c

```

This is a mate to hatomt10.in except that everything is He at low density

---

## **heiont10 *continuous emission from HeII***

```

title continuous emission from HeII
c
c commands controlling continuum =====
laser 5 ryd
ionization 1
c
c commands for density & abundances =====
hden 8
set eden 10
element abundance helium 2
init "hheonly.ini"
c
c commands controlling geometry =====
stop thickness -20
c
c other commands for details =====
constant temperature 4
iterate
c increase continuum resolution by factor of ten
set continuum resolution 0.1
set nFnu diffuse outward, diffuse inward
c
c commands controlling output =====
print continuum
print line faint 1
print line column punch continuum heiont10.con"last no units microns
punch continuum "heiont10.con" last no title units microns punch spectrum heiont10.spc"last
no units microns
punch spectrum "heiont10.spc" last no title units microns
c
c heiont10.in
c class limit
c =====
c

```

This tests the He II continuous emission. The helium abundance is very large so that He II overwhelms other emission sources. The resolution of the continuum mesh is increased so that we get a better representation of

the continuous emission.

---

## helike\_ar *He-like argon emission*

```

title He-like argon emission
c
c commands controlling continuum =====
laser 600
ionization -1
c
c commands for density & abundances =====
hden 7
c include lots of the element we are testing and set the ionization
c to be dominated by the correct ion stage
element argon abundance -2
element argon ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element nitrogen off
element carbon off
element neon off
element magnesium off
element silicon off
element sulphur off
element oxygen off
element iron off
c
c commands controlling geometry =====
set dr -12
stop zone 1
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print he-like departure argon
print line faint -2
c This is Lyman alpha
normalize to "Ar17" 3.949A
c
c 3.994A is 2^3S to ground
c Totl is 2^3P to ground
c 3.365A is 3^1P to ground
c 21.54A is 3^3P to 2^3S
c 22.24A is 3^3D to 2^3P
c
c helike_ar.in
c class limit
c =====
c

```

test He-like emission for argon

---

## helike\_c *he-like carbon emission*

```

title he-like carbon emission
c
c commands controlling continuum =====
laser 50
ionization -1
c
c commands for density & abundances =====
hden 7
c include lots of the element we are testing and set the ionization
c to be dominated by the correct ion stage
element carbon abundance -2
element carbon ionization -5 -5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element nitrogen off
element oxygen off
element neon off
element magnesium off
element silicon off
element sulphur off
element argon off
element iron off
c
c commands controlling geometry =====
set dr -12
stop zone 1
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print he-like departure carbon
print line faint -2
c This is Lyman alpha
normalize to "C 5" 40.27A
c
c
c helike_c.in
c class limit
c =====
c

```

test he-like carbon emission

---

## helike\_co *He-like cobalt emission*

```

title He-like cobalt emission
atom he-like levels 10
atom he-like collapsed levels 20
print he-like departure cobalt
constant temper 5
iterate
hden 7
set dr -12
init file "ism.ini"
c must turn on since ism.ini turned it off
element cobalt on

```

```

c then set abundance and ionization after turning it on
element cobalt abundance -2
element cobalt ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5
-5 -5 -5 -5 -5 0 -5
c now turn lots of elements off to save time
element helium off
element nitrogen off
element carbon off
element neon off
element magnesium off
element silicon off
element sulphur off
element oxygen off
element argon off
element iron off
laser 1300
ionization -1
stop zone 1
print line faint -2
c
c helike_co.in
c class limit
c =====
c

```

test emission of He-like Co

---

## helike\_cu *He-like copper emission*

```

title He-like copper emission
atom he-like levels 10
atom he-like collapsed levels 20
print he-like departure copper
constant temper 5
iterate
hden 7
set dr -12
init file "ism.ini"
c must turn on since ism.ini turned it off
element copper on
c then set abundance and ionization after turning it on
element copper abundance -2
element copper ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5
-5 -5 -5 -5 -5 -5 0 -5
c now turn lots of elements off to save time
element helium off
element nitrogen off
element carbon off
element neon off
element magnesium off
element silicon off
element sulphur off
element oxygen off
element argon off
element iron off
laser 1300
ionization -1
stop zone 1
print line faint -2
c
c helike_cu.in

```

```

c class limit
c =====
c

```

test emission of He-like Cu

---

## **helike\_fe** *he-like iron emission*

```

title he-like iron emission
c
c commands controlling continuum =====
laser 1300
ionization -1
c
c commands for density & abundances =====
hden 7
set dr -12
c include lots of the element we are testing and set the ionization
c to be dominated by the correct ion stage
element iron abundance -2
element iron ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5
-5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element nitrogen off
element carbon off
element neon off
element magnesium off
element silicon off
element sulphur off
element oxygen off
element argon off
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print line faint -2
print he-like departure iron
c This is Lyman alpha
normalize to "Fe25" 1.850A
c
c
c helike_fe.in
c class limit
c =====
c

```

check He-like emission for iron

---

## **helike\_mg** *he-like magnesium emission*

```

title he-like magnesium emission
c
c commands controlling continuum =====
laser 250
ionization -1
c
c commands for density & abundances =====
hden 7
set dr -12
element magnesium abundance -2
element magnesium ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element nitrogen off
element carbon off
element neon off
element silicon off
element oxygen off
element sulphur off
element argon off
element iron off
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print he-like departure magnesium
print line faint -2
c This is Lyman alpha
normalize to "Mg11" 9.169A
c
c
c helike_mg.in
c class limit
c =====
c

```

test He-like Mg emission

---

## **helike\_n *He-like nitrogen emission***

```

title He-like nitrogen emission
c
c commands controlling continuum =====
laser 75
ionization -1
c
c commands for density & abundances =====
hden 7
element nitrogen abundance -2
element nitrogen ionization -5 -5 -5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element carbon off

```



```

element oxygen off
element neon off
element magnesium off
element silicon off
element sulphur off
element argon off
element iron off
c
c commands controlling geometry =====
stop zone 1
set dr -12
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print he-like departure nitrogen
print line faint -2
c This is Lyman alpha
normalize to "N 6" 28.79A
c
c helike_n.in
c class limit
c =====
c

```

test He-like emission for N

---

## **helike\_ne** *he-like neon emission*

```

title he-like neon emission
c
c commands controlling continuum =====
laser 170
ionization -1
c
c commands for density & abundances =====
hden 7
c include lots of the element we are testing and set the ionization
c to be dominated by the correct ion stage
element neon abundance -2
element neon ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element nitrogen off
element carbon off
element oxygen off
element magnesium off
element silicon off
element sulphur off
element argon off
element iron off
c
c commands controlling geometry =====
stop zone 1
set dr -12
c
c other commands for details =====

```

```

atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print he-like departure neon
print line faint -2
c This is Lyman alpha
normalize to "Ne 9" 13.45A
c
c
c helike_ne.in
c class limit
c =====
c

```

test He-like emission for oxygen

---

## **helike\_ni** *he-like nickel emission*

```

title he-like nickel emission
atom he-like levels 10
atom he-like collapsed levels 20
print he-like departure nickel
constant temper 5
iterate
hden 7
set dr -12
init file "ism.ini"
c must turn on since ism.ini turned it off
element nickel on
c then set abundance and ionization after turning it on
element nickel abundance -2
element nickel ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5
-5 -5 -5 -5 -5 -5 0 -5
c now turn lots of elements off to save time
element helium off
element nitrogen off
element carbon off
element neon off
element magnesium off
element silicon off
element sulphur off
element oxygen off
element argon off
element iron off
laser 1300
ionization -1
stop zone 1
print line faint -2
c
c helike_ni.in
c class limit
c =====
c

```

Test He-like Ni emission.

---

## **helike\_o** *he-like oxygen ion vs. Bautista & Kallman 2000 Table 1, column 3*

```

title he-like oxygen ion vs. Bautista & Kallman 2000 Table 1, column 3
c
c commands controlling continuum =====
laser 100
ionization -1
c
c commands for density & abundances =====
hden 7
c include lots of the element we are testing and set the ionization
c to be dominated by the correct ion stage
element oxygen abundance -2
element oxygen ionization -5 -5 -5 -5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element nitrogen off
element carbon off
element neon off
element magnesium off
element silicon off
element sulphur off
element argon off
element iron off
c
c commands controlling geometry =====
stop zone 1
set dr -12
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print he-like departure oxygen
print line sort wavelength
print line faint -5
normalize to "o 7" 21.60A
c
c helike_o.in
c class limit
c =====
c

```

test He-like emission for oxygen

---

## **helike\_si** *He-like silicon emission*

```

title He-like silicon emission
c
c commands controlling continuum =====
laser 350
ionization -1
c

```

```

c commands for density & abundances =====
hden 7
c include lots of the element we are testing and set the ionization
c to be dominated by the correct ion stage
element silicon abundance -2
element silicon ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 0 -5
init file "ism.ini"
element helium off
element nitrogen off
element carbon off
element neon off
element magnesium off
element oxygen off
element sulphur off
element argon off
element iron off
c
c commands controlling geometry =====
stop zone 1
set dr -12
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 5
iterate
c
c commands controlling output =====
print he-like departure silicon
print line faint -2
c This is Lyman alpha
normalize to "Si13" 6.648A
c
c
c helike_si.in
c class limit
c

```

test He-like emission for silicon

---

## helike\_zn *He-like zinc emission*

```

title He-like zinc emission
atom he-like levels 10
atom he-like collapsed levels 20
print he-like departure zinc
constant temper 5
iterate
hden 7
set dr -12
init file "ism.ini"
c must turn on since ism.ini turned it off
element zinc on
c then set abundance and ionization after turning it on
element zinc abundance -2
element zinc ionization -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5 -5
-5 -5 -5 -5 -5 -5 -5 -5 0 -5
c now turn lots of elements off to save time
element helium off
element nitrogen off
element carbon off

```

```

element neon off
element magnesium off
element silicon off
element sulphur off
element oxygen off
element argon off
element iron off
laser 1300
ionization -1
stop zone 1
print line faint -2
c
c helike_zn.in
c class limit
c =====
c

```

test emission for He-like Zn

---

## ***hhe\_otspp plane parallel conservation and H-like emission for H, He***

```

title plane parallel conservation and H-like emission for H, He
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
hden 5
init file "hheonly.ini"
c
c commands controlling geometry =====
c expanding sphere to stop Ly $\alpha$  from destroying HeI 23S
sphere
stop efrac 0.05 %stop when 5 percent ionized, so just beyond H I-front
c
c other commands for details =====
diffuse ots
atom h-like element hydrogen levels resolved 10
c need to iterate since we will assert helium triplet lines
c that depend on having stable solution
iterate
constant temper 7500
no induced processes (OK)
c
c commands controlling output =====
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "hhe_otspp.ovr" last
punch dr "hhe_otspp.dr" last
c
c helium lines
c hhe_otspp.in
c class limit
c =====
c

```

This tests the total emission from a spherical pure H + He Stromgren sphere using the outward only approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

## ***hhe\_otssp spherical conservation and H-like emission for H and He***

```

title spherical conservation and H-like emission for H and He
c
c commands controlling continuum =====
blackbody 40000
ionization -2
c
c commands for density & abundances =====
hden 5
init file "hheonly.ini"
c
c commands controlling geometry =====
diffuse ots
atom h-like levels resolved 10
radius 13
stop efrac 0.05 %stop when 5 percent ionized
sphere static
c
c other commands for details =====
constant temper 7500
iterate
no level2
no induced processes (OK)
c
c commands controlling output =====
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "hhe_otssp.ovr" last
punch dr "hhe_otssp.dr" last
c
c helium lines
c hhe_otssp.in
c class limit
c =====
c

```

This tests the total emission from a spherical pure H + He-like Stromgren sphere using the outward only approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked. The geometry is plane paralel.

---

## ***hhe\_outpp plane parallel conservation emission for H, He gas***

```

title plane parallel conservation emission for H, He gas
c
c commands controlling continuum =====
blackbody 40000
ionization -2
c
c commands for density & abundances =====
hden 5
init file "hheonly.ini"
c
c commands controlling geometry =====
c open sphere to stop Ly $\alpha$  from destroying HeI 23S
sphere
stop efrac 0.05 %stop when 5 percent ionized
c
c other commands for details =====

```

```

c need this since will assert helium triplet lines that depend on stable soln
iterate
constant temper 7500
atom h-like element hydrogen levels resolved 10
no level2
no induced processes (OK)
c
c commands controlling output =====
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "hhe_outpp.ovr" last
punch dr "hhe_outpp.dr" last
c
c hhe_outpp.in
c class limit
c =====
c

```

This tests the total emission from a spherical pure hydrogen Stromgren sphere using the outward only approximation. The conservation of the total number of ionizing photons, and the emitted spectrum, are all checked.

---

### ***hhe\_outppff plane parallel filling factor pure H, He gas***

```

title plane parallel filling factor pure H, He gas
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
hden 5
init file "hheonly.ini"
c
c commands controlling geometry =====
filling factor -1
stop efrac 0.05 %stop when 5 percent ionized
sphere
c
c other commands for details =====
c need this since will assert helium triplet lines that depend on stable soln
iterate
constant temper 7500
no level2
no induced processes (OK)
atom h-like element hydrogen levels resolved 10
c
c commands controlling output =====
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "hhe_outppff.ovr" last
punch dr "hhe_outppff.dr" last
c
c helium lines
c hhe_outppff.in
c class limit
c =====
c

```

This is a plane-parallel constant temperature cloud with only hydrogen and helium. The gas has a filling factor of 0.1. Induced processes are turned off and a large H atom is used so that the hydrogen recombination

spectrum will be close to Case B. The calculation stops beyond the hydrogen ionization front, because of the stop efrac command (it needs this since this is a constant temperature calculation, so the usual lower-temperature stopping criterion does not apply). The asserts confirm that energy is conserved and that the hydrogen spectrum is correct.

---

## **hhe\_outsp** *spherical conservation and H-like emission for H, He*

```

title spherical conservation and H-like emission for H, He
c
c commands controlling continuum =====
ionization -2
blackbody 40000
c
c commands for density & abundances =====
hden 5
init file "hheonly.ini"
c
c commands controlling geometry =====
radius 13
stop efrac 0.05 %stop when 5 percent ionized
sphere static
c
c other commands for details =====
iterate
diffuse outward only
constant temper 7500
no level2
no induced processes (OK)
atom h-like element hydrogen levels resolved 10
c
c commands controlling output =====
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "hhe_outsp.ovr" last
punch dr "hhe_outsp.dr" last
c
c helium lines
c
c hhe_outsp.in
c class limit
c =====
c

```

This tests a spherical cloud with only hydrogen and helium. Diffuse fields are transferred with the outward only approximation. The asserts check that the ionizing radiation is conserved.

---

## **hii\_blister** *Lexington 1995 dust-free hii blister region*

```

title Lexington 1995 dust-free hii blister region
c
c commands controlling continuum =====
c a simple blackbody
blackbody 40000
c the flux of H-ionizing photons
phi(h) 13.0
c add the CMB after hot star shape & luminosity specified

```



```

CMB
c
c commands for density & abundances =====
hden 4
init file="ism.ini"
c HII region abundance but no dust
abundances hii region no grains
abundances hel c1 n1 o1 nel mg=.0001 si.01 s1 cl1 ar1
continue fe.001
c
c commands controlling geometry =====
sphere
c
c other commands for details =====
c this is to desaturate Lya to prevent excited state photoionization
c from being important - the other codes did not include this important
c physical process
turbulence 10 km/s
c must iterate since optically thick
iterate
c
c commands controlling output =====
normalize to "Ca B" 4861
print line faint .01
print line sum
He 1 5876
totl 2326
C 2 1335
TOTL 1909
N 2 6584
N 2 6548
N 3 57.21m
TOTL 7325
TOTL 3727
O 3 51.80m
O 3 5007
O 3 4959
Ne 2 12.81m
Ne 3 15.55m
Ne 3 3869
Ne 3 3968
S 3 18.67m
S 3 9532
S 3 9069
S 4 10.51m
end of lines
punch overview last "hii_blister.ovr"
punch dr last "hii_blister.dr"
punch results last "hii_blister.rlt"
print line optical depths
c
c
c hii_blister.in
c class hii
c =====

```

This is one of the test cases from the Lexington Meeting suite of nebulae (Ferland et al. 1995). It is a grain-free hii\_blister HII region, similar to inner regions of the Orion Nebula, except for the absence of grains. The set of lines entered with the print line sum command lists the most powerful coolants in this model. This is one of the tabulated quantities in the Lexington Meeting, and is a fundamental test of energy conservation in the code. The ratio of the sum of these lines to H $\beta$  is equivalent to the Stoy ratio, used for determining stellar temperatures.

The "dielec kludge 0" command is to turn off my estimates of the DR rates for those elements that had none. This was only to allow comparison with other calculations that did not make similar estimates. For an actual calculation I would not include this command, since the guesses are better than nothing.

the turbulence is to stop the balmer lines from becoming optically thick since few other codes include an actual H atom, but use case b instead. The Orion HII region does have an observed turbulence of about 8 km/s.

This calculation stops near the H<sup>+</sup> - H<sup>0</sup> ionization front, where the temperature falls below the default lowest temperature of 4000 K. This model would have continued into the PDR had a lower temperature been specified with the STOP LOWEST TEMP command.

## **hii\_coolstar *dust free cool HII region model, Lexington 1995***

```

title dust free cool HII region model, Lexington 1995
c
c commands controlling continuum =====
black body, T = 20000 K
q(h) 49
c add the CMB after hot star shape & luminosity specified
CMB
c
c commands for density & abundances =====
hden = 2
init file="ism.ini"
abund He-1 C-3.6576 N-4.39794 O-3.481146 ne-4.30103 mg-8
continue si-8 s-5.04576 cl-7 ar-8 fe-8
c
c commands controlling geometry =====
radius = 18.477121
sphere
c this is to go deep to pick up all H recombs
stop efract -2
stop temperature 1,000
c
c other commands for details =====
c must iterate since fine structure lines are optically thick
iterate
c assume there is neutral gas beyond what we compute
double
c
c commands controlling output =====
plot continuum range .1
print line faint .01
c this is the sum of lines in table 2 of the Lexington meeting
print line sum
N 2 6584
N 2 6548
TOTL 3727
Ne 2 12.81m
S 2 6720
S 3 18.67m
S 3 33.47m
S 3 9532
S 3 9069
end of lines
punch overview last "hii_coolstar.ovr"
punch dr last "hii_coolstar.dr"

```

```

punch results last "hii_coolstar.rlt"
punch continuum last units microns "hii_coolstar.con"
punch lines, cumulative, "hii_coolstar.cum"
totl 4861
o 3 5007
totl 3727
o 1 6300
end of lines
c
c Hbeta 4.93E36, L(total)4.30xHbeta
c hii_coolstar.in
c class hii
c =====
c

```

This is one of the test cases from the Lexington Meeting suite of nebulae. It is a grain-free HII region ionized by a very cool star. Hydrogen is ionized but not helium so this tests the transport of the H Lyman continuum. The set of lines is entered with the print line sum command to test energy conservation.

---

## ***hii\_icf HII region with negative He/H ICF***

```

title HII region with negative He/H ICF
c
c commands controlling continuum =====
c this is the result of this command
c table star mihalas 46000
c and makes it possible to run these orion sims without
c installing the stellar atmosphere files
table read "star_mihalas_46000.dat"
ionization parameter -2.
c add the CMB after hot star shape & luminosity specified
CMB
c
c commands for density & abundances =====
hden 3
init file "ism.ini"
abundances ism no grains
grains no qheat single
metals and grains 0.320
c
c commands controlling geometry =====
stop efrac -3
stop temperature 15
sphere
c
c other commands for details =====
failures 1
c
c commands controlling output =====
punch overview "hii_icf.ovr" last
punch dr "icr.dr" last
c
c hii_icf.in
c class hii
c =====

```

This is an example of an H II region irradiated by a hard stellar continuum - one of the Mihalas NLTE stars. The hard continuum produces a negative He/H ionization correction factor, as discussed in Ballantyne, Ferland & Martin (2000). >>refer HeI icf by Ballantyne,D.R., Ferland, G.J., & Martin, P.G., 2000, ApJ 536,

773-777

---

## hii\_paris *New'Paris meeting HII region*

```
title "New" Paris meeting HII region
c "standard" HII region model of the Pequignot Meudon Conferance
c
c commands controlling continuum =====
black body, T = 40000 K radius = 12.113943
c add the CMB after hot star shape & luminosity specified
CMB
c
c commands for density & abundances =====
hden = 2
init file="ism.ini"
abund he -1 C-3.6576 N-4.39794 O-3.481146 ne-4.30103 mg-8
continue si-8 s-5.04576 cl=-7 ar-8 fe-8
c
c commands controlling geometry =====
radius = 18.477121
sphere
c next two to make sure we pick up all possible H recombination
* stop temperature 600
set temperature floor 1000
c make sure sim goes very deep into cold gas
stop temperature 100K
stop efrac -2
c
c other commands for details =====
c must iterate since fine structure lines are optical thick
atom h-like element hydrogen levels resolved 10
iterate
c
c commands controlling output =====
normalize to "Ca B" 4861
print column densities
print line optical depths
plot continuum range .1
print line faint 0.01
print line sum
He 1 5876
totl 2326
TOTL 1909
N 2 121.7m
N 2 6584
N 2 6548
N 3 57.21m
TOTL 3727
O 3 5007
O 3 4959
O 3 51.80m
O 3 88.33m
Ne 2 12.81m
Ne 3 15.55m
Ne 3 3869
Ne 3 3968
S 2 6720
S 3 18.67m
S 3 33.47m
S 3 9532
S 3 9069
```

```

S 4 10.51m
end of lines
punch overview last "hii_paris.ovr"
punch hydrogen 21 cm last "hii_paris.21cm"
punch results last "hii_paris.rlt"
punch dr last "hii_paris.dr"
c this is to confirm that this command works
punch lines emissivity "hii_paris.str" last
TOTL 4861
end lines
c
table lines "LineList_HII.dat"
c hii_paris.in
c class hii
c =====
c

```

This is one of the "standard" models computed at the Paris and Lexington meetings on photoionization and shock calculations. a table in hazy compares the predictions of the current version of CLOUDY with predictions of a few of the other codes. It is necessary to iterate since some fine structure lines are optically thick. The set of lines entered with the print line sum command is used to obtain the total luminosity in detected lines, a measure of the Stoy temperature.

Checks:

- Hb close to case B, Q(H) 4861, intensities.
- Enter answers in Table Error! Reference source not found..

## **hlike\_c *H-like C VI case B***

```

title H-like C VI case B
c
c commands controlling continuum =====
table agn
ionization parameter 3
c
c commands for density & abundances =====
hden = 8
init file="ism.ini"
element carbon abundance 1
set eden 9
c
c commands controlling geometry =====
set dr -18
stop zone 1
c
c other commands for details =====
case b hummer no photoionization
no induced processes (OK)
constant temperature = 1.e5
c
c commands controlling output =====
punch results "hlike_c.rlt"
c
c hlike_c.in
c class limit
c =====
c

```

This test case compares the predictions of the multi-level H-like CVI atom with the Storey and Hummer (1995) results. The set dr command sets the zone thickness to 1 cm. The case b command sets Lyman line optical depths to very large values.

Checks:

- Neutral fractions
  - H\* emissivity
  - Relative line intensities
- 

## **hlike\_o** *H-like O VIII case B*

```

title H-like O VIII case B
c
c commands controlling continuum =====
table agn
ionization parameter 4
c
c commands for density & abundances =====
hden = 8
init file="ism.ini"
element oxygen abundance 1
set eden 9
c
c commands controlling geometry =====
c set following so the n^2 dr is unity
set dr -18
stop zone 1
c
c other commands for details =====
case b hummer no photoionization
no induced processes (OK)
constant temperature = 1.e5
c
c commands controlling output =====
punch results "hlike_o.rlt"
c
c caseb_o.in
c class limit
c =====
c

```

This test case compares the predictions of the multi-level H-like O VIII atom with the Storey and Hummer (1995) results. The set dr command sets the zone thickness to 1 cm. The case b command sets Lyman line optical depths to very large values.

Checks:

- Neutral fractions
  - H\* emissivity
  - Relative line intensities
- 

## **igm\_lalpha** *Ly alpha forest cloud*

```

title Ly alpha forest cloud
c
c commands controlling continuum =====
c ionized by cosmic background

```

```

c plus quasar/starburst continuum
cmb z=2
table hm05 z = 2
c
c commands for density & abundances =====
hden -2
metals -1.5 ;reduce all heavy elements by 1.5 dex
c
c commands controlling geometry =====
stop neutral column density 15
double ; mimic two-sided photoionization
c
c other commands for details =====
iterate to convergence; must iterate since optically thin
c
c commands controlling output =====
print line faint -1
punch overview last "igm_lalpha.ovr"
punch dr last "igm_lalpha.dr"
punch results last "igm_lalpha.rlt"
c
c on 05 aug 29 changed from old background command to
c cmb 2 plus hm05 z=2 background - all results changed substantially
c igm_lalpha.in
c class igm
c =====
c

```

This example demonstrates the behavior of the code in the low-density limit. The ionizing source is the cosmic background at a redshift of  $z = 2$ . The cloud is assumed to be optically thin to ionizing radiation, in keeping with the Gunn-Peterson test, so the double command is included, and an iteration is performed to converge the optical depth scale. Continuum fluorescent excitation of lines is important because the cloud is optically thin.

---

## ***igm\_primal cloud with primordial abundances exposed to background at $Z=10$***

```

title cloud with primordial abundances exposed to background at Z=10
c
c commands controlling continuum =====
background, redshift= 10
c background due to agn from Haardt & Madau 1996 ApJ, 461, 20
c this sets both continuum shape and intensity, actually for z=2
table HM96 old
c
c other commands for details =====
c put in a velocity field to stop Lya from causing stability problems
turbulence 20 km/s
c do a second iteration to establish that it is optically thick downstream
iterate
c this sim should run cleanly - do not accept problems
failures 2
c
c commands for density & abundances =====
c this log of the hydrogen density
hden 5
c stored set of primordial abundances
abundances primordial

```

```

c
c commands controlling geometry =====
c we want to shielded face to not be exposed to empty space, but rather
c be just very deep in a dense cloud - this prevents lines from freely
c radiating from the shielded face
double optical depths
c set a lower temperature stopping criterion - default is 4000K
stop temperature 30
c
c commands controlling output =====
c don't want to print lots of faint hydrogen lines
print line faint -3
c some output files with info we want to save
punch temperature history "igm_primal.tem.his"
punch ionizing continuum "igm_primal.ion"
punch hydrogen conditions "igm_primal.mol"
punch heating "igm_primal.het"
punch dr "igm_primal.dr"
punch cooling "igm_primal.col"
punch overview "igm_primal.ovr"
punch results last "igm_primal.rlt"
c
c igm_primal.in
c class hii pdr igm
c =====

```

This is a high redshift cloud irradiated by the cosmic background and AGN light.  
H- absorption of the z=10 cosmic background is a MAJOR heating agent  
This model is very sensitive to treatment of Lya destruction, since resulting  
J-bar is major destruction process for H-, which is the H2 formation route.  
Lya ots oscillations could develop and are damped by not reevaluating H-  
photo rate after first n times, as in hmo1e.c

---

## ***igm\_z3 redshift 1000 recombination epoch***

```

title redshift 1000 recombination epoch
c
c commands controlling continuum =====
c a true blackbody radiation field, in strick thermo equilibrium
blackbody 3000 STE
c
c commands for density & abundances =====
hden 0
abundances primordial
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
case b
iterate
c
c commands controlling output =====
normalize lines to "Ca B" 4861
print ages
print departure coef
print line faint -5
c
c igm_z3.in
c class limit igm
c =====

```



c

This is a model of the universe near the recombination epoch, at a redshift of a thousand. The gas is exposed to a true blackbody at 3000 K, and the abundances are primordial.

---

## ***ism cloud irradiated by ism background***

```

title cloud irradiated by ism background
c
c commands controlling continuum =====
c this uses the Black continuum with no H-ionizing radiation,
c and with background cosmic rays
table ism
extinguish by a column of 22
c need cosmic rays to provide ionization to the chemistry
cosmic rays, background
c
c commands for density & abundances =====
hden 0
init file="ism.ini"
abundances ism
c
c commands controlling geometry =====
c set sphere since matter in all directions
sphere
stop temperature linear 10
stop thickness 0.1 linear parsecs
c
c other commands for details =====
c this prevents continuum pumping from exciting H0
case b
c this is to converge optical depths
iterate
c this should have no effect since T is larger than 100 K
set temperature floor 100K
c
c commands controlling output =====
c this model has no heavy element molecules, because H2 does not
c really fully form, since Lyman bands are not self-shielded
normalize to 157.6m "C 2"
print line pump
punch overview "ism.ovr" last
punch dr "ism.dr"
punch heating "ism.het"
punch coolign "ism.col"
c
c ism.in
c class ism
c =====

```

TODO - look at temperature structure - it has jitter at about the level of convergence, up and down. Temp jitter caused by eden jitter. This model is nearly isothermal, jitter measures noise in solver, and is great chance to pin this down.

This is a test of the behavior of the code in the extreme of photoionization by a relatively hard continuum, at low densities. The continuum is the galactic background, attenuated by a column density of  $10^{22}$  cm<sup>-2</sup>. Ionization by galactic background cosmic rays is included. Case b appears since this region is deep in the ISM, and the Lyman lines are quite thick. This example checks whether the ionization balance, thermal

balance, and electron density sum, are performed correctly in this limit.

Checks:

- Numerical stability of solution
- Thickness exact

## ***ism\_cosmicray background cosmic ray ionization by suprathermal electrons only***

```

title background cosmic ray ionization by suprathermal electrons only
c
c commands controlling continuum =====
c background CR H0 ionization rate should be 2.5e-17 s^-1,
c Williams et al. ApJ, 503, 689
cosmic rays, background
c this will be VERY faint black 50000
ionization parameter -25
black 50000
c
c commands for density & abundances =====
hden 5
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
iterate
set temperature floor 1000 K
c turn off molecules to test only ions
no molecules
c must turn off charge transfer and its heating, which would dominate,
c since we want to test cr heating and ionization
no charge transfer
no ctheat
c must turn of photoionization by recombining species,
c since we want to test cr heating and ionization
no photoionization
c any time photoionization is turned off, it is best to also
c turn off induced processes, which include pumping in the lyman lines
no induced processes
c
c commands controlling output =====
print ages
c
c ism_cosmicray.in
c class limit
c =====
c

```

This test conditions of cosmic ray ionization. Molecules and charge transfer are disabled so that analytical estimates can be made.

## ***ism\_grid interstellar cloud irradiated by ism background***

```

title interstellar cloud irradiated by ism background
c
c commands controlling continuum =====
c this uses the Black continuum with no H-ionizing radiation,
c and with background cosmic rays
table ism
extinguish by a column of 22
cosmic rays, background
c
c commands for density & abundances =====
hden 0 vary
grid from -3 to 2 in 0.5 dex steps
init file="ism.ini"
c this turns on ism abundances and grains
abundances ism
c
c commands controlling geometry =====
c set sphere since matter in all directions
sphere
stop temperature linear 10
stop thickness 0.1 linear parsecs
stop zone 1
c
c other commands for details =====
c this prevents continuum pumping from exciting H0
case b
c
c commands controlling output =====
c this model has no heavy element molecules, because H2 does not
c really fully form, since Lyman bands are not self-shielded
normalize to 157.6m "C 2"
punch grid "ism_grid.grd" last
punch averages "ism_grid.avr" last
temperature hydrogen 1
end of averages
punch overview "ism_grid.ovr" last
punch dr "ism_grid.dr"
punch heating "ism_grid.het"
punch coolign "ism_grid.col"
punch xspec mtable "ism_grid.fit" range 0.1 3 keV
punch temperature "ism_grid.tem"
c
c this is a special version of the assert that does a series of models
c sert tempe hydro 1 9942 .05 grid 8575 6468 3606 776 304 169 103 66.2 44.1 31.3
c
c ism_grid.in
c class ism
c =====

```

this shows an S-curve calculation - make plot showing density as X-axis and gas pressure (nT) as y-axis

the gas is ionized by the galactic background. the density varies between  $1e-3$  and  $100 \text{ cm}^{-3}$ . this is the full range found in the diffuse ism. The components that are produced are CNM - cold neutral medium, density  $\sim 40 \text{ cm}^{-3}$  WNM -  $n \sim 0.5 \text{ cm}^{-3}$ , WIM - warm ionized medium,  $n \sim 0.25 \text{ cm}^{-3}$  HIM - hot ionized medium,  $n \sim 1e-3 \text{ cm}^{-3}$ , calculation DOES NOT reproduce observed temperature of HIM - we get  $\sim 1e4\text{K}$  but observed is  $\sim 1e6 \text{ K}$ . HIM is shock, not photo, ionized

---

***ism\_hot\_brems generate continuum due to hot ism in high  $Z,z$  starburst***

```

title generate continuum due to hot ism in high Z,z starburst
c
c commands controlling continuum =====
c this sets up coronal equilibrium for temperature and continuum
c cloud is predominantly collisionally ionized
coronal equilibrium, t = 1.e6 K
c
c commands for density & abundances =====
c log of hydrogen density (cm-3), by default a constant density model
hden 1
c chemical composition for a well-evolved star cluster with high Z
c the starburst command generates scale factors that will be used to
c multiply the old solar composition, used in ver 84. this is maintain
c record of physical changes
abundances old solar 84
abundances starburst Z=10 Zsun
c
c commands controlling geometry =====
c the log of the total hydrogen column density, cm-2
stop column density 21
c
c other commands for details =====
c continuum is plotted in Hzay, we me must iterate to predict this
iterate
c
c commands controlling output =====
c normalize to 0 7 Lya
normalize to "0 7" 21.60
c this shortens the printout somewhat
print line faint 1
c output options punch continuum last !sm_hot_brems.con'no title, units keV
punch continuum last "ism_hot_brems.con" no title, units keV
punch sulphur ionization "ism_hot_brems.sul"
punch overview last "ism_hot_brems.ovr"
c this is only to document the problem described above
punch dr last "ism_hot_brems.dr"
c
c ism_hot_brems.in
c class ism
c =====
c

```

This model generates a large column constant density cloud similar to the hot phase of the interstellar medium. The continuum is punched to generate one of the figures in Part 2 of Hazy.

There is a strange feature between  $1.7e-3\text{\AA}$  and  $2.2e-3\text{\AA}$  that is the N emission (head starting at  $1.7e-3\text{\AA}$  with O absorption at  $2.2e-3\text{\AA}$ ). This model is strongly enriched in heavies so many metal edges, esp O, are optically thick.

>>TODO 1 the guess of the thickness of the first zone is badly too small, because this model is collisionally ionized, and it used Stromgren length - better to use collisional balance and dr - as result of this the model takes far too many zones

## **ism\_jura *check rate H2 forms on grain surfaces***

```

title check rate H2 forms on grain surfaces
c

```

```

c commands controlling continuum =====
c first continuum is FIR hot grain continuum produced in
c unmodeled HII Region
blackbody, t = 75 K
intensity 2.7 (total)
c this is hot star continuum
black 30000
intensity 1.90 range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
cosmic rays, background
c
c commands for density & abundances =====
hden 0
grains ism, abundance log 0.16 no qheat
init file="ism.ini"
abundances he -1.01 c -3.52 n-8 o-3.30 ne-8 mg-5.89
continue si -6.10 s -5.10 cl=-7 ar-8 fe -6.60
c
c commands controlling geometry =====
sphere
c stop when gas is fully neutral
stop efrac -10
c stop when gas is cold
stop temperature 10 linear
stop zone 1
c
c other commands for details =====
c set the gas temperature
constant temperature 100 K
case b
c
c commands controlling output =====
normalize to "C 2" 157.6m
c uv lines are strongly pumped by stellar continuum, break out contribution
print line pump
print line optical depths
print ages
punch hydrogen 21 cm last "ism_jura.21cm"
c
c ism_jura.in
c class limit
c =====
c

```

This model started out life as the Tielens & Hollenbach 1985 pdr. The density was set to unity and the incident radiation field adjusted so that the two default grains have temperatures near 100K. The model asserts that the H<sub>2</sub> formation rate on grain surfaces is close to the  $\gg$ refer H<sub>2</sub> grain physics Jura, M., 1975, ApJ, 197, 575 rate.

---

## ***ism\_opacity generate standard ISM opacity curve***

```

title generate standard ISM opacity curve
c
c commands controlling continuum =====
table agn
ionization parameter -3
c
c commands for density & abundances =====
grains no qheat

```

```

c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
iterate
c
c commands controlling output =====
punch total opacity last "ism_opacity.opc" units kev
c
c ism_opacity.in
c class limit
c =====
c

```

This example creates the file `ism_opacity.opc` which tabulates the total opacity of the gas as a function of energy. These plots are used in ISM studies to understand the transmission characteristics along a line of sight. The opacity depends on the dust to gas ratio, the gas phase abundances, and the level of ionization, all of which can be changed by altering parameters given above.

The model is of a 1 cm thick parcel of gas which is optically thin in the Lyman continuum and Lyman lines. As a result the hydrogen emission line spectrum is close to case C. The model iterates so that the predicted ionization and emission know about this.

## ***ism\_set\_cr\_rate background cosmic ray ionization by suprathemal electrons only***

```

title background cosmic ray ionization by suprathemal electrons only
c
c commands controlling continuum =====
cosmic ray rate -16
c this will be VERY faint black 50000
ionization parameter -25
black 50000
c
c commands for density & abundances =====
hden 5
element hydrogen ionization 1 0.00000001
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
constant temperature 1000 K
c turn off molecules to test only ions
no molecules
c must turn off charge transfer and its heating, which would dominate,
c since we want to test cr heating and ionization
no charge transfer
no ctheat
c must turn of photoionization by recombining species,
c since we want to test cr heating and ionization
no photoionization
c any time photoionization is turned off, it is best to also
c turn off induced processes, which include pumping in the lyman lines
no induced processes

```

```

c
c commands controlling output =====
print ages
c
c ism_set_cr_rate.in
c class limit
c =====
c

```

This test conditions of cosmic ray ionization. Molecules and charge transfer are disabled so that analytical estimates can be made.

---

## **limit\_casea\_h\_den13 case A**

```

title case A
c
c Seaton, M.J, 1959 MN 119, 90,
c  $4\pi j(\beta) 5.56E-26$ 
c for this model total H-beta=4.745
c  $b(2)=3.73E-3$  (3)= $3.69E-2$  (4)=0.091 (5)=0.145 (6)=0.193
c n.b. very different results if not l-mixed
c
c commands controlling continuum =====
black body, T = 50000 K
ionization parameter -2
c
c commands for density & abundances =====
hden = 13;
init file="honly.ini"
c
c commands controlling geometry =====
stop lyman optical depth -6
stop zone 1
set dr 0
c
c other commands for details =====
no induced processes (OK) ;N.B. case A is a fiction; no included transitions
atom h-like collisions off ; must turn off all collisions
constant temperature = 10000 K
iterate
c
c commands controlling output =====
c this should make the predicted totl Hbeta near unity for case a
normalize to "CaBo" 4861 = 2.23
print populations h-like
print departure coefficients
c
c limit_casea_h_den13.in
c class limit
c =====
c

```

Case A is a mathematical fiction; when the Lyman lines are optically thin continuum pumping must be important if the gas is ionized. Fluorescence is turned off with the no induced processes command. The density is set to a very high value ( $10^{15} \text{ cm}^{-3}$ ) so that the 2s-2p states are well l-mixed, in keeping with standard case A assumptions. As a result, collisional excitation would dominate the level populations, and hydrogen collisions must be turned off with the hydrogen collisions off command. The Ly\* optical depth is set to a small value. The set dr command sets the zone thickness to 1 cm. The abundances are set to a very small

value so that the electron density is equal to the hydrogen density.

Checks:

- Departure coefficients for H, He levels
- Neutral fractions
- H\* emissivity

## **limit\_casea\_h\_den\_temp** *Test model H in Case A limit*

```

title Test model H in Case A limit
c grid over nebular range of density and temperature to check Case A H
c
c commands controlling continuum =====
c create H+
laser 2 Ryd
ionization parameter -2
c
c commands for density & abundances =====
init "honly.ini"
hden 2. vary
grid 2 6 2
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
no scattering escape
case A hummer no Pdest no photoionization
no induced processes
constant temperature 4. vary
grid 5,000K to 20,000K in 5,000K steps
c
c commands controlling output =====
atom H-like print levels
c
c fixit the wavelength range is adjusted to avoid the Lyman lines
c the Lyman a / 2 photon ratio is wrong at high Tempetaure -
c is there a temperature dependent branching ratio to n=2 out
c of the collapsed level?
c
c limit_casea_h_den_temp.in
c class limit
c =====
c

```

this tests the predicted H I and He I spectra in the Case B limit. The grid is over both density and temperature.

## **limit\_caseb\_h\_hs87** *Case B from Hummer and Storey"*

```

title "Case B from Hummer and Storey"
init "honly.ini"
laser 1.1
ionization parameter -1
constant temperature 4
case b hummer no photoionization no pdest
atom h-like element hydrogen levels resolved 30

```



```

no induced processes
no level2
no scattering escape
set dr -20
hden 4
stop zone 1
iterate
print populaions h-like
print departure h-like
c
c
c limit_caseb_h_hs87.in
c class limit

```

---

## **limit\_caseb\_h\_lot *log density case B, T=500 log n=2***

```

title log density case B, T=500 log n=2
c
c commands controlling continuum =====
black body, T = 2.e5 K
ionization parameter -1
c
c commands for density & abundances =====
hden = 2
init file="honly.ini"
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature = 500
c want H and he fully stripped so we can check H He spectra
case b hummer no photoionization no pdest
no induced processes (OK)
iterate
c
c commands controlling output =====
print diffuse continua
normalize to "Ca B" 4861
c
c does not agree with Hummer & Storey, as explained by Ferguson & Ferland
c until 03 jun 02, used small atom, which had smaller differences
c turn on large atom to get best answer
atom h-like levels resolved 20
c
punch results "limit_caseb_h_lot.rlt" last
c

c limit_caseb_h_lot.in c class limit c ===== c

```

This tests the ionization and emission line spectrum for H case B at a low density and temperature.

---

## **limit\_caseb\_h\_n8 *h\_caseb\_n8 high density case B***

```

title h_caseb_n8 high density case B

```

```

c
c commands controlling continuum =====
black body, T = 2.e5 K
ionization parameter -1
c
c commands for density & abundances =====
hden = 8
init file="honly.ini"
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
c this will not affect the main model, but will test CR heating rate
cosmic ray background
case b hummer no photoionization no Pdest
no scattering escape
no induced processes (OK)
constant temperature = 20000
iterate
c
c commands controlling output =====
normalize to "Ca B" 4861
punch results "limit_caseb_h_n8.rlt"
punch ionizing continuum "limit_caseb_h_n8.cion"
c
c
c limit_caseb_h_n8.in
c class limit
c =====
c

```

This test case compares the predictions of the multi-level hydrogen atom with the Storey and Hummer (1995) results. The set dr command sets the zone thickness to 1 cm. The case b command sets Lyman line optical depths to very large values.

Checks:

- Neutral fractions
- H\* emissivity
- Relative line intensities

high density causes disagreement with HS - collisions

---

## **limit\_caseb\_he2\_den8 *limit\_caseb\_he2\_den8 He II case B***

```

title limit_caseb_he2_den8 He II case B
c
c commands controlling continuum =====
table agn
ionization parameter 0
c
c commands for density & abundances =====
hden = 8
element helium abundance 1
init file="ism.ini"
c
c commands controlling geometry =====
set dr 0
stop zone 1

```

```

c
c other commands for details =====
atom h-like element helium levels resolved 15
case b hummer no photoionization no Pdest
no induced processes (OK)
constant temperature = 50,000K
c
c commands controlling output =====
punch results "limit_caseb_he2_den8.rlt"
c
c
c limit_caseb_he2_den8.in
c class limit
c =====
c

```

This test case compares the predictions of the multi-level hydrogen atom with the Storey and Hummer (1995) results. The set dr command sets the zone thickness to 1 cm. The case b command sets Lyman line optical depths to very large values.

Checks:

- Neutral fractions
- H\* emissivity
- Relative line intensities

## **limit\_caseb\_he\_den** *Test model H and He atoms in Case B limit*

```

title Test model H and He atoms in Case B limit
c grid over nebular range of density and temperature to check Case B H & He
c
c commands controlling continuum =====
laser 5 Ryd
ionization parameter -2
c
c commands for density & abundances =====
init "hheonly.ini"
hden 2. vary
grid 2 6 2
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
normalize to "He 2" 4686
case b hummer no Pdest no photoionization
no scattering escape
c prevent H I excited levels from being predicted
atom H-like element hydrogen levels 3
constant temperature 4.
c
c commands controlling output =====
atom H-like print levels
c
c
c limit_caseb_he_den.in
c class limit
c =====
c

```

this tests the predicted He II spectra in the Case B limit.

this effectively turns off hydrogen to avoid the problem with every other heII line lying beneath an HI line. this is done by reducing the number of levels for H I.

this asserts the values are within 9% for the standard  $T = 1e4K$  and a range of densities. Actually they are all nearly within a few percent except at the lowest temperature of 5,000K.

## **limit\_caseb\_he\_den4\_temp4** *the best we can do to predict the HeI emission spectrum*

```

title the best we can do to predict the HeI emission spectrum
c
c commands controlling continuum =====
laser 2
ionization -1
c
c commands for density & abundances =====
hden 9000 linear
set eden 4
element helium abundance -1
init file "hheonly.ini"
c
c commands controlling geometry =====
stop zone 1
set dr -6.95424
c
c other commands for details =====
atom he-like levels 30
atom he-like collapsed levels 70
atom he-like gbar vriens
constant temper 4
case b no photoionzation no pdest
iterate
c
c commands controlling output =====
normalise to "He 1" 4471
print departure he-like helium
print line faint 0.01
c punch fits last "limit_caseb_he_den4_temp4.fit" punch diffuse continuum
limit_caseb_he_den4_temp4.dif"last no units microns
punch diffuse continuum "limit_caseb_he_den4_temp4.dif" last no title units microns
punch continuum limit_caseb_he_den4_temp4.con"last no units microns
punch continuum "limit_caseb_he_den4_temp4.con" last no title units microns
c
c
c limit_caseb_he_den4_temp4.in
c class limit
c =====
c

```

This is close to the best and most complete model of He I that the code can do. it is not hacked to try to reproduce HS?

## **limit\_caseb\_he\_den\_temp** *Test model He atoms in Case B limit*

```

title Test model He atoms in Case B limit
c grid over nebular range of density and temperature to check Case B H & He
c
c commands controlling continuum =====
laser 5 Ryd
ionization parameter -2
c
c commands for density & abundances =====
init "hheonly.ini"
hden 2. vary
grid 2 6 2
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
normalize to "He 2" 4686
case b hummer no Pdest no photoionization
no scattering escape
c prevent H I excited levels from being predicted
atom H-like element hydrogen levels 3
constant temperature 4. vary
grid 5,000K to 20,000K in 5,000K steps
c
c commands controlling output =====
atom H-like print levels
c
c
c limit_caseb_he_den_temp.in
c class limit
c =====
c

```

this tests the predicted He II spectra in the Case B limit.

this effectively turns off hydrogen to avoid the problem with every other heII line lying beneath an HI line. this is done by reducing the number of levels for H I.

this asserts the values are within 15% for a range of density and temperature. Actually they are all nearly within a few percent except at the lowest temperature. The error greater than 10% occurs at the lowest temperature of 5,000K.

## **limit\_caseb\_hhe\_den** *Test model H and He atoms in Case B limit*

```

title Test model H and He atoms in Case B limit
c grid over nebular range of density and temperature to check Case B H & He
c
c commands controlling continuum =====
c create H+ and He+
laser 2.0 Ryd
ionization parameter -2
c
c commands for density & abundances =====
init "hheonly.ini"
hden 2. vary
grid 2 6 2
c
c commands controlling geometry =====
stop zone 1

```

```

c
c other commands for details =====
no scattering escape
case b hummer no Pdest no photoionization
constant temperature 4.
c
c commands controlling output =====
atom H-like print levels
c
c
c limit_caseb_hhe_den.in
c class limit
c =====
c

```

this tests the predicted H I and He I spectra in the Case B limit.

the grid is over density at the standard temperature of 1e4 K.

---

### **limit\_caseb\_hhe\_den\_temp** *Test model H and He atoms in Case B limit*

```

title Test model H and He atoms in Case B limit
c grid over nebular range of density and temperature to check Case B H & He
c
c commands controlling continuum =====
c create H+ and He+
laser 2 Ryd
ionization parameter -2
c
c commands for density & abundances =====
init "hheonly.ini"
hden 2. vary
grid 2 6 2
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
no scattering escape
case b hummer no Pdest no photoionization
constant temperature 4. vary
grid 5,000K to 20,000K in 5,000K steps
c
c commands controlling output =====
atom H-like print levels
c
c
c limit_caseb_hhe_den_temp.in
c class limit
c =====
c

```

this tests the predicted H I and He I spectra in the Case B limit. The grid is over both density and temperature.

---

### **limit\_casec\_h\_den2** *H only optically thin in Lyman continuum*

```

title H only optically thin in Lyman continuum
c
c commands controlling continuum =====
blackbody 40000
ionization -2
c
c commands controlling geometry =====
stop thickness 16
c
c commands for density & abundances =====
hden 2
init file "honly.ini"
c
c other commands for details =====
iterate
constant temper 10000
no level2
c
c commands controlling output =====
print diffuse continuum
print line inward
normalize to "Ca B" 4861 = 1 % normalize to exact result
punch overview "limit_cassec_h-den2.ovr" last
punch fine optical depths last "limit_cassec_h-den2.fin" range 0.9 1.02 every 1
punch optical depths last "limit_cassec_h-den2.opt"
punch continuum "limit_cassec_h-den2.con" last units microns
punch dr "limit_cassec_h-den2.dr" last
c
c limit_cassec_h-den2.in
c class limit
c =====
c

```

This is a pure hydrogen cloud that is optically thin in the Lyman continuum. The asserts check the emission in several H I lines and continua. This should be close to what really happens in a low column density cloud exposed to a continuum source that does not have strong Lyman lines. (The continuum source used is a pure blackbody, and so has no lines). So this is an example of "Case C" emission >>refer H case C Ferland, G.J. 1999, PASP, 111, 1524

---

## **limit\_cassec\_h\_den5 case C**

```

title case C
c
c commands controlling continuum =====
black body, T = 50000 K
ionization parameter -2
c
c commands for density & abundances =====
hden = 5
init file="hheonly.ini"
c
c commands controlling geometry =====
set dr -10
stop zone 1
c
c other commands for details =====
iterate
stop lyman optical depth -6
constant temperature = 10000 K

```

```

c
c commands controlling output =====
normalize to "Ca B" 4861
print h-like populations
print departure coefficients
c
c limit_cassec_h_den5.in
c class limit
c =====
c

```

This is Case C, what really happens when optically thin gas is irradiated by a continuum with Lyman line continuum fluorescence allowed.

Checks:

- Departure coefficients for H, He levels
- Neutral fractions
- H\* emissivity Case C is described in

>>refer H case C Ferland, G.J. 1999, PASP, 111, 1524

## **limit\_compton\_hi\_t** *test high-T Compton energy exchange*

```

title test high-T Compton energy exchange
c
c commands controlling continuum =====
c as hot as STE bb can be on IEEE 32-bit cpu
c equilibrium temperature should also be 2.51e7
c since gas radiated by true blackbody will equilibrate
c at its temperature
blackbody 7.4 STE
c
c commands for density & abundances =====
hden 10
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
atom h-like element oxygen levels 5
atom h-like element oxygen levels collapsed 30
set temperature convergence 0.0001
iterate
print h-like departure element oxygen
c
c commands controlling output =====
c
c limit_compton_hi_t.in
c class limit
c =====
c

```

This is the highest Compton temperature that can be computed in LTE on an IEEE 32-bit processor. This tests the code in the high-temperature Compton limit. Temperatures as high as  $10^{10}$  K can be computed successfully on CPUs with longer word lengths, such as a Cray or the new 64 bit processors.

Checks:



- The equilibrium temperature should be exactly 107.4 K (2.51239107 K).

---

## **limit\_compton\_lo\_t *test low-T Compton energy exchange***

```

title test low-T Compton energy exchange
c
c commands controlling continuum =====
black linear 3 lte
brems 5
ionizat -5
c
c commands for density & abundances =====
hden -10
init file "hheonly.ini"
eden 0
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
set temperature convergence 0.0001
iterate 3
c
c commands controlling output =====
c
c limit_compton_lo_t.in
c class limit
c =====
c

```

This tests the code in the low temperature Compton limit. The gas is illuminated by a 3 K blackbody in thermodynamic equilibrium. The equilibrium temperature should be exactly 3 K. It is necessary to add an extra component of free electrons to test the code in this limit with the eden command.

---

## **limit\_compton\_mid\_t *mid-T Compton energy exchange***

```

title mid-T Compton energy exchange
c test of thermal equil in limit_compton_mid_t limit; temp should EXACTLY equal 2E5K
c check continuum partition;
c energy range, photon densities, luminosities, follow
c 0.25-1. Q=26.6470 L=15.8190 c 1-1.807 Q=26.8867 L=16.3766
c 1.807-4 Q=27.3042 L=17.0948 c 4-20 Q=27.2041 L=17.3179
c 20 -- Q=22.9038 L=22.9038 c total lumin 17.5597
c nufnu(912A) = 1.8029E+16
c
c commands controlling continuum =====
black body t = 2.e5 K lte
c
c commands for density & abundances =====
hden = 6
init file "hheonly.ini"
c
c commands controlling geometry =====
stop zone 1
stop lyman continuum optical depth = -6

```

```

set dr 0
c
c other commands for details =====
set temperature convergence .0001 % decrease error on heat-cool match
c
c commands controlling output =====
print line faint .1
print departure coef
c
c limit_compton_mid_t.in
c class limit
c =====
c

```

This tests the behavior of the code in the Compton limit. The incident continuum is a blackbody in strict thermodynamic equilibrium. Strict thermodynamic equilibrium is expected for all constituents of the gas. The input stream also lists the expected photon fluxes for the incident continuum; this tests the normalization of the continuum, and its distribution. Grains are included to confirm their behavior in the LTE limit. The set dr command sets the zone thickness to 1 cm.

Checks:

- Luminosity, photon flux, over various energy intervals, 4\*J at 912143.
- Electron temperature exactly 2\*105 K.
- Grain temperature forced to 2\*105 K by radiative processes.

## **limit\_conserve** *test that energy is limit\_conserv*

```

title test that energy is limit_conserv
* trace temperature convergence 0 2
c
c commands controlling continuum =====
c primary continuum is a hot blackbody
black body, temp=5.5
intensity total -4
c add cosmic background
backbround
c
c commands controlling geometry =====
c don't stop the calculation until all energy used up
c in order to check energy conservation
stop temperature 3 linear
c cloud will go very deep with temperature near 3K without this
stop thickness 20
c
c commands for density & abundances =====
no molecules
metals -0.5
hden 1.0
constant pressure
c
c other commands for details =====
iterate
c want a clean calculation
failures 3
c this allows for more molecular gas beyond what we model
double optical depths
c must increase number of levels since so optically thick

```

```

atom co levels 30
c
c commands controlling output =====
punch pressure "limit_conserve.pre"
punch pressure history "limit_conserve.his"
punch cooling "limit_conserve.col"
punch heating "limit_conserve.het"
punch dr "limit_conserve.dr"
punch overview "limit_conserve.ovr1"
punch overview last "limit_conserve.ovr"
punch results "limit_conserve.rlt"
punch element carbon "limit_conserve.car"
punch element neon "limit_conserve.ne"
punch element magnesium "limit_conserve.mag"
punch element silicon "limit_conserve.sil"
punch molecules "limit_conserve.mol"
c
c limit_conserve.in
c class hii pdr
c =====

```

This checks that energy is limit\_conserved. The code always checks that it did not radiate more energy than was absorbed. This calculation extends well past the photo-dissociation zone into fully molecular gas, so that all of the incident radiation is absorbed. Grains, CMB, & CRs are not present so that only the incident radiation field powers the gas.

Small changes can affect this model to surprising extents because of the presence of a major thermal front at the H0 - H+ transition region.

## **limit\_eden** *Martin Gaskell's funny model*

```

title Martin Gaskell's funny model
c used to test that electron density convergence is ok
c hydrogen line spectrum strongly pumped by continuum
c all elements > H have HIGH abundances
c ionization of elements Fe, Mg, Si strongly affected
c by charge transfer
c
c commands controlling continuum =====
black 4000
lumin 27.2
c
c commands for density & abundances =====
hden 5.138
abundances all 1000
c
c commands controlling geometry =====
radius 15
stop zone 1
set dr 0
c
c other commands for details =====
constant temper 5500
c
c commands controlling output =====
print line faint -3
c
c limit_eden.in
c class limit

```

```
c =====
c
```

This is mainly a test of the ability of the code to converge a model with a very strange electron density. The electrons are mainly contributed by heavy elements, and the gas is only slightly ionized.

Ionization of elements Fe, Mg, Si strongly affected by charge transfer with other heavy elements.

Checks:

- Electron density is correct.
- Hydrogen line spectrum strongly pumped by continuum.

## ***limit\_h\_induc constant temper black body limit from Ferland and Rees 1988***

```
title constant temper black body limit from Ferland and Rees 1988
c tests whether departure coef are forced to unity by induced processes
c
c commands controlling continuum =====
black body, t = 50000 lte
c
c commands for density & abundances =====
hden 10
init file="hheonly.ini"
c
c commands controlling geometry =====
stop lyman continuum optical depth -6
set dr -10
stop zone 1
c
c other commands for details =====
iterate
constant temper 50000
c
c commands controlling output =====
print populations h-like
print heating
print departure coef
print line faint -1
c
c limit_h_induc.in
c class limit
c =====
c
```

This example tests whether induced processes force level populations of hydrogen to LTE when they are irradiated by a blackbody in strict thermodynamic equilibrium. The density is low enough value for radiation to dominate the rate equations coupling levels with each other and the continuum. The expectation is for all departure coefficients to equal unity.

Checks:

- Departure coefficients exactly unity.
- Grain temperatures are exactly  $5 \times 10^4$  K.

---

## **limit\_hi\_ion** *very high ionization parameter limit*

```

title very high ionization parameter limit
c commands controlling continuum =====
table agn
ionization parameter 15
c commands controlling geometry =====
hden 0
stop zone 1
c other commands for details =====
iterate
c
c commands controlling output =====
punch heating "limit_hi_ion.het"
punch cooling "limit_hi_ion.col"
c
c
c func_hi_ion.in
c class function
c =====
c

```

This tests a limit of very high ionization

---

## **limit\_laser\_1** *test of H ionization in optically thin limit*

```

title test of H ionization in optically thin limit
c
c commands controlling continuum =====
laser 1.5 Ryd
phi(h) 10
c
c commands for density & abundances =====
hden 1
init file="hheonly.ini"
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
constant temperature = 4
iterate to convergence
c
c commands controlling output =====
c
c limit_laser_1.in
c class limit
c =====
c

```

This checks the calculation of the hydrogen photoionization equilibrium. The continuum is a laser peaked at 1.5 Ryd, where the hydrogen photoionization cross section is  $2.09 \times 10^{-18}$  cm<sup>2</sup>.

Checks:

- The hydrogen neutral fraction is nearly  $2.00 \times 10^{-4}$  (not exact since laser has finite width).
- Hb emissivity close to high density case A. The predicted TOTL 4861 intensity should be nearly 2.2 times the expected case B intensity.

H cross section is  $2.09 \times 10^{-18} \text{ cm}^2$ , rec coef is  $4.18 \times 10^{-14}$  answer is neutral fraction  $2.00 \times 10^{-4}$  also checks that only 3 iterations needed

---

## **limit\_laser\_2** *test of H and HeI ionization in optically thin limit*

```

title test of H and HeI ionization in optically thin limit
c
c commands controlling continuum =====
laser 2.0 Ryd
intensity -0.3604
c
c commands for density & abundances =====
hden 1
init file="hheonly.ini"
abundances all -10
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature = 4
iterate to convergence
c
c commands controlling output =====
c
c limit_laser_2.in
c class limit
c =====
c

```

This checks the calculation of the hydrogen and helium photoionization equilibrium. The continuum is a laser peaked at 2.0 Ryd, and so can only ionize hydrogen and atomic helium.

Checks:

- The hydrogen neutral fraction is nearly  $\text{H}_0/\text{H}^+ = 4.51 \times 10^{-4}$  (not exact since laser has finite width).
- Hb emissivity close to high density case A. The predicted TOTL 4861 intensity should be nearly 2.2 times the expected case B intensity.
- Helium ionization should be  $\text{He}_0/\text{He}^+ = 6.61 \times 10^{-4}$ .

H cross section is  $0.927 \times 10^{-18} \text{ cm}^2$ , rec coef is  $4.18 \times 10^{-13}$  answer is  $\text{H}_0/\text{H}^+ = 4.51 \times 10^{-4}$  HeI cross section is  $6.54 \times 10^{-18} \text{ cm}^2$ , rec coef is  $4.32 \times 10^{-13}$  answer is  $\text{He}_0/\text{He}^+ = 6.61 \times 10^{-5}$

---

## **limit\_laser\_200** *ionization in Auger-dominated limit*

```

title ionization in Auger-dominated limit
c
c commands controlling continuum =====

```

```

laser 200 Ryd
phi(h) 10
c
c commands for density & abundances =====
hden 1
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
iterate to convergence
constant temperature = 5
c
c commands controlling output =====
c
c limit_laser_200.in
c class limit
c =====
c

```

This checks the calculation of ionization equilibrium. The continuum is a laser peaked at 200 Ryd. It asserts ionization of C, O, and Fe. their ionization is dominated by the Auger effect.

Checks: Auger OK

---

### **limit\_laser\_200\_low *test ionization in Auger-dominated limit***

```

title test ionization in Auger-dominated limit
c
c commands controlling continuum =====
laser 200 Ryd
phi(h) 5
c
c commands for density & abundances =====
hden 1
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
iterate to convergence
constant temperature = 4
c
c commands controlling output =====
c
c
c
c
c
c limit_laser_200_low.in
c class limit
c =====
c

```

This checks the calculation of ionization equilibrium. The continuum is a laser peaked at 200 Ryd. It asserts ionization of C, O, and Fe. their ionization is dominated by the Auger effect.

Checks: Auger OK

---

### **limit\_laser\_3 test H and He ionization in optically thin limit**

```

title test H and He ionization in optically thin limit
c
c commands controlling continuum =====
laser 4.3 Ryd
phi(h) 10
c
c commands for density & abundances =====
hden 1
abundances all -10
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
iterate to convergence
constant temperature = 4
c
c commands controlling output =====
c
c limit_laser_3.in
c class limit
c =====
c

```

This checks the calculation of the hydrogen and helium photoionization equilibrium. The continuum is a laser peaked at 4.3 Ryd, where it can fully ionize both hydrogen and helium.

Checks:

- The hydrogen neutral fraction is nearly  $4.18 \times 10^{-4}$  (not exact since laser has finite width).
- Helium ion: The ratio  $\text{He}^+/\text{He}^{++}$  should be  $1.69 \times 10^{-3}$  and the ratio  $\text{He}^0/\text{He}^+$  should be  $2.86 \times 10^{-4}$ .
- Hb emissivity should be close to high-density case A. The predicted TOTL 4861 intensity should be nearly 2.2 times the expected case B intensity.

H cross section is  $1.0\text{E-}18 \text{ cm}^2$ , rec coef is  $4.18\text{E-}13$  answer is  $n(\text{H}^0)/n(\text{H}^+)=4.18\text{e-}3$  HeI cross section is  $1.51\text{E-}18 \text{ cm}^2$ , rec coef is  $4.32\text{e-}13$  answer is  $n(\text{He}^0)/n(\text{He}^+)=2.86\text{e-}4$ , so  $\text{He}^0/\text{He} = 4.83\text{e-}7$  HeII cross section is  $1.30\text{E-}18 \text{ cm}^2$ , rec coef is  $2.20\text{e-}12$  answer is  $n(\text{He}^+)/n(\text{He}^{2+})=1.69\text{e-}3$

---

### **limit\_lowd0 test low density limit**

```

title test low density limit
c this and limit_lowdm6 should get same results
c
c commands controlling continuum =====
table agn
ionization parameter -2
c

```



```

c commands for density & abundances =====
hden 0
abundances old solar 84
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
iterate
c
c commands controlling output =====
normalize to "H 1" 4861 0.86
print line sort wavelength range 100 1m
punch cooling last "limit_lowd0.col"
punch continuum last "limit_lowd0.con"
c
c all asserts for both sims are in this file
init file="lowd.dat"
c
c limit_lowd0.in
c class limit
c =====
c

```

this test case is paired with lowdm6.in both tests read in the same set of asserts, those contained in the file lowd.dat, and they should get exactly the same answer

this is also the test of the print lines intensity command

## **limit\_lowden** *test optically thin model that extends to very low densities*

```

title test optically thin model that extends to very low densities
c
c commands controlling continuum =====
table agn
ionization parameter 0
c
c commands for density & abundances =====
hden -2 to the -2 power
init file="ism.ini"
c
c commands controlling geometry =====
sphere
c this will take density down to 1e-8, below stated limit
radius 10 13
c
c other commands for details =====
c
c commands controlling output =====
punch overview "limit_lowden.ovr" last
c
c limit_lowden.in
c class limit
c =====
c

```

This model is optically thin, with density falling off as inverse square law, so ionization and temperature

should be nearly constant. if outer radius increased by 2 dex problem with level3 will appear, several li seq lines (OVI, NeVIII) will fluctuate when density about  $1e-9$

We do not assert H lines since the cloud is optically thin and takes at least three iterations to converge optical depth scale, That is not the purpose of this sim

---

## **limit\_lowdm6 *test low density limit***

```

title test low density limit
c this and limit_lowd0 should get same results
c
c commands controlling continuum =====
table agn
ionization parameter -2
c
c commands for density & abundances =====
hden -6
abundances old solar 84
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
iterate
c
c commands controlling output =====
normalize to "H 1" 4861 0.86
print line sort wavelength range 100 1m
punch cooling last "limit_lowdm6.col"
punch continuum last "limit_lowdm6.con"
c
c all asserts are in this file
init file="lowd.dat"
c
c limit_lowdm6.in
c class limit
c =====
c

```

this test case is paired with lowd0.in both tests read in the same set of asserts, those contained in the file lowd.dat, and they should get exactly the same answer

this also tests the print line sort range command

---

## **limit\_lowion\_low *test conditions of very low ionization matrix/simple solver***

```

title test conditions of very low ionization matrix/simple solver
c
c commands controlling continuum =====
blackbody 50000
ionization parameter -30
c

```

```

c commands for density & abundances =====
hden 2
init file "hheonly.ini"
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
atom h-like matrix lowt
atom he-like matrix lowt
no molecules
constant temperature 100
c
c commands controlling output =====
c
c limit_lowion_low.in
c class limit
c =====
c

```

The lowion\_pops.in and limit\_lowion\_low.in models form a pair that have identical boundary conditions but use the two different hydrogenic level populations solvers. The results should agree. lowion\_pops.in uses the full solution with the associate matrix inversion. This can fail under conditions of extreme low ionization due to numerical instabilities and roundoff. The solver used in limit\_lowion\_low.in is much simpler and will work for any conditions.

The model is almost totally molecular.

## **limit\_lowion\_pops *test conditions of very low ionization matrix/simple solver***

```

title test conditions of very low ionization matrix/simple solver
c
c commands controlling continuum =====
blackbody 50000
ionization parameter -30
c
c commands for density & abundances =====
hden 2
init file "hheonly.ini"
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
atom h-like matrix pops
atom he-like matrix pops
no molecules
constant temperature 100
c
c commands controlling output =====
c
c
c limit_lowion_pops.in
c class limit
c =====
c

```

The `limit_lowion_pops.in` and `lowion_low.in` models form a pair that have identical boundary conditions but use the two different hydrogenic level populations solvers. The results should agree. `limit_lowion_pops.in` uses the full solution with the associate matrix inversion. This can fail under conditions of extreme low ionization due to numerical instabilities and roundoff. The solver used in `lowion_low.in` is much simpler and will work for any conditions.

The model is almost totally molecular.

### `limit_lte_h_t50_cion` *collisionally ionized H in LTE limit*

```

title collisionally ionized H in LTE limit
c
c commands controlling continuum =====
black body, T = 50000 K
ionization parameter -8
c
c commands for density & abundances =====
hden = 20
set eden 20
init file="hheonly.ini"
c
c commands controlling geometry =====
stop lyman optical depth -6
set dr -10
stop zone 1
c
c other commands for details =====
constant temperature = 50000 K
c want pure collisional model
no induced processes (OK)
atom h-like collisional excitation off
atom h-like collisions l-mixing off
atom he-like collisional excitation off
atom he-like collisions l-mixing off
atom he-like collapsed levels 20
iterate
c
c commands controlling output =====
print h-like departure coefficients
print he-like departure coefficients
print populations h-like
print departure coefficients
c
c limit_lte_h_t50_cion.in
c class limit
c =====
c

```

This is the limiting case pure hydrogen collisional ionization, There are no excitation or l-mixing collisions, so this tests whether collisional ionization - three body recombination works in detailed balance.

### `limit_lte_h_t50_coll` *collisionally excited H in LTE limit*

```

title collisionally excited H in LTE limit
c

```

```

c commands controlling continuum =====
black body, T = 50000 K
ionization parameter -8
c
c commands for density & abundances =====
hden = 20
init file="hheonly.ini"
c
c commands controlling geometry =====
stop lyman optical depth -6
stop zone 1
set dr 0
set eden 20
c
c other commands for details =====
no induced processes (OK)
atom h-like collisions l-mixing off
atom he-like collapsed levels 20
atom he-like collisions l-mixing off
constant temperature = 50000 K
iterate
c
c commands controlling output =====
print h-like departure coefficients
print h-like populations
print he-like departure coefficients
print populations h-like
c
c limit_lte_h_t50_coll.in
c class limit
c =====
c

```

This checks that the model H atom goes to LTE at high densities.

chng 06 aug 24, had not included collisional ionization, and so he-like departure coefficients were very large, around 202. comments said there were problems. turned on collisional ionization, no problems noted

chng 06 jul 22 with RP changes in high-n n-changing collisions the rates are now much smaller - needed to change density to be far higher and several quantities changed. at lower density ( $1e18 \text{ cm}^{-3}$ ) the populations are very unphysical and runaway maser now occurs. this is only a homework problem and intended to only test n-changing collisions. with higher density this test is done.

## **limit\_lte\_he1\_coll** *He atom at high densities*

```

title He atom at high densities
c
c commands controlling continuum =====
laser 2
ionization -3
c
c commands for density & abundances =====
hden 18
set eden 18
element helium abundance -1
init file "hheonly.ini"
c
c commands controlling geometry =====
stop zone 1

```

```

set dr -10
c
c other commands for details =====
atom he-like levels 10
atom he-like collapsed levels 20
constant temper 20000 K
iterate
c
c commands controlling output =====
print he-like departure helium
normalise to "He 1" 4471
print line faint -2
c
c limit_lte_he1_coll.in
c class limit
c =====
c

```

test whether he-like ion populations go to lte in high density limit. The level populations should be in LTE, and the departure coefficients should be unity.

---

### **limit\_lte\_he1\_ste** *He atom in LTE at high densities*

```

title He atom in LTE at high densities
c
c commands controlling continuum =====
blackbody 50000 STE
c
c commands for density & abundances =====
hden 18
* set eden 18
c a high he abundance so that it dominates
element helium abundance 1
init file "hheonly.ini"
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
atom he-like levels 7
atom he-like collapsed levels 30
iterate
c
c commands controlling output =====
print he-like departure helium
print h-like departure hydrogen
normalise to "He 1" 4471
print line faint -2
punch heating "limit_lte_he1_ste.het"
punch cooling "limit_lte_he1_ste.col"
punch dr "limit_lte_he1_ste.dr"
c
c limit_lte_he1_ste.in
c class limit
c =====
c

```

test whether a gas dominated by He goes to LTE in high-density limit. The level populations should be in LTE, the departure coefficients should be unity, and the temperature equal to the BB temp.

---

## limit\_lte\_hhe\_coll\_t50 *H, He in LTE at high densities*

```

title H, He in LTE at high densities
c test from Ferland and Rees 88, collisions drive H to LTE
c collisions should drive all departure coef to unity
c
c commands controlling continuum =====
black body t = 50000
ionization parameter -5
c
c commands for density & abundances =====
hden = 19
init file="hheonly.ini"
set eden 20
c
c commands controlling geometry =====
set dr -10
stop zone 1
stop lyman optical depth -6
c
c other commands for details =====
constant temperature = 50000
iterate
c
c commands controlling output =====
print departure coefficients h-like
print populations h-like
print departure coefficients h-like helium
print populations h-like helium
print departure coefficients he-like
print populations he-like
punch dr "limit_lte_hhe_coll_t50.dr" last
c
c limit_lte_hhe_coll_t50.in
c class limit
c =====
c

```

This model is a test of the behavior of hydrogen and helium in the high density, collision dominated, limit. The temperature is preset, the hydrogen density is set to a very high value, and the ionization parameter is very low. The resulting model is collision dominated, so this case checks that the collision physics occurs in detailed balance. The predicted departure coefficients should all equal unity. The set dr command sets the zone thickness to 1 cm.

### Checks:

- Hydrogen departure coefficients exactly unity.
- Helium departure coefficients near unity. (Density not high enough to bring helium departure coefficients exactly to unity.)
- H-, H2, H2+ H3+, and HeH+ departure coefficients exactly unity. `

---

## limit\_lte\_hhe\_induc *radiation dominated H, He gas goes to STE*

```

title radiation dominated H, He gas goes to STE
c
c commands controlling continuum =====

```

```

black body, t = 50000 STE
c
c commands for density & abundances =====
hden 4
init file="hheonly.ini"
c
c commands controlling geometry =====
stop lyman continuum optical depth -6
set dr -10
stop zone 1
c
c other commands for details =====
iterate
set temperature convergence 0.002
c
c commands controlling output =====
print populations h-like
print heating
print departure coef
print line faint -1
c
c limit_lte_hhe_induc.in
c class limit
c =====
c

```

This is a H, He-only gas that is optically thin in the Lyman continuum. It is irradiated by a blackbody in strict thermodynamic equilibrium. The tests confirm that the gas temperature equilibrates close to the black body temperature.

---

## **limit\_lte\_hhe\_ste** *H, He in STE*

```

title H, He in STE
c from Ferland and Rees 1988
c this tests whether thermal processes go to STE
c
c commands controlling continuum =====
black body, t = 50000 STE
c
c commands for density & abundances =====
hden 10
init file="hheonly.ini"
c
c commands controlling geometry =====
stop lyman optical depth -6
set dr -10
stop zone 1
c
c other commands for details =====
iterate
c
c commands controlling output =====
print heating
print populations h-like
print departure coef h-like
print departure coef he-like
c
c limit_lte_hhe_ste.in
c class limit
c =====

```



c

This is the ultimate test of the behavior of the code in the strict thermodynamic equilibrium limit. The temperature is not held constant, so the resulting equilibrium temperature determines whether cooling processes are treated properly in the detailed balance limit. The equilibrium temperature should be exactly  $5 \times 10^4$  K, and all departure coefficients should equal unity. A small amount of grains are included to check that the grain thermal balance is handled properly in this limit.

Checks:

- Electron temperature exactly  $5 \times 10^4$  K.
- Departure coefficients unity.

## **limit\_lte\_hminus *H- goes to LTE***

```

title H- goes to LTE
c
c commands controlling continuum =====
c this is a second way to get STE and tests the energy density option
blackbody 5000 energy density 5000K
c
c commands for density & abundances =====
hden 10
c
c commands controlling geometry =====
stop zone 1
set dr 0
c
c other commands for details =====
iterate
constant temperature 5000
c
c commands controlling output =====
print populations h-like hydrogen
print departure coef
c
c limit_lte_hminusin
c
c class limit
c =====
c

```

This checks that the negative hydrogen ion goes to thermodynamic equilibrium when irradiated by a blackbody in thermodynamic equilibrium. It was originally presented in >>refer H- test Ferland, G. J., & Persson, S. E. 1989, ApJ, 347, 656

## **limit\_lte\_metal *STE with metals***

```

title STE with metals
c this tests whether thermal processes go to STE
c
c commands controlling continuum =====
c a dilution factor of 1 is one way to get STE
black body, t = 20000 dilution 1

```

```

c
c commands for density & abundances =====
hden 10
abundances starburst 5
c
c commands controlling geometry =====
stop zone 1
stop lyman optical depth -6
set dr -10
c
c other commands for details =====
iterate
c
c commands controlling output =====
print departure coef
print populations h-like
punch heating "limit_lte_metal.het" last
punch cooling "limit_lte_metal.col" last
c
c limit_lte_metal.in
c class limit
c =====
c

```

This checks that the code goes to strict thermodynamic equilibrium for the case of a metal rich gas exposed to a true black body. The many heavy element lines should dominate cooling, so this is a test that the multilevel atoms go to LTE in the radiation-dominated limit.

Checks:

- Temperature should equilibrate at 20000 K.
- Departure coefficients should equal unity.

## ***limit\_recoil\_ion test compton recoil ionization of hydrogen***

```

title test compton recoil ionization of hydrogen
c
c commands controlling continuum =====
laser 2000 ryd
ionization parameter -5
c
c commands for density & abundances =====
hden 0
init "honly.ini"
c
c commands controlling geometry =====
set dr 14
stop zone 10
c
c other commands for details =====
iterate
no secondary ionization
constant temperature 1000
c
c commands controlling output =====
c
c
c limit_recoil_ion.in
c class limit

```

```
c =====
c
```

H ionization is totally due to recoil ionization in this model. The assert checks the final hydrogen ionization.

---

## **limit\_strom** *pure-H Stromgren sphere*

```
title pure-H Stromgren sphere
c the answer is R(Stromgren) = 4.16E17 cm
c
c commands controlling continuum =====
blackbody 50000 K
q(h) 49
c
c commands for density & abundances =====
hden 4
init file "honly.ini"
abundances all -10
c
c commands controlling geometry =====
sphere static
radius 16
stop efrac -0.5
c
c other commands for details =====
no level2
turbulence 20 km/s
constant temper 7500
iterate
c
c commands controlling output =====
punch overview last "limit_strom.ovr"
punch results last "limit_strom.rlt"
punch dr last "limit_strom.dr"
c
c limit_strom.in
c class limit
c =====
c
```

This case checks that the code computes the geometry and emissivity correctly for a pure hydrogen spherical shell. The low temperature is chosen to avoid collisional ionization. The model stops at the Ho-H<sup>+</sup> ionization front. The turbulence is to prevent the Balmer lines from becoming optically thick.

Checks

- Outer radius should be 4.16391017 cm.
  - Predicted H $\beta$ , case B H $\beta$ , and Q(H) H $\beta$ , all agree.
- 

## **limit\_supra** *secondary ionization dominated gas*

```
title secondary ionization dominated gas
c like SN envelope
c
c commands controlling continuum =====
c this continuum will be VERY faint
table agn
```

```

ionization parameter -25
c secondary ionizations will dominate
set csupra -5
c
c commands for density & abundances =====
hden 5
abundances old solar 84
c
c commands controlling geometry =====
stop temperature 20
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature 1000 K
iterate convergence
c
c commands controlling output =====
print ages
c check that no uninitialized DR rates are printed (r2757)
print arrays carbon only
punch heating "limit_supra.het"
punch ionizing continuum "limit_supra.ion" last
punch molecules "limit_supra.mol"
c
c
c limit_supra.in
c class limit
c =====
c

```

This model computes the ionization within cool gas that is totally ionized by suprathermal secondary electrons.

Charge transfer heating is VERY important in this simulation.

## ***limit\_vbhum compare with Van Blerkom and Hummer exact RT results***

```

title compare with Van Blerkom and Hummer exact RT results
c
c commands controlling continuum =====
blackbody 50000 K
phi(h) 12.30103
c commands for density & abundances =====
hden 4
init "honly.ini"
abundances all -10
c
c commands controlling geometry =====
stop efrac -0.5
c
c other commands for details =====
diffuse ots
no level2 lines
constant temper 4
iterate
c
c commands controlling output =====

```

```

punch overview last "limit_vbhum.ovr"
punch dr last "limit_vbhum.dr"
c
c
c limit_vbhum.in
c
c class limit
c =====
c

```

This is a test of the treatment of the diffuse fields, their transfer, and their effects on the ionization structure of a nebula. The comparison is made against the exact calculation published by Van Blerkom and Hummer (1967). The geometry is open, that is, similar to that assumed in most BLR calculations.

>>refer H ionization Van Blerkom, D., & Hummer, D. G. 1967, MNRAS, 137, 353

The diffuse ots command is entered in order to reproduce the Van Blerkom and Hummer results. The default assumption, outward only, does not agree as well. I changed the default from OTS to outward only to be in better agreement with predictions by Harrington and Rubin at the Lexington meeting. They have not checked whether their codes are in agreement with the Van Blerkom and Hummer paper.

Checks:

- Neutral fraction at illuminated face  $5.8 \times 10^{-4}$ .
- Location of ionization front at  $7.8 \times 10^{16}$  cm.
- 34TOTL 486134 and 34CA B 486134 agree; both slightly lower than 34Q(H) 486134.
- Answers with OTS agree with 1967 results.

test hydrogen ground state rec effic against vb+h exact results this is their case e) - "zero condition" their answer for H0/Htot at the illuminated edge is approx  $5.8E-4$ , and a Stromgren radius of approximately  $7.7E16$  cm

## **limit\_veryveryfast** *very fast simulation for Purify/valgrind*

```

title very fast simulation for Purify/valgrind
c
c commands controlling continuum =====
table agn
ionization parameter -2
c
no file opacity
set dr 0
c
c commands for density & abundances =====
hden 9
init file "hheonly.ini"
element oxygen on
c
c commands controlling geometry =====
stop zone 1
c
c other commands for details =====
no level 2
constant temperature 4
atom h-like levels very small
atom he-like levels very small

```

```

c
c commands controlling output =====
punch dr last "limit_veryveryfast.dr"
c
c
c limit_veryveryfast
c class limit
c =====
c

```

This is meant to be a very fast calculation to use when running extensive debug-enabled runtimes.

---

## **nlr\_lex00** *NLR model for Lexington 2000 Meeting*

```

title NLR model for Lexington 2000 Meeting
c
c commands controlling continuum =====
phi(h) 12.47712
c shape of continuum is interpolated table
interpolate (0 -10) (0.08 -10) (0.1 1) (3676 -4.935) (4700 -11)
continue (4750 -20) (7.4e6 -30)
c
c commands controlling geometry =====
stop column 22
c
c commands for density & abundances =====
hden 4
init file="c84.ini"
abundances -1 c-3.52 n-4 o-3.097 ne-4 na-9
continue mg-4.523 al-8 si-4.523 s-4.824 a-9 ca-8 fe-8 ni-9
c
c other commands for details =====
iterate
c
c commands controlling output =====
* normalize to "totl" 4861 of 0.923
print line faint .01
print diffuse continuum
c add sum of these lines to printout for Stoy ratio
print line sum
TOTL 1216
He 1 5876
TOTL 1909
TOTL 1549
N 2 6584
N 2 6548
N 3 57.21m
totl 1750
totl 1486
o 1 6300
o 1 63.17m
TOTL 3727
totl 1665
O 3 51.80m
O 3 5007
O 3 4959
totl 1402
Ne 3 15.55m
Ne 3 3869
Ne 3 3968
Ne 4 2424

```

```

Ne 5 3426
Ne 5 3346
totl 2798
si 2 34.81m
s 2 6720
S 3 18.67m
S 3 9532
S 3 9069
S 4 10.51m
end of lines
c
punch overview last "nlr_lex00.ovr"
punch dr last "nlr_lex00.dr"
punch results last "nlr_lex00.rlt"
c
table lines "LineList_NLR.dat"
c nlr_lex00.in
c class nlr
c =====

```

This is one of the test cases from the Lexington (1993) Meeting suite of nebulae. It is a grain-free NLR model.

---

## **nlr\_liner** *NLR liner model*

```

title NLR liner model
c a constant pressure (gas+radiation) model of a nlr_liner cloud
c in the spirit of Ferland and Netzer 1983
c
c commands controlling continuum =====
table agn
ionization parameter -3
c
c commands for density & abundances =====
hden 6
init file "ism.ini"
abundances old solar 84
metals 0.3
c
c commands controlling geometry =====
stop thickness 16.3
stop column density 23
constant pressure
c
c other commands for details =====
iterate to convergence ;must iterate to get radiation pressure correctly
c
c commands controlling output =====
punch overview last "nlr_liner.ovr"
punch element nitrogen last "nlr_liner.nit"
punch dr last "nlr_liner.dr"
punch results last "nlr_liner.rlt"
c
c nlr_liner.in
c class nlr
c =====

```

This is a model somewhat like the Liner parameters proposed by Ferland and Netzer (1983). A second iteration is performed to allow the calculation of the line radiation pressure.

>>refer nlr\_liner model Ferland, G. J., & Netzer, H. 1983, ApJ, 264, 105

---

## **nlr\_liner\_grains** *liner model with grains*

```

title liner model with grains
c
c a constant pressure (gas+radiation) model of a liner cloud
c in the spirit of Ferland and Netzer 1983
c
c commands controlling continuum =====
ionization parameter -3
table agn
c
c commands for density & abundances =====
hden 6
c this turns on both grains and depleted abundances
abundances ism no grains
grains ism no qheat single
init file "ism.ini"
c
c commands controlling geometry =====
stop column density 23
c
c other commands for details =====
constant pressure
iterate to convergence ;must iterate to get radiation pressure correctly
c
c commands controlling output =====
punch overview last "nlr_liner_grains.ovr"
punch dr last "nlr_liner_grains.dr"
punch continuum last "nlr_liner_grains.con" units microns
punch results last "nlr_liner_grains.rlt"
c
c nlr_liner_grains.in
c class nlr
c =====
c

```

This is a model somewhat like the Liner parameters proposed by Ferland and Netzer (1983). A second iteration is performed to allow the calculation of the line radiation pressure. It includes grains and is so more realistic.

>>refer liner model Ferland, G. J., & Netzer, H. 1983, ApJ, 264, 105

---

## **nlr\_paris** *Paris meeting NLR model*

```

title Paris meeting NLR model
c
c commands controlling continuum =====
interpolate (0 -10) (0.073 -10) (0.074 1) (7352 -4)
continue (7400 -15) (7.353e6 -20)
ionization parameter -2
c
c commands controlling geometry =====
stop lyman continuum optical depth 4
c

```



```

c commands for density & abundances =====
hden 3
init file="c84.ini"
abund -1 C-3.5229 N-4. 0-3.22185 ne-3.82391 na--8 mg-4.5229 al--8
continue si-4.5229 s-4.82391 ar-8 ca--8 fe-7 ni-8
c
c other commands for details =====
c this large group of lines are turned off so that the model runs
c 30 percent faster - lines are mainly fuv and x-ray, which we will
c not track in this model
no level2
iterate
atom h-like element helium levels resolved 10
c
c commands controlling output =====
punch overview last "nlr_paris.ovr"
punch dr last "nlr_paris.dr"
punch results last "nlr_paris.rlt"
c
c nlr_paris.in
c class nlr
c =====

```

This is the NLR model presented in the Meudon meeting on model nebulae. The init file is entered to make the code behave more like version 84.

Checks:

- init file works

## **nova\_dqher** *cold nova shell*

```

title cold nova shell
c Ferland et al. 1984 DQ Her
c model of room temperature ionized cloud arund old nova DQ Her
c roughly that of
c tests behaviour of code at very low temperatures
c
c commands controlling continuum =====
c flat continuum, followed by nu^-2 power law
interpolate (0 0) (0.3 0) (8.e6 -14.8)
luminosity total 34
c
c commands for density & abundances =====
hden = 2.
abundances nova
grains no qheat single
c
c commands controlling geometry =====
filling factor 0.667
covering factor 0.667
radius 16.5682 , thickness = 16.14613
c need this since gas is so cool
stop temp = 100
c
c other commands for details =====
# >>chng 06 aug 02, add cosmic rays, should always have been included
cosmic rays background
iterate
turbulence 300

```

```

age 70 years %it exploded in the 1930's
atom h-like element hydrogen levels resolved 15
atom h-like element helium levels resolved 15
c
c commands controlling output =====
normalize to 4861 label="CaBo" intensity 1.169
print line all
punch overview last "nova_dqher.ovr"
punch results last "nova_dqher.rlt"
punch dr last "nova_dqher.dr"
punch transmitted continuum "nova_dqher.trn" last
c
c nova_dqher.in
c class nova
c =====
c

```

This tests the code39s behavior in the limit posed by the metal rich low density nebula surrounding DQ Her (Ferland et al. 1984).

Checks:

- Thickness exact
- Thermal stability High-Z gas ionization at low temperature

## **nova\_photos** *dense nova photosphere*

```

title dense nova photosphere
c
c commands controlling continuum =====
c this is the result of this command
c table star kurucz 35000; use old atlas atmosphere
c and makes it possible to run sims without
c installing the stellar atmosphere files
c set SED of central star
table read "star_kurucz_35000.dat"
c set luminosity of central star
absolute bolometric magnitude -8.1
c
c commands for density & abundances =====
hden 10 ;log of total hydrogen density (cm-3) at illuminated face
abundances old solar 84 ; use old solar abundances for record keeping
c
c commands controlling geometry =====
sphere ; model is a full sphere covering the central star
radius 12 to 14; log of inner and outer radii (cm)
c
c other commands for details =====
iterate to convergence
c
c commands controlling output =====
print line faint -1
punch overview "nova_photos.ovr" last
punch results "nova_photos.rlt" last
punch dr "nova_photos.dr" last
punch heat "nova_photos.het" last
punch cool "nova_photos.col" last
c
c nova_photos.in
c class nova
c =====
c

```

this model is intensely affected by continuum pumping of atoms. The hydrogen ionization is by Lyman line pumping, followed by photoionization from excited states.

---

## **optimize\_phymir** *test phymir optimizers*

```

title test phymir optimizers
c spectrum computed with hden 5, temp 4
c subplex, phymir, powell
c
c commands controlling continuum =====
black 40000 K
ionization par -3
c
c commands for density & abundances =====
hden 4 vary
init file "hheonly.ini"
element oxygen on
c
c commands controlling geometry =====
Case B
stop zone 1
c
c other commands for details =====
optimize tolerance 0.03
constant temper 4.3 vary
optimize phymir
optimize iteration 100
c the following spectrum was derived by running this sim at hden 5 and t=4
c resulting eden was 5.039 and T=4, which we shall assert that we find
c >>chng 06 apr 03, backdate to old [OII] As and rederive spectrum
optimize lines
o ii 3729 0.1432
o ii 3726 0.4055
o ii 7323 0.5281
o ii 7332 0.4257
o 3 5007 3.6189
totl 4363 0.0342
end of lines
c
c commands controlling output =====
print line faint -2
c
c
c optimize_phymir.in
c class optimizer
c =====
c

```

This checks whether the optimizer can recover a known solution. The line spectrum was calculated at  $T = 1e4$  K and  $n_H = 1e5 \text{ cm}^{-3}$ , and resulted in the given electron density. The model `optimize_subplex.in` is a copy of this file.

---

## **optimize\_subplex** *test subplex optimizer*

```

title test subplex optimizer
c spectrum computed with hden 5, temp 4

```

```

c subplex, phymir, amoeba, powell
c
c commands controlling continuum =====
black 40000 K
ionization par -3
c
c commands for density & abundances =====
hden 4 vary
init file "hheonly.ini"
element oxygen on
c
c commands controlling geometry =====
Case B
stop zone 1
c
c other commands for details =====
constant temper 4.3 vary
optimize subplex
optimize iteration 100
optimize tolerance 0.01
c the following spectrum was derived by running this sim at hden 5 and t=4
c resulting eden was 5.039 and T=4, which we shall assert that we find
c >>chng 06 apr 03, backdate to old [OII] As and rederive spectrum
optimize lines
o ii 3729 0.1432
o ii 3726 0.4055
o ii 7323 0.5281
o ii 7332 0.4257
o 3 5007 3.6189
totl 4363 0.0342
end of lines
c
c commands controlling output =====
print line faint -2
c
c
c optimize_subplex.in
c class optimizer
c =====
c

```

This checks whether the optimizer can recover a known solution. The line spectrum was calculated at  $T = 1e4$  K and  $n_H = 1e5 \text{ cm}^{-3}$ , and resulted in the given electron density. The model `optimize_phymir.in` is a copy of this file.

---

## **orion\_hii\_open *conditions similar to Orion nebula blister***

```

title conditions similar to Orion nebula blister
c orion hii region with open geometry
c
c commands controlling continuum =====
c SED of central star
table read "star_kurucz_39600.dat"
c its flux of H-ionizing photons
phi(h) 13.0
c
c this is to include velocity field in radiative transfer
turbulence 8 km/sec no pressure
c
c commands for density & abundances =====

```

```

hden 4
abundances hii region no grains
grains orion
c
c commands controlling geometry =====
cosmic rays, background
constant pressure
c this is the account for continued matter beyond i front
double optical depths
c
c other commands for details =====
iterate 2
c
c commands controlling output =====
print line faint -5
c print lines as surface brightness
print lines surface brightness arcsec
print line optical depths
print diffuse continuum
c want to print line intensities are surface brightness, per arcsec^2
print lines surface brightness arcsec
print diffuse continuum
punch overview last "orion_hii_open.ovr"
punch dr last "orion_hii_open.dr"
punch continuum last "orion_hii_open.con" units microns
punch transmitted continuum last "orion_hii_open.trn"
punch overview last "orion_hii_open.ovr"
punch fine opacity last "orion_hii_open.opc" range 0.7 0.9 ryd
punch element nitrogen last no hash "orion_hii_open.nit"
punch line emissivity "orion_hii_open.ems" last no hash emergent
totl 5199
o 1 6300
end of lines
punch linelist "orion_hii_open.lin" "LineList_HII.dat" last no hash emergent absolute
c
c orion_hii_open.in
c class hii
c =====

```

This is a model similar in spirit to the blister geometry HII region model computed by Baldwin et al. (1991). Size-resolved Orion grains are included. The constant pressure command does a hydrostatic equilibrium structure. The predicted emission-line spectrum is affected by the reddening of the internal grains. The resulting t2 analysis produces artificial results as a result. This has an open geometry, the original BFM paper was a closed geometry. (This makes little difference). Background cosmic rays are also included although these should have little effect on warm ionized gas. The emission line spectrum is given in surface brightness units, as in the BFM paper.

---

## **orion\_hii\_pdr *constant pressure H<sup>+</sup> region/pdr***

```

title constant pressure H+ region/pdr
c
c commands controlling continuum =====
c the incident continuum is two parts
c star and flux of photons striking it
c this is the photosphere of the OVI star, its temperature and Q(H)
c this is the result of this command
c table star kurucz 39600 K
c and makes it possible to run these orion sims without
c installing the stellar atmosphere files

```

```
table read "star_kurucz_39600.dat"
Q(H) 49
c this adds the observed hot brems, its temperature and the flux of
c photons striking the cloud
brems 6
phi(h) 10
c add cosmic rays, which are important in the pdr
cosmic rays, background
c
c commands controlling geometry =====
c this sets the separation from the star and the face of the cloud
radius 17.4507
c this turns off the stop temperature option
c the sim will not stop due to temperature
stop temperature off
c this is typical of a gmc, larson 1981
stop total hydrogen column density 22
c this sets the thickness of the HII region & PDR
stop thickness 0.5 linear parsec
c assume constant pressure
c constant gas pressure
constant pressure
c this will result in a milli gauss B-field in molecular region
magnetic field -5 gauss
c the orion environment full covers the star, so turn on sphere
sphere
c we have a spherical geometry but want to simulate observing
c through a spectrometer's slit. use the aperture
c command for this
aperture beam
c
c other commands for details =====
c mimic existence of unmodeled molecular gas
double
c iterate since lines optically thick
iterate
c the observed microturbulence, partially a flow, so not included in pressure
turbulence 8 km/sec no pressure
c set the line width so lines appear on the punch continuum
set punchLwidth 10 km/s
c there should be no failures at all
failures 3
c
c this model goes quite deep, needs lots of zones
set nend 2200
c
c commands for density & abundances =====
c this is the log of the initial H density, cm-3
hden 4
c this will speed up the calculation a bit
init file="ism.ini"
c this uses HII region abundances, but no grains
abundances hii region no grains
c this turns on orion grains, but leaves quantum heating off to save time
grains orion no qheat
c turn on PAHs, with an abundance that depends on H0 fraction,
c as suggested by long-slit observations of Orion bar,
c with an abundance 3x larger than default built into the code
grains pah function 3 no qheat
c
c commands controlling output =====
normalize to "O 1" 63.17m
c print lots of faint CO lines
print line faint -4
punch overview last "orion_hii_pdr.ovr"
```

```

punch heating "orion_hii_pdr.het"
punch cooling "orion_hii_pdr.col"
punch dr last "orion_hii_pdr.dr"
punch results last "orion_hii_pdr.rlt"
punch continuum last "orion_hii_pdr.con" units microns
punch hydrogen 21 cm last "orion_hii_pdr.21cm"
punch hydrogen lya last "orion_hii_pdr.lya"
punch element silicon last "orion_hii_pdr.sil"
punch grain extinction last "orion_hii_pdr.grnext"
punch grain charge last "orion_hii_pdr.grnchr"
punch grain potential last "orion_hii_pdr.grnpot"
punch grain temperature last "orion_hii_pdr.grntem"
c
c
c orion_hii_pdr.in
c class hii pdr
c =====

```

This extends BFM from the H<sup>+</sup> region into the PDR as in Abel et al 2005. This is the correct way to do a PDR calculation.

>>refer Orion model Baldwin, J., Ferland, G. J., >>refercon Martin, P. G., Corbin, M., Cota, S., Peterson, >>refercon B. M., & Slettebak, A. 1991, ApJ, 374, 580

>>refer physics HII/PDR Abel, N.P., Ferland, G.J., Shaw, G. & >>refercon van Hoof, P.A.M. 2005, ApJS, 161, 65

## **orion\_hii\_pdr\_fast *constant gas pressure H<sup>+</sup> region/PDR***

```

title constant gas pressure H+ region/PDR
c much faster due to physics disabled by fast.ini
c
c commands controlling continuum =====
c the incident continuum is two parts
c star and flux of photons striking it
c this is the result of this command
c table star kurucz 39600 K
c and makes it possible to run these orion sims without
c installing the stellar atmosphere files
table read "star_kurucz_39600.dat"
Q(H) 49
c add cosmic rays, which are important at depth
cosmic rays, background
c plus hot brems
brems 6
phi(h) 10
c
c commands for density & abundances =====
hden 4
init file="fast.ini"
abundances hii region no grains
c single sized grain with no quantum heating
grains orion no qheat single
c
c commands controlling geometry =====
radius 17.4507
sphere
c let it go into the molecular cloud
stop temperature 10 linear

```

```

c this is roughly the thickness of OMC1
stop thickness 1 linear parsec
c mimic existence of unmodeled molecular gas with double command
double
c
c other commands for details =====
failures 3
iterate
c make constant gas pressure since fast turned off line radiation pressure
constant gas pressure
c the observed microturbulence
turbulence 8 km/sec
c we have a spherical geometry but want to simulate observing
c through a spectrometer's slit. use the aperture
c command for this
aperture beam
c
c commands controlling output =====
normalize to "CaBo" 4861
print line faint -4
punch overview last "orion_hii_pdr_fast.ovr"
punch heating "orion_hii_pdr_fast.het"
punch cooling "orion_hii_pdr_fast.col"
punch dr last "orion_hii_pdr_fast.dr"
punch results last "orion_hii_pdr_fast.rlt"
punch continuum last "orion_hii_pdr_fast.con" units microns
punch ionizing continuum last "orion_hii_pdr_fast.ion"
punch hydrogen 21 cm last "orion_hii_pdr_fast.21cm"
punch hydrogen lya last "orion_hii_pdr_fast.lya"
punch element silicon last "orion_hii_pdr_fast.sil"
punch element sulphur last "orion_hii_pdr_fast.sul"
c
c
c orion_hii_pdr_fast.in
c class hii pdr
c =====
c

```

Orion HII region and PDR, simialr to orion\_hii\_pdr but much faster because of fast.ini

---

## **orion\_hii\_pdr\_pp *the Orion HII Region / PDR / Molecular cloud with an open geometry***

```

title the Orion HII Region / PDR / Molecular cloud with an open geometry
c
c commands controlling continuum =====
c the incident continuum has two parts
c kurucz continuum with flux of photons striking cloud
c this is the the OVI star, its temperature and phi(H)
c this is the result of this command
c table star kurucz 39600 K
c and makes it possible to run these orion sims without
c installing the stellar atmosphere files
table read "star_kurucz_39600.dat"
phi(H) 13
c
c this adds the observed hot brems
c its temperature (as log of T) and the flux of
c photons striking the cloud
brems 6

```



```
phi(h) 10
c
c cosmic rays are important for pdr chemistry
cosmic rays, background
c
c commands controlling geometry =====
c this turns off the stop temperature option
c so the sim will not stop due to temperature
stop temperature off
c this sets the thickness of the HII region & PDR
stop thickness 0.5 linear parsec
c this is typical of a gmc, larson 1981
stop total hydrogen column density 22
c this will result in a milli gauss B-field in molecular region
magnetic field -5 gauss
c assume constant pressure
constant pressure
set nend 2000
c
c other commands for details =====
failures 3
c mimic existence of unmodeled molecular gas
double
c iterate since lines optically thick
iterate
c set microturbulence in equipartition with B field
turbulence equipartition
c set the line width so lines appear on the punch continuum
set punchLwidth 10 km/s
c
c commands for density & abundances =====
c this is the log of the initial H density, cm-3
hden 4
c this will speed up the calculation a bit
init file="ism.ini"
c this uses HII region abundances, but no grains
abundances hii region no grains
c this uses orion grains
grains orion
c turn on PAHs, with an abundance that depends on H0 fraction,
c as suggested by long-slit observations of Orion bar,
c with an abundance 3x larger than default built into the code
grains pah function 3
c
c commands controlling output =====
c print lots of faint CO lines
print line faint -6
c normalize to Ha
normalize to "H 1" 6563
punch overview last "orion_hii_pdr_pp.ovr"
punch temperature last "orion_hii_pdr_pp.tem"
punch overview "orion_hii_pdr_pp.ovr1"
punch molecules last "orion_hii_pdr_pp.mol"
punch molecules "orion_hii_pdr_pp.mol1"
punch heating "orion_hii_pdr_pp.het"
punch cooling "orion_hii_pdr_pp.col"
punch dr last "orion_hii_pdr_pp.dr"
punch results last "orion_hii_pdr_pp.rlt"
punch continuum last "orion_hii_pdr_pp.con" units microns
punch hydrogen 21 cm last "orion_hii_pdr_pp.21cm"
punch hydrogen lya last "orion_hii_pdr_pp.lya"
punch element silicon last "orion_hii_pdr_pp.sil"
punch pressure last "orion_hii_pdr_pp.pre"
punch pressure "orion_hii_pdr_pp.prel"
punch grain extinction last "orion_hii_pdr_pp.grnext"
```

```

punch grain charge last "orion_hii_pdr_pp.grnchr"
punch grain potential last "orion_hii_pdr_pp.grnpot"
punch grain temperature last "orion_hii_pdr_pp.grntem"
punch grain temperature "orion_hii_pdr_pp.grntem1"
c
c
c orion_hii_pdr_pp.in
c class hii pdr
c =====

```

similar to orion\_hii\_pdr except for plane parallel geometry

---

## **pdr\_co\_fully *H2 and CO in fully molecular limit***

```

title H2 and CO in fully molecular limit
c this is a pair with pdr_co_fully_noneq - that tests non equilibrium chem
c
c commands controlling continuum =====
blackbody 5000
luminosity total solar linear 2
brems 6
luminosity total solar log -2.7
c
c commands for density & abundances =====
c set the density and composition, ism with grains
hden 10
abundances ism no qheat
c
c other commands for details =====
c this is an important ionization source
cosmic ray background
c code will stop at 4000 K by default, lower this temp so we
c can do fully molecular gas
stop temperature 10K linear
c
c commands controlling geometry =====
c the radius and thickness
radius 15.8
stop thickness 6
turbulence 3 km / s
c
c commands controlling output =====
punch overview "pdr_co_fully.ovr"
punch molecules "pdr_co_fully.mol"
punch dr "pdr_co_fully.dr"
punch heating "pdr_co_fully.het"
punch cooling "pdr_co_fully.col"
punch grain charge "pdr_co_fully.grnchr"
punch grain drift "pdr_co_fully.grndft"
punch grain temperature "pdr_co_fully.grntem"
c
c
c gas has more he++ than He+ due to fast charge transfer with co
c neutralizing he+
c pdr_co_fully.in
c class pdr
c =====

```

test code in fully molecular limit this is a pair with pdr\_co\_fully\_noneq - that tests non equilibrium chem

---

## **pdr\_co\_fully\_noneq *H2 and CO are fully molecular, non-equilibrium case***

```

title H2 and CO are fully molecular, non-equilibrium case
c this is a pair with pdr_co_fully.in, which does not include non-equil chem
c
c commands controlling continuum =====
blackbody 5000
luminosity total solar linear 2
brems 6
luminosity total solar log -2.7
c
c commands for density & abundances =====
c set the density and composition, ism with grains
hden 10
abundances ism no qheat
c
c other commands for details =====
c this is an important ionization source
cosmic ray background
c code will stop at 4000 K by default, lower this temp so we
c can do fully molecular gas
stop temperature 10K linear
c
c commands controlling geometry =====
c the radius and thickness
radius 15.8
stop thickness 6
turbulence 3 km / s
c this turns on non-equilibrium chemistry as suggested by Federman et al. 1996
set chemistry non equilibrium
c
c commands controlling output =====
punch overview "pdr_co_fully_noneq.ovr"
punch molecules "pdr_co_fully_noneq.mol"
punch dr "pdr_co_fully_noneq.dr"
punch heating "pdr_co_fully_noneq.het"
punch cooling "pdr_co_fully_noneq.col"
punch grain charge "pdr_co_fully_noneq.grnchr"
punch grain drift "pdr_co_fully_noneq.grndft"
punch grain temperature "pdr_co_fully_noneq.grntem"
c
c
c gas has more he++ than He+ due to fast charge transfer with co
c neutralizing he+
c pdr_co_fully_noneq.in
c class pdr
c =====

```

test code in fully molecular limit with Federman non-equilibrium chem this is a pair with pdr\_co\_fully.in, which does not include non-equil chem

---

## **pdr\_coolbb *illumination by cool STE blackbody***

```

title illumination by cool STE blackbody
c gasis fully molecular, grains should be in STE
c

```

```

c commands controlling continuum =====
c illuminate with 120 blackbody in thermodynamic equilibrium
blackbody 120 STE
CMB
c
c commands for density & abundances =====
c density
hden 7
c ism dust and grains
abundances ISM no grains
c dust to gas ratio is 10x standard
grains 1.0
c lowers metals and grains to 1/10 standard, so metals are
c 1/10 solar while dust to gas ratio is galactic
metals and grains 0.1
c
c other commands for details =====
c set background cosmic rays to allow chemistry
cosmic ray background
iterate
c
c commands controlling geometry =====
radius 17.5
sphere
stop temperature 20K linear
stop column density 25.5
c
c commands controlling output =====
punch overview "pdr_coolbb.ovr"
punch molecules "pdr_coolbb.mol"
punch heating "pdr_coolbb.het"
punch cooling "pdr_coolbb.col"
punch grain temperature "pdr_coolbb.grntem"
c
c
c pdr_coolbb.in
c class pdr
c =====

```

This is a dense ISM exposed to a cool blackbody. This tests the ability to converge onto a VERY molecular environment. Nearly all C and H are in CO and H<sub>2</sub>. Cosmic rays are the main source of heat and ionization.

---

## **pdr\_dense\_persei *dense phase zeta persei cloud***

```

title dense phase zeta persei cloud
c
c density and abundances =====
c density of dense phase from Table 2 of Le Petit paper =====
hden 4.3
c
c abundances from Table 1 of Le Petit paper =====
element carbon abundance 0.000132 linear
element helium abundance 0.10 linear
element oxygen abundance 0.00032 linear
element nitrogen abundance 0.000075 linear
element sulphur abundance 0.0000186 linear
element silicon abundance 0.000029 linear
c set abundance of all other elements to zero =====
element copper off
element magnesium off

```

```

element manganese off
element sodium off
element chlorine off
element vanadium off
element potassium off
element phosphorous off
element calcium off
element iron off
element zinc off
element neon off
element argon off
element fluorine off
element aluminum off
element boron off
element lithium off
element beryllium off
element scandium off
element nickel off
element titanium off
element chromium off
element cobalt off
c do not use Federman rates for this model =====
set federman chemistry off
c use standard ism grain size distribution =====
grains ism
c
c command controlling the continuum, for this model is Draine 1978 field ====
table draine 0.5 linear
c make sure no H-ionizing radiation strikes the cloud
extinguish 24
c
c fix the temperature to 20 K =====
constant temperature 20
c stop at a radius of 4.3e-4 parsecs =====
stop thickness 0.00043 parsecs linear
c Le Petit model does not consider ices, so turn this off =====
no grain molecules
c turn on cosmic rays =====
cosmic rays background
c Set cosmic ray ionization rate to Table 2 of Le Petit paper =====
set csupra -15.6
c Allow calculation to go extend into cold environment =====
stop temperature linear 3
c
c commands controlling output =====
punch overview "pdr_dense_persei.ovr"
punch dr "pdr_dense_persei.dr"
punch molecules "pdr_dense_persei.mol"
punch heating "pdr_dense_persei.het"
punch cooling "pdr_dense_persei.col"
c
c pdr_dense_persei.in
c class pdr
c =====

```

This is the dense phase model presented by LePetit, Roueff, and Herbst in order to reproduce C2 and C3 column densities observed along the line of sight to zera persei. This is our attempt at reproducing their calculation. This is the dense molecular phase, not the phase that produces H3+

// >>refer model pdr Le Petit, F., Roueff, E., & Herbst, E. 2004, // >>refercon A&A, 417, 993

If you do a thermal equilibrium calculation by removing the constant temperature command the kinetic temperature will be about three times larger than assumed in their paper.

## **pdr\_HTT91 *low-density PDR from HTT91***

```
title low-density PDR from HTT91
c >>refer test model Hollenbach, D., Takahashi, T., &
c >>refcon Tielens, A.G.G.M., 1991, ApJ,377, 192-209
c
c commands controlling continuum =====
c first continuum is FIR hot grain continuum produced in
c unmodeled HII Region
c this is hot star continuum
black 30000
intensity 0.2 range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
cosmic rays, background
c
c commands for density & abundances =====
hden 3
init file="ism.ini"
abundances he -1.01 c -3.52 n-8 o-3.30 ne-8 mg-5.89
continue si -6.10 s -5.10 cl-7 ar-8 fe -6.60
grains orion, abundance log 0.16 no qheat
c
c commands controlling geometry =====
sphere
c stop when gas is fully neutral
stop efrac -10
c this is to stop at an intended place, since results would be
c very dependent on details if we stop on temperature
stop thickness 18.954
c stop when gas is cold
stop temperature 10 linear
c add this to mimic unmodelled neutral gas
double optical depths
c
c other commands for details =====
iterate 2
case b
turbulence 1.5 km/sec
c
c commands controlling output =====
normalize to 157.6m "C 2"
c uv lines are strongly pumped by stellar continuum, break out contribution
print line pump
print line optical depths
print ages
punch overview last "pdr_HTT91.ovr"
punch dr "pdr_HTT91.dr"
punch molecules last "pdr_HTT91.mol"
punch results last "pdr_HTT91.rlt"
punch continuum units microns last "pdr_HTT91.con"
punch heating last "pdr_HTT91.het"
punch cooling last "pdr_HTT91.col"
c
c pdr_HTT91.in
c class pdr
c =====
c
```

This is the Hollenbach et al 1991 Low-density PDR The case b command appears because the Lyman lines are vastly optically thick in this environment. If the command is removed the Lyman lines will be optically

thin, which will result in fast fluorescent excitation of the (nearly totally neutral) hydrogen atom.

there is very little CO in this model since it is not thick enough for the UV pumping lines to become optically thick

## **pdr\_leiden\_f1** *Leiden meeting model 1*

```

title Leiden meeting model 1
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 5
extinguish 24
c
c commands for density & abundances =====
c hydrogen density
hden 3.
grains ism 1.16 no qheat
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c
c commands controlling output =====
print line optical depths
punch overview "pdr_leiden_f1.ovr"
punch continuum "pdr_leiden_f1.con"
punch leiden lines "pdr_leiden_f1.lin"
punch leiden "pdr_leiden_f1.lei"
punch dr "pdr_leiden_f1.dr"
punch molecules "pdr_leiden_f1.mol"
punch grain physical conditions "pdr_leiden_f1.grn"
punch chemistry rates CO "pdr_leiden_f1.rat"
punch H2 destruction "pdr_leiden_f1.H2d"
punch H2 creation "pdr_leiden_f1.H2c"
punch secondaries "pdr_leiden_f1.sec"
c
c
c pdr_leiden_f1.in
c class pdr
c =====
c

```

This sim has some interesting properties. The grain temp is so low that O freezes onto grains as H<sub>2</sub>O. This is so efficient that CO does not become well formed - the CO/C ratio only reaches 0.3. This is one of the tests in Rollig et al. 2007, A&A, 467, 187

## **pdr\_leiden\_f2 *Leiden meeting model 2***

```

title Leiden meeting model 2
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use
c half the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c commands for density & abundances =====
grains ism 1.16 no qheat
c hydrogen density
hden 3.
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c
c commands controlling output =====
punch overview "pdr_leiden_f2.ovr"
punch continuum "pdr_leiden_f2.con"
punch molecules "pdr_leiden_f2.mol"
punch leiden lines "pdr_leiden_f2.lin"
punch leiden "pdr_leiden_f2.lei"
punch dr "pdr_leiden_f2.dr"
punch grain physical conditions "pdr_leiden_f2.grn"
c
c
c pdr_leiden_f2.in
c class pdr
c =====
c

c This is one of the tests in Rollig et al. 2007, A&A, 467, 187

```

---

## **pdr\_leiden\_f3 *Leiden meeting model 3***

```

title Leiden meeting model 3
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 5
extinguish 24
c
c commands for density & abundances =====

```



```

grains ism 1.16 no wheat
c hydrogen density
hden 5.5
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c
c commands controlling output =====
punch overview "pdr_leiden_f3.ovr"
punch continuum "pdr_leiden_f3.con"
punch leiden lines "pdr_leiden_f3.lin"
punch leiden "pdr_leiden_f3.lei"
punch dr "pdr_leiden_f3.dr"
punch grain physical conditions "pdr_leiden_f3.grn"
c
c
c pdr_leiden_f3.in
c class pdr
c =====
c

c This is one of the tests in Rollig et al. 2007, A&A, 467, 187

```

---

## **pdr\_leiden\_f4** *Leiden meeting model 4*

```

title Leiden meeting model 4
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c
c commands for density & abundances =====
grains ism 1.16 no wheat
c
c commands controlling geometry =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c hydrogen density
hden 5.5
c
c commands controlling output =====
punch overview "pdr_leiden_f4.ovr"

```

```

punch continuum "pdr_leiden_f4.con"
punch leiden lines "pdr_leiden_f4.lin"
punch leiden "pdr_leiden_f4.lei"
punch dr "pdr_leiden_f4.dr"
punch grain physical conditions "pdr_leiden_f4.grn"
c
c
c pdr_leiden_f4.in
c class pdr
c =====
c

```

c This is one of the tests in Rollig et al. 2007, A&A, 467, 187

---

## **pdr\_leiden\_hack\_f1 *Leiden meeting model 1 with hacks***

```

title Leiden meeting model 1 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 5
extinguish 24
c
c commands for density & abundances =====
c hydrogen density
hden 3.
grains ism 1.16 no qheat
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_f1.lin"
punch leiden "pdr_leiden_hack_f1.lei"
punch dr "pdr_leiden_hack_f1.dr"
punch grain physical conditions "pdr_leiden_hack_f1.grn"
punch overview "pdr_leiden_hack_f1.ovr"
c
c
c
c
c pdr_leiden_f1_hack.in
c class pdr
c =====
c

```

c This is one of the tests in Rollig et al. 2007, A&A, 467, 187 this includes hacks to get "standard answer"

---

## **pdr\_leiden\_hack\_f2** *Leiden meeting model 2 with hacks*

```

title Leiden meeting model 2 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use
c half the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c commands for density & abundances =====
grains ism 1.16 no qheat
c hydrogen density
hden 3.
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_f2.lin"
punch leiden "pdr_leiden_hack_f2.lei"
punch dr "pdr_leiden_hack_f2.dr"
punch grain physical conditions "pdr_leiden_hack_f2.grn"
punch overview "pdr_leiden_hack_f2.ovr"
c
c
c pdr_leiden_hack_f2.in
c class pdr
c =====
c

c This is one of the tests in Rollig et al. 2007, A&A, 467, 187 includes hacks to get standard answer.

```

---

## **pdr\_leiden\_hack\_f3** *Leiden meeting model 3 with hacks*

```

title Leiden meeting model 3 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 5
extinguish 24
c
c commands for density & abundances =====
grains ism 1.16 no qheat
c hydrogen density

```

```

hden 5.5
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_f3.lin"
punch leiden overview "pdr_leiden_hack_f3.ovr"
punch leiden "pdr_leiden_hack_f3.lei"
punch dr "pdr_leiden_hack_f3.dr"
punch grain physical conditions "pdr_leiden_hack_f3.grn"
c
c
c pdr_leiden_hack_f3.in
c class pdr
c =====
c

```

c This is one of the tests in Rollig et al. 2007, A&A, 467, 187 include hacks to get standard answer.

---

## **pdr\_leiden\_hack\_f4 *Leiden meeting model 4 with hacks***

```

title Leiden meeting model 4 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c
c commands for density & abundances =====
grains ism 1.16 no qheat
c
c commands controlling geometry =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
c This command defines the grain temperature to be a constant 20 Kelvin
constant grain temperature 20
c This sets the temperature to a constant 50 Kelvin
constant temperature 50 linear
c hydrogen density
hden 5.5
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_f4.lin"
punch leiden "pdr_leiden_hack_f4.lei"
punch dr "pdr_leiden_hack_f4.dr"
punch grain physical conditions "pdr_leiden_hack_f4.grn"

```

```
punch overview "pdr_leiden_hack_f4.ovr"
c
c
c pdr_leiden_hack_f4.in
c class pdr
c =====
c
```

This is one of the tests in Rollig et al. 2007, A&A, 467, 187 includes hacks needed to get standard answer.

---

## **pdr\_leiden\_hack\_v1** *Leiden meeting model 5 with hacks*

```
title Leiden meeting model 5 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c is half the requested value
c they want for the model is actually twice the value below
table draine 5
extinguish 24
c
c commands for density & abundances =====
c hydrogen density
hden 3.
grains ism 1.16 no qheat
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
* element nitrogen on
* element nitrogen scale -1
* element silicon on
* element silicon scale -1
* element sulphur on
* element sulphur scale -1
* element chlorine on
* element chlorine scale -1
c Define the abundance of PAHs so we can reproduce the same photoelectric
c heating curve as BT94
c We have also scaled the grain abundance to achieve the desired AV/NH ratio
* grains PAH no qheat 8
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_v1.lin"
punch leiden "pdr_leiden_hack_v1.lei"
punch dr "pdr_leiden_hack_v1.dr"
punch grain temperature "pdr_leiden_hack_v1.grn"
punch overview "pdr_leiden_hack_v1.ovr"
c
c
c pdr_leiden_hack_v1.in
c class pdr
c =====
c
```

This is one of the tests in Rollig et al. 2007, A&A, 467, 187 includes hacks to get standard answer.

---

### **pdr\_leiden\_hack\_v2** *Leiden meeting model 6 with hacks*

```

title Leiden meeting model 6 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c commands for density & abundances =====
c Define the abundance of PAHs so we can reproduce the same photoelectric
c heating curve as BT94
c We have also scaled the grain abundance to achieve the desired AV/NH ratio
* grains PAH no qheat 40
grains ism 1.16 no qheat
c hydrogen density
hden 3.
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_v2.lin"
punch leiden "pdr_leiden_hack_v2.lei"
punch dr "pdr_leiden_hack_v2.dr"
punch grain temperature "pdr_leiden_hack_v2.grn"
punch overview "pdr_leiden_hack_v2.ovr"
punch heating "pdr_leiden_hack_v2.het"
punch cooling "pdr_leiden_hack_v2.col"
punch h2 rates "pdr_leiden_hack_v2.h2rat"
c
c
c pdr_leiden_hack_v2.in
c class pdr
c =====
c

```

This is one of the tests in Rollig et al. 2007, A&A, 467, 187 includes hacks to get standard answer.

---

### **pdr\_leiden\_hack\_v3** *Leiden meeting model 7 with hacks*

```

title Leiden meeting model 7 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below

```

```

table draine 5
extinguish 24
c
c commands for density & abundances =====
c Define the abundance of PAHs so we can reproduce the same photoelectric
c heating curve as BT94
c We have also scaled the grain abundance to achieve the desired AV/NH ratio
* grains PAH no qheat 5.7
grains ism 1.16 no qheat
c hydrogen density
hden 5.5
c
c commands controlling geometry =====
c
c other commands for details =====
c this sim has more than one thermal solution - force into lowest one,
c the one found by the rest of the group - without this we will find a
c soln at 1400 K
force temperature to 60K
failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_v3.lin"
punch leiden "pdr_leiden_hack_v3.lei"
punch dr "pdr_leiden_hack_v3.dr"
punch grain temperature "pdr_leiden_hack_v3.grn"
punch overview "pdr_leiden_hack_v3.ovr"
punch heating "pdr_leiden_hack_v3.het"
punch cooling "pdr_leiden_hack_v3.col"
c
c
c pdr_leiden_hack_v3.in
c class pdr
c =====
c

```

This is one of the tests in Rollig et al. 2007, A&A, 467, 187 includes hacks to get standard answer

---

## **pdr\_leiden\_hack\_v4** *Leiden meeting model 8 with hacks*

```

title Leiden meeting model 8 with hacks
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c commands for density & abundances =====
c Define the abundance of PAHs so we can reproduce the same photoelectric
c heating curve as BT94
c We have also scaled the grain abundance to achieve the desired AV/NH ratio
* grains PAH no qheat 40
* grains ism 0.936 no qheat
grains ism 1.16 no qheat
set PAH Bakes
c hydrogen density

```

```

hden 5.5
c
c commands controlling geometry =====
c
c other commands for details =====
* failures 3
c use leiden initialization file
init file="pdr_leiden_hack.ini"
c
c commands controlling output =====
punch leiden lines "pdr_leiden_hack_v4.lin"
punch leiden "pdr_leiden_hack_v4.lei"
punch dr "pdr_leiden_hack_v4.dr"
punch grain temperature "pdr_leiden_hack_v4.grn"
punch overview "pdr_leiden_hack_v4.ovr"
punch heating "pdr_leiden_hack_v4.het"
punch cooling "pdr_leiden_hack_v4.col"
punch transmitted continuum "pdr_leiden_hack_v4.con"
c
c
c pdr_leiden_hack_v4.in
c class pdr
c =====
c

```

This is one of the tests in Rollig et al. 2007, A&A, 467, 187 includes hacks to get standard answer.

---

## **pdr\_leiden\_v1 *Leiden meeting model 5***

```

title Leiden meeting model 5
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c is half the requested value
c they want for the model is actually twice the value below
table draine 5
extinguish 24
c
c commands for density & abundances =====
c hydrogen density
hden 3.
grains ism 1.16 no qheat
c add PAHs
grains PAH no qheat 3 function
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c
c commands controlling output =====
punch overview "pdr_leiden_v1.ovr"
punch leiden lines "pdr_leiden_v1.lin"
punch leiden "pdr_leiden_v1.lei"
punch dr "pdr_leiden_v1.dr"
punch grain temperature "pdr_leiden_v1.grn"
c

```



```

c
c
c pdr_leiden_v1.in
c class pdr
c =====
c

```

This is one of the tests in Rollig et al. 2007, A&A, 467, 187

---

## **pdr\_leiden\_v2 *Leiden meeting model 6***

```

title Leiden meeting model 6
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c commands for density & abundances =====
c add PAHs and grains
grains PAH no qheat 3 function
grains ism 1.16 no qheat
c hydrogen density
hden 3.
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c
c commands controlling output =====
punch overview "pdr_leiden_v2.ovr"
punch leiden lines "pdr_leiden_v2.lin"
punch leiden "pdr_leiden_v2.lei"
punch dr "pdr_leiden_v2.dr"
punch grain temperature "pdr_leiden_v2.grn"
c
c
c pdr_leiden_v2.in
c class pdr
c =====
c

```

c This is one of the tests in Rollig et al. 2007, A&A, 467, 187

---

## **pdr\_leiden\_v3 *Leiden meeting model 7***

```

title Leiden meeting model 7
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====

```

```

c Use the Draine 1978 field, for a semi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 5
extinguish 24
c
c commands for density & abundances =====
c add PAHs and grains
grains PAH no qheat 3 function
grains ism 1.16 no qheat
c hydrogen density
hden 5.5
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c
c commands controlling output =====
punch overview "pdr_leiden_v3.ovr"
punch leiden lines "pdr_leiden_v3.lin"
punch leiden "pdr_leiden_v3.lei"
punch dr "pdr_leiden_v3.dr"
punch grain temperature "pdr_leiden_v3.grn"
punch heating "pdr_leiden_v3.het"
punch cooling "pdr_leiden_v3.col"
c
c
c pdr_leiden_v3.in
c class pdr
c =====
c

```

---

## **pdr\_leiden\_v4** *Leiden meeting model 8*

```

title Leiden meeting model 8
c This is one of the tests in Rollig et al. 2007, A&A, 467, 187
c
c commands controlling continuum =====
c Use the Draine 1978 field, for asemi-infinite slab we have to use half
c the requested value, so the actual value
c they want for the model is actually twice the value below
table draine 50000
extinguish 24
c
c commands for density & abundances =====
grains PAH no qheat 3 function
grains ism 1.16 no qheat
c hydrogen density
hden 5.5
c
c commands controlling geometry =====
c
c other commands for details =====
failures 3
c use leiden initialization file
init file="pdr_leiden.ini"
c
c commands controlling output =====

```

```

punch overview "pdr_leiden_v4.ovr"
punch leiden lines "pdr_leiden_v4.lin"
punch leiden "pdr_leiden_v4.lei"
punch dr "pdr_leiden_v4.dr"
punch grain temperature "pdr_leiden_v4.grn"
punch transmitted continuum "pdr_leiden_v4.con"
c
c
c pdr_leiden_v4.in
c class pdr
c =====
c

```

---

## **pdr\_orion\_veil** *Orion's veil*

```

title Orion's veil
c
c commands controlling continuum =====
q(h) 50.000000
table star rubin
c AGN & CMB background
background z=0
cosmic rays background
c as a test this was tried - atomic temperature raised fm 100 to 140
c but nothing really wrong
* cosmic rays equipartition
c the observed magnetic field, roughly 100 micro Gauss
magnetic field -4
c
c commands for density & abundances =====
hden 3.1
init "ism.ini"
abundances orion no grains
grains orion no qheat
c
c commands controlling geometry =====
stop neutral hydrogen column density 21.6
radius 19.1
stop temperature linear 5
c this will not be hit, but put it in to insure that we stop if
c gas becomes totally molecular
stop av 10
c
c other commands for details =====
failures 3
* uncomment to include large H2 molecule - results will CHANGE!
* atom h2
turbulence 0.5 km/sec
sphere
c c this simulates Lyman lines in the stellar continuum
c atom h-like lyman pumping off
c
c commands controlling output =====
print line optical depths
punch PDR "pdr_orion_veil.pdr"
punch feii column densities "pdr_orion_veil.fe2col"
punch hydrogen 21 cm "pdr_orion_veil.21cm"
punch molecules "pdr_orion_veil.mol"
punch overveiw "pdr_orion_veil.ovr"
punch column densities "pdr_orion_veil.col"
punch heating "pdr_orion_veil.het"

```

```

punch cooling "pdr_orion_veil.col"
punch dr "pdr_orion_veil.dr"
punch H2 rates "pdr_orion_veil.h2rat"
punch pressure "pdr_orion_veil.pre"
punch wind "pdr_orion_veil.wnd"
print line optical depths
punch continuum units microns "pdr_orion_veil.con" last
c
c pdr_orion_veil.in
c class pdr
c =====
c

```

This is Abel et al 2004 model of Orion's veil, the layer of gas in front of the Orion Nebula. >>refer Orion Veil  
Abel, N. P., Brogan, C. L., Ferland, G. J., >>refercon O'Dell, C. R., Shaw, G., & Troland, T. H. 2004

---

## **pdr\_th85ism** *TH85 PDR with ISM grains*

```

title TH85 PDR with ISM grains
c
c commands controlling continuum =====
c cosmic radiation background
background z=0
c cosmic ray background ionization and heating
cosmic rays, background
c first continuum is FIR hot grain continuum produced in
c unmodeled HII Region
blackbody, t = 75 K
intensity 2.7 (total)
c this is the attenuated hot stellar continuum
black 30000
intensity 2.204 range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c
c commands controlling geometry =====
c simulate effects of gas we do not model
double
c their turbulence
turbulence 2.7 km/sec
c stopping criteria
c this should be the one actually used
stop AV 33.2 extended
c stop when gas is fully neutral
stop efrac -10
c stop when gas is cold
stop temperature 10 linear
c stop at thickness of 16.5 so that mole limit does not
c stop this calculation - that would make results very
c detail dependent
stop thickness 19
c
c commands for density & abundances =====
hden 5.362
c this turns off some elements, and processes we don't need
c so needs to come before the elements
init file="ism.ini"
abundances he -1.01 c -3.52 n-8 o-3.30 ne-8 mg-5.89
continue si -6.10 s -5.10 cl-7 ar-8 fe -6.60
c use ism abundances but turn of quantum heating

```

```

grains ism, abundance log 0.16 no qheat
c
c other commands for details =====
c must iterate since many FIR MIR lines are optically thick
iterate
c this is a closed geometry, in Orion, veil covers 2pi sr
sphere
c make Lyman lines very optically thick, which stops pumping by
c Balmer continuum
case b
c this should run cleanly - turn down number of allow conv fails
failures 3
c
c commands controlling output =====
normalize to "C 2" 157.6m
c want the spectrum to be relative to this [C II] line
c uv lines are strongly pumped by stellar continuum, break out contribution
print line pump
print line optical depths
print ages
punch overview last "pdr_th85ism.ovr"
punch pdr last "pdr_th85ism.pdr"
punch molecules last "pdr_th85ism.mol"
punch dr last "pdr_th85ism.dr"
punch hydrogen physical conditions last "pdr_th85ism.hyd"
punch results last "pdr_th85ism.rlt"
punch continuum units microns last "pdr_th85ism.con"
punch heating last "pdr_th85ism.het"
punch cooling last "pdr_th85ism.col"
punch hydrogen 21 cm last "pdr_th85ism.21cm"
punch element magnesium last "pdr_th85ism.mag"
punch grain extinction last "pdr_th85ism.grnext"
punch grain potential last "pdr_th85ism.grnpot"
punch grain temperature last "pdr_th85ism.grntem"
punch grain charge last "pdr_th85ism.grnchr"
c
table lines "LineList_PDR.dat"
c pdr_th85ism.in
c class pdr
c =====
c

```

This is the Tielens and Hollenbach (1985a, b) standard model of the Orion photodissociation region (PDR). Specifically, this is my attempt at their Table 2 of paper 2, ApJ 291, p749. The case b command appears because the Lyman lines are vastly optically thick in this environment. If the command is removed the Lyman lines will be optically thin, which will result in fast fluorescent excitation of the (nearly totally neutral) hydrogen atom.

// >>refer model pdr Tielens, A. G. G. M., & Hollenbach, D. 1985a, ApJ, 291, 722 // >>refer model pdr Tielens, A. G. G. M., & Hollenbach, D. 1985b, ApJ, 291, 746

---

## **pdr\_th85ism\_cgto** *TH85 pdr with ISM grains and C > O*

```

title TH85 pdr with ISM grains and C > O
c
c commands controlling continuum =====
c cosmic radiation background
background z=0
c cosmic ray background ionization and heating

```

```
cosmic rays, background
c first continuum is FIR hot grain continuum produced in
c unmodeled HII Region
blackbody, t = 75 K
intensity 2.7 (total)
c this is the attenuated hot stellar continuum
black 30000
intensity 2.204 range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c
c commands controlling geometry =====
c simulate effects of gas we do not model
double
c their turbulence
turbulence 2.7 km/sec
c stopping criteria
c this should be the one actually used
stop AV 33.2 extended
c stop when gas is fully neutral
stop efrac -10
c stop when gas is cold
stop temperature 10 linear
c stop at thickness of 16.5 so that mole limit does not
c stop this calculation - that would make results very
c detail dependent
stop thickness 19
c
c commands for density & abundances =====
hden 5.362
c this turns off some elements, and processes we don't need
c so needs to come before the elements
init file="ism.ini"
c this is the TH85 mix except C is raised by 0.2 dex,
c and O lowered by his amount, so that C/O > 1
abundances he -1.01 c -3.4 n-8 o-3.50 ne-8 mg-5.89
continue si -6.10 s -5.10 cl-7 ar-8 fe -6.60
c use ism abundances but turn of quantum heating
grains ism, abundance log 0.16 no qheat
c
c other commands for details =====
c must iterate since many FIR MIR lines are optically thick
iterate
c this is a closed geometry, in Orion, veil covers 2pi sr
sphere
c make Lyman lines very optically thick, which stops pumping by
c Balmer continuum
case b
c this should run cleanly - turn down number of allow conv fails
failures 3
c
c commands controlling output =====
normalize to "C 2" 157.6m
c want the spectrum to be relative to this [C II] line
c uv lines are strongly pumped by stellar continuum, break out contribution
print line pump
print line optical depths
print ages
punch overview last "pdr_th85ism_cgto.ovr"
punch pdr last "pdr_th85ism_cgto.pdr"
punch molecules last "pdr_th85ism_cgto.mol"
punch dr last "pdr_th85ism_cgto.dr"
punch hydrogen physical conditions last "pdr_th85ism_cgto.hyd"
punch results last "pdr_th85ism_cgto.rlt"
punch continuum units microns last "pdr_th85ism_cgto.con"
```

```

punch heating last "pdr_th85ism_cgto.het"
punch cooling last "pdr_th85ism_cgto.col"
punch hydrogen 21 cm last "pdr_th85ism_cgto.21cm"
punch element magnesium last "pdr_th85ism_cgto.mag"
punch grain extinction last "pdr_th85ism_cgto.grnext"
punch grain potential last "pdr_th85ism_cgto.grnpot"
punch grain temperature last "pdr_th85ism_cgto.grntem"
punch grain charge last "pdr_th85ism_cgto.grnchr"
c
c pdr_th85ism_cgto.in
c class pdr
c =====
c

```

This is the Tielens and Hollenbach (1985a, b) standard model of the Orion photodissociation region (PDR) but with the C abundance increased by 0.2 dex, and O lowered by this amount, so that C/O>1. Specifically, this is their Table 2 of paper 2, ApJ 291, p749. The case b command appears because the Lyman lines are vastly optically thick in this environment. If the command is removed the Lyman lines will be optically thin, which will result in fast fluorescent excitation of the (nearly totally neutral) hydrogen atom.

```

// >>refer model pdr Tielens, A. G. G. M., & Hollenbach, D. 1985a, ApJ, 291, 722 // >>refer model pdr
Tielens, A. G. G. M., & Hollenbach, D. 1985b, ApJ, 291, 746

```

---

## **pdr\_th85orion** *TH85 PDR with Orion grains*

```

title TH85 PDR with Orion grains
c
c commands controlling continuum =====
c cosmic background
background
c galactic cosmic ray background
cosmic rays, background
c first continuum is FIR hot grain continuum produced in
c unmodeled HII Region
blackbody, t = 75 K
intensity 2.7 (total)
c this is the attenuated hot stellar continuum
black 30000
intensity 2.204 range 0.4412 to 1 Ryd
c this will remove all ionizing radiation
extinguish 24 0
c
c commands for density & abundances =====
hden 5.362
init file="ism.ini"
grains orion, abundance log 0.16 no qheat
abundances he -1.01 c -3.52 n-8 o-3.30 ne-8 mg-5.89
continue si -6.10 s -5.10 cl-7 ar-8 fe -6.60
c
c commands controlling geometry =====
c simulate effects of gas we do not model
double
sphere
c stopping criteria
c this should be the one actually used
c >>chn 04 mar 13, from 10 to 36.3, bug in AV had stopped at this depth,
c so change it to this to keep model the same
stop AV 36.3 extended
c stop when gas is fully neutral

```

```

stop efrac -10
c stop when gas is cold
stop temperature 10 linear
c stop at thickness so that would make results very
c detail dependent
stop thickness 19
c
c other commands for details =====
turbulence 2.7 km/sec
iterate
failures 3
c
c this is done to not allow pumping and subsequent photoionization
c of H by Balmer continuum in keeping with std pdr assumptions
case b
c
c commands controlling output =====
normalize to "C 2" 157.6m
c uv lines are strongly pumped by stellar continuum, break out contribution
print line pump
print line optical depths
print ages
c
punch overview "pdr_th85orion.ovr" no hash
punch pdr last "pdr_th85orion.pdr"
c this only outputs major molecules
punch molecules last "pdr_th85orion.mol"
punch dr "pdr_th85orion.dr"
c this is to get all of H molecules
punch hydrogen physical conditions last "pdr_th85orion.hyd"
punch results last "pdr_th85orion.rlt"
punch continuum units microns last "pdr_th85orion.con"
punch ionizing continuum last "pdr_th85orion.ion"
punch heating last "pdr_th85orion.het"
punch cooling last "pdr_th85orion.col"
punch hydrogen 21 cm last "pdr_th85orion.21cm"
c these are for the grains
punch grain extinction last "pdr_th85orion.grnext"
punch grain potential last "pdr_th85orion.grnpot"
punch grain temperature last "pdr_th85orion.grntem"
punch grain charge last "pdr_th85orion.grnchr"
c
c
c pdr_th85orion.in
c class pdr
c =====

```

This is the Tielens and Hollenbach (1985a, b) standard model of the Orion photodissociation region (PDR). Specifically, this is my attempt at their Table 2 of paper 2, ApJ 291, p749. The case b command appears because the Lyman lines are vastly optically thick in this environment. If the command is removed the Lyman lines will be optically thin, which will result in fast fluorescent excitation of the (nearly totally neutral) hydrogen atom.

```
// >>refer model pdr Tielens, A. G. G. M., & Hollenbach, D. 1985a, ApJ, 291, 722 // >>refer model pdr
Tielens, A. G. G. M., & Hollenbach, D. 1985b, ApJ, 291, 746
```

---

## **pn\_fluc *Paris PN with density fluctuations***

```
title Paris PN with density fluctuations
```



```

c
c commands for density & abundances =====
fluctuations 16.5 4 3
c this sets up the code like version c84, with fewer chemical elements
init file="c84.ini"
c this sets the abundances of the elements that we will use
abund -1 C-3.523 N-4. O-3.222 ne-3.824 na=-10 mg-4.523 al=-10
continue si-4.523 s-4.824 ar-10 ca=-10 fe-10 ni=-10
c
c commands controlling continuum =====
black body, T = 150000 K, radius = 10
black 5
luminosity total 38
c
c commands controlling geometry =====
sphere
radius 17
c
c other commands for details =====
c
c commands controlling output =====
punch overview last "pn_fluc.ovr"
punch element nitrogen last "pn_fluc.nit"
punch dr last "pn_fluc.dr"
punch results last "pn_fluc.rlt"
c
c pn_fluc.in
c class pn function
c =====

```

The boundary conditions are similar to those for the Paris meeting PN, a homogeneous grain-free PN. This model has density fluctuations, with values chosen so that the mean  $n_e n_p V$  are the same for the nebulae with and without fluctuations.

Checks:

- Check that the fluctuations command works.
- How do results compare with homogeneous Paris pn?

## **pn\_ots *Paris PN with ots***

```

title Paris PN with ots
c recompute "standard" PN model of the Pequignot Meudon Conference
c
c commands controlling continuum =====
black body, T = 150000 K, radius = 10
c
c commands for density & abundances =====
hden = 3.4771213
init file="c84.ini"
abund -1 C-3.523 N-4. O-3.222 ne-3.824 na=-10 mg-4.523 al=-10
continue si-4.523 s-4.824 ar-10 ca=-10 fe-10 ni=-10
c
c commands controlling geometry =====
radius = 17
sphere
c
c other commands for details =====
diffuse ots

```

```

no level2
iterate
atom h-like element hydrogen levels resolved 10
atom h-like element helium levels resolved 15
c
c commands controlling output =====
punch overview last "pn_ots.ovr"
punch element nitrogen last "pn_ots.nit"
punch results last "pn_ots.rlt"
punch dr last "pn_ots.dr"
c
c pn_ots.in
c class pn
c =====
c

```

This is the on-the-spot version of the Paris Planetary Nebula.

Checks:

- Q(H) total 4861 luminosity is close to expected value
- Line spectrum similar to that predicted by default conditions.

## **pn\_paris** *Paris meeting PN*

```

title Paris meeting PN
c standard" PN model of the Pequignot Meudon Conferance
c
c commands controlling continuum =====
black body, T = 150000 K, radius = 10
c
c commands for density & abundances =====
hden = 3.4771213
init file="ism.ini"
abund he -1 C-3.523 N-4. O-3.222 ne-3.824 mg-4.523
continue si-4.523 s-4.824 cl-7 ar-10 fe-10
c
c commands controlling geometry =====
radius = 17
sphere
c
c other commands for details =====
age 10000 years
c
c commands controlling output =====
normalize to "Ca b" 4861
print line optical depths
print ages
print column densities
punch overview "pn_paris.ovr"
punch transmitted continuum "pn_paris.trn"
punch continuum units microns "pn_paris.con"
punch element carbon "pn_paris.car"
punch physical conditions "pn_paris.phy"
punch element nitrogen "pn_paris.nit"
punch dr "pn_paris.dr"
punch temperature "pn_paris.tem"
punch lines emissivity "pn_paris.ems"
h 1 6563
s 2 6720
end of lines
punch averages "pn_paris.avr" last

```

```

temperature hydrogen 1
end of averagers
c
c pn_paris.in
c class pn
c =====
c

```

This is one of the "standard" models computed at the Paris meeting on photoionization and shock calculations. A table in Hazy compares the predictions of the current version of CLOUDY with predictions of a few of the other codes.

The model is meant to represent a planetary nebula ionized by a very hot central object. As a result there is a great deal of He<sup>++</sup> and the associated line and continuum emission is very important.

An age of 10000 years was set. This tests logic in the age command.

## **pn\_paris\_fast** *Paris PN, fast version*

```

title Paris PN, fast version
c standard" PN model of the Pequignot Meudon Conference
c
c commands controlling continuum =====
black body, T = 150000 K, radius = 10
c
c commands for density & abundances =====
hden = 3.4771213
init file="fast.ini"
abund he -1 C-3.523 N-4. O-3.222 ne-3.824 mg-4.523
continue si-4.523 s-4.824 ar-10
c
c commands controlling geometry =====
sphere
radius = 17
c
c other commands for details =====
c
c commands controlling output =====
normalize to "Ca b" 4861
punch overview "pn_paris_fast.ovr"
punch physical conditions "pn_paris_fast.phy"
punch element nitrogen "pn_paris_fast.nit"
punch dr "pn_paris_fast.dr"
c
c pn_paris_fast.in
c class pn
c =====
c

```

This is one of the "standard" models computed at the Paris meeting on photoionization and shock calculations. A table in Hazy compares the predictions of the current version of CLOUDY with predictions of a few of the other codes.

The model is meant to represent a planetary nebula ionized by a very hot central object. As a result there is a great deal of He<sup>++</sup> and the associated line and continuum emission is very important.

## **pn\_sqrden** *PN with density propt $R^{-2}$ and filling factor*

```

title PN with density propt  $R^{-2}$  and filling factor
c this is an optically thin model, which will have a broad range
c of density but the same ionization parameter across the model.
c the physical condtions should not vary by much across the structure
c
c commands controlling continuum =====
c continuum is hot blackbody at  $\sim$ Ledd for solar mass
blackbody, t=5 luminosity=38
c
c commands for density & abundances =====
c the density will fall off as the inverse square of the radius, hence the name
hden 5 -2
init file "ism.ini"
c use default PN abundances but turn off quantum heating since
c we are not going to save the predicted near IR continuum
abundances planetary no qheat
c
c commands controlling geometry =====
c small filling factor so that region will be optically thin
filling factor -2
c set inner and outer radius
radius 16 18
sphere
c
c other commands for details =====
iterate
c
c commands controlling output =====
print line faint 1
punch overview "pn_sqrden.ovr" last
punch dr "pn_sqrden.dr" last
c
c pn_sqrden.in
c class pn
c =====

```

### Checks:

- Zone thickness budgeting handled OK.
  - physical conditions nearly constant across computed structure
- 

## **stars\_atlas** *Atlas stellar atmosphere*

```

title Atlas stellar atmosphere
c
c commands controlling continuum =====
c Approximate model of a hot star ionizing its surroundings.
c Assume H density of  $1e+6$  per cc. This run tests the new Kurucz model option.
c I choose the  $T_{\text{eff}}$  and  $\log(g)$  values arbitrarily. It implies a star
c with luminosity  $32300 L_{\text{sun}}$ , radius  $6.4745 R_{\text{sun}}$ , and mass  $3.745 M_{\text{sun}}$ .
table star atlas 30400.0 4.2
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0

```

```

grains ism
c
c commands controlling geometry =====
radius -3.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
print continuum
c
punch overview last "stars_atlas.ovr"
punch results last "stars_atlas.rlt"
punch transmitted continuum last "stars_atlas.trn" no header
punch continuum last "stars_atlas.con" units microns
c
c stars_atlas.in
c class stars
c =====

```

This is a test that the code can correctly access the large block of Atlas model atmosphere continua described by Kurucz (1991). Kevin Volk provided it as part of his original coding of these stellar atmosphere files. Constant temperature is assumed since this tests shape of stellar continuum rather than thermal physics.

Checks: table star atlas command works.

---

## **stars\_atlas\_3d** *Atlas stellar atmosphere, 3D interpolation*

```

title Atlas stellar atmosphere, 3D interpolation
c
c commands controlling continuum =====
c Approximate model of a hot star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the atlas 3dim model option.
c I choose the T_eff and log(g) values arbitrarily.
table star atlas 3-dim 38400.0 4.82 -1.42
luminosity 4.509202522 solar
table star atlas odfnw 3-dim 38400.0 4.82 0.42
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -3.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_atlas_3d.ovr"

```

```

punch results last "stars_atlas_3d.rlt"
punch transmitted continuum last "stars_atlas_3d.trn" no header
punch continuum last "stars_atlas_3d.con" units microns
c
c stars_atlas_3d.in
c class stars
c =====

```

This is a test that the code can correctly access the large 3-dim block of Atlas model atmosphere continua.

Checks: table star atlas 3dim command works.

## **stars\_atlas\_all *interpolate on Atlas grids***

```

title interpolate on Atlas grids
c
c commands controlling continuum =====
c Include stars with a wide range of metallicities to test access to these grids.
table star atlas Z+1.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z+0.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z+0.3 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z+0.2 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z+0.1 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z+0.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-0.1 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-0.2 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-0.3 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-0.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-1.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-1.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-2.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-2.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-3.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-3.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-4.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-4.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas Z-5.0 38400.0 4.82
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84

```

```

hden 6.0
grains ism
c
c commands controlling geometry =====
radius -2.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_atlas_all.ovr"
punch results last "stars_atlas_all.rlt"
punch transmitted continuum last "stars_atlas_all.trn" no header
punch continuum last "stars_atlas_all.con" units microns
c
c stars_atlas_all.in
c class stars
c =====

```

This is a test that the code can correctly access the 2D grids of various metallicity of Atlas model atmosphere continua.

Checks: access to the various atlas grids

---

## **stars\_atlas\_odfnew\_all *access various Atlas ODFNEW grids***

```

title access various Atlas ODFNEW grids
c
c commands controlling continuum =====
c Include stars with a wide range of metallicities to test access to these grids.
table star atlas odfnew Z+0.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas odfnew Z+0.2 38400.0 4.82
luminosity 4.509202522 solar
table star atlas odfnew Z+0.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas odfnew Z-0.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas odfnew Z-1.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas odfnew Z-1.5 38400.0 4.82
luminosity 4.509202522 solar
table star atlas odfnew Z-2.0 38400.0 4.82
luminosity 4.509202522 solar
table star atlas odfnew Z-2.5 38400.0 4.82
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -2.0 parsec
sphere

```

```

stop zone 1
set dr 0
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_atlas_odfnew_all.ovr"
punch results last "stars_atlas_odfnew_all.rlt"
punch transmitted continuum last "stars_atlas_odfnew_all.trn" no header
punch continuum last "stars_atlas_odfnew_all.con" units microns
c
c stars_atlas_odfnew_all.in
c class stars
c =====

```

This is a test that the code can correctly access the Atlas ODFNEW model atmosphere continua.

Checks: access to the various atlas odfnew grids

---

## **stars\_bstar2006 *tlusty bstar grid***

```

title tlusty bstar grid
c
c commands controlling continuum =====
c Approximate model of a B star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the new Tlusty model option.
c I choose the T_eff and log(g) values arbitrarily.
table star tlusty bstar 20400.0 4.2
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -3.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
c
c in this model the Lyman lines are pumped significantly by the stellar continuum;
c disable this process to avoid critical dependence on the precise shape of the
continuum.
no induced processes
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_bstar2006.ovr"
punch results last "stars_bstar2006.rlt"
punch transmitted continuum last "stars_bstar2006.trn" no header
c
c stars_bstar2006.in
c class stars

```



c =====

This is a test that the code can correctly access the large block of Tlusty B-star model atmosphere continua.

Checks: table star tlusty bstar command works.

## **stars\_bstar2006\_3d *tlusty bstar grid***

```

title tlusty bstar grid
c
c commands controlling continuum =====
c Approximate model of a B star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the new Tlusty model option.
c I choose the T_eff, log(g) and log(Z) values arbitrarily.
table star tlusty bstar 3-dim 20400.0 4.2 -0.573
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -3.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
c
c in this model the Lyman lines are pumped significantly by the stellar continuum;
c disable this process to avoid critical dependence on the precise shape of the
continuum.
no induced processes
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_bstar2006_3d.ovr"
punch results last "stars_bstar2006_3d.rlt"
punch transmitted continuum last "stars_bstar2006_3d.trn" no header
c
c stars_bstar2006_3d.in
c class stars
c =====

```

This is a test that the code can correctly access the large 3-dim block of Tlusty B-star model atmosphere continua.

Checks: table star tlusty bstar 3-dim command works.

## **stars\_bstar2006\_all *tlusty bstar grid***

```

title tlusty bstar grid
c
c commands controlling continuum =====
c Approximate model of 6 B stars ionizing their surroundings.
c Assume H density of 1e+6 per cc. This run tests all the various
c metallicities contained in the BSTAR 2006 grids. I choose the T_eff
c and log(g) values arbitrarily.
table star tlusty bstar Z+0.3 20000.0 4.5
luminosity 4.509202522 solar
table star tlusty bstar Z+0.0 21400.0 4.5
luminosity 4.509202522 solar
table star tlusty bstar Z-0.3 15000.0 4.2
luminosity 4.509202522 solar
table star tlusty bstar Z-0.7 20400.0 4.2
luminosity 4.509202522 solar
table star tlusty bstar Z-1.0 27500.0 3.0
luminosity 4.509202522 solar
table star tlusty bstar Z-INF 18000.0 4.75
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -2.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
c
c in this model the Lyman lines are pumped significantly by the stellar continuum;
c disable this process to avoid critical dependence on the precise shape of the
continuum.
no induced processes
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_bstar2006_all.ovr"
punch results last "stars_bstar2006_all.rlt"
punch transmitted continuum last "stars_bstar2006_all.trn" no header
c
c stars_bstar2006_all.in
c class stars
c =====

```

This is a test that the code can correctly access the large block of Tlusty B-star model atmosphere continua.

Checks: table star tlusty bstar metallicity option works.

## **stars\_costar1 *costar interpolation mode 1***

```

title costar interpolation mode 1
c
c commands controlling continuum =====
table star costar, 40000 K

```

```

ionization parameter -2
table star costar, 31000 K, index 4
ionization parameter -2
c
c commands for density & abundances =====
hden 2
grains orion single no qheat
abundances ism no grains
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
* >>chng 06 jul 01, add this to stars set since want to test shape
* of continuum and resulting ion dist, not thermal equilibrium of nebula
constant temperature 6000K
c
c commands controlling output =====
punch continuum "stars_costar1.con" units microns last
c
c stars_costar1.in
c class stars
c =====
c

```

This is a test of whether the code can read the CoStar stellar atmospheres. It checks the resulting ionization to confirm that the continuum shape was ok. Constant temperature is assumed since this tests shape of stellar continuum rather than thermal physics.

---

## **stars\_costar2 *costar interpolation mode 2***

```

title costar interpolation mode 2
c
c commands controlling continuum =====
table star costar 35575.4 4.2367
ionization parameter -2
table star costar 34468.5 3.86765
ionization parameter -2
table star costar 41741.2 4.02
ionization parameter -2
table star costar 45000 4.1593
ionization parameter -2
table star costar 27500 3.3
ionization parameter -2
c
c commands for density & abundances =====
hden 2
grains orion single no qheat
abundances ism no grains
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
c
c commands controlling output =====
punch continuum "stars_costar2.con" units microns last
c

```

```

c stars_costar2.in
c class stars
c =====
c

```

This is a test of whether the code can read the CoStar stellar atmospheres. It checks the resulting ionization to confirm that the continuum shape was ok.

---

### **stars\_costar3 *costar interpolation mode 3***

```

title costar interpolation mode 3
c
c commands controlling continuum =====
table star costar zams 40.0 Msol 2.85e6
ionization parameter -2
table star costar zams 40.0 Msol 4e6
ionization parameter -2
table star costar zams 50.9 Msol 2.85e6
ionization parameter -2
table star costar zams 42.9 Msol 1.6e6
ionization parameter -2
table star costar zams 70.5 Msol 2.1e6
ionization parameter -2
c
c commands for density & abundances =====
hden 2
grains orion single no qheat
abundances ism no grains
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
c
c commands controlling output =====
punch continuum "stars_costar3.con" units microns last
c
c stars_costar3.in
c class stars
c =====
c

```

This is a test of whether the code can read the CoStar stellar atmospheres. It checks the resulting ionization to confirm that the continuum shape was ok.

---

### **stars\_costar4 *costar interpolation mode 4***

```

title costar interpolation mode 4
c
c commands controlling continuum =====
table star costar age 2.85e6 40.0
ionization parameter -2
table star costar age 4e6 40.0
ionization parameter -2
table star costar age 2.85e6 50.9

```

```

ionization parameter -2
table star costar age 1.6e6 42.9
ionization parameter -2
table star costar age 2.1e6 70.5
ionization parameter -2
c
c commands for density & abundances =====
hden 2
grains orion single no qheat
abundances ism no grains
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
c
c commands controlling output =====
punch continuum "stars_costar4.con" units microns last
c
c stars_costar4.in
c class stars
c =====
c

```

This is a test of whether the code can read the CoStar stellar atmospheres. It checks the resulting ionization to confirm that the continuum shape was ok.

---

## **stars\_costarhalo** *costar halo abundances*

```

title costar halo abundances
c
c commands controlling continuum =====
table star costar, halo abundances, 40000 K
ionization parameter -2
c
c commands for density & abundances =====
hden 2
abundances ism
c
c commands controlling geometry =====
set dr 0
stop zone 1
c
c other commands for details =====
* >>chng 06 jul 01, add this to stars set since want to test shape
* of continuum and resulting ion dist, not thermal equilibrium of nebula
constant temperature 6900K
c
c commands controlling output =====
punch continuum last "stars_costarhalo.con" units microns
c
c
c stars_costarhalo.in
c class stars
c =====
c

```

This test checks that the code can read the CoStar stellar atmospheres. It checks the resulting ionization to make sure that the continuum shape is ok. Constant temperature is assumed since this tests shape of stellar

continuum rather than thermal physics.

---

## **stars\_kurucz79** *Kurucz 1979 SED*

```

title Kurucz 1979 SED
c
c commands controlling continuum =====
c Approximate model of a hot star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the Kurucz model option.
c I choose the T_eff and log(g) values arbitrarily. It implies a star
c with luminosity 32300 L_sun, radius 6.4745 R_sun, and mass 3.745 M_sun.
table star kurucz 30400.0
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -3.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_kurucz79.ovr"
punch results last "stars_kurucz79.rlt"
punch transmitted continuum last "stars_kurucz79.trn" no header
punch continuum last "stars_kurucz79.con" units microns
c
c stars_kurucz79.in
c class stars
c =====

```

This is a test that the code can correctly access the Kurucz model atmosphere continua described by Kurucz (1979). Kevin Volk provided it as part of his original coding of these stellar atmosphere files. Constant temperature is assumed since this tests shape of stellar continuum rather than thermal physics.

Checks: table star kurucz command works.

---

## **stars\_mihalas** *Mihalas SED*

```

title Mihalas SED
c
c commands controlling continuum =====
c Approximate model of a hot star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the new Kurucz model option.
c I choose the T_eff and log(g) values arbitrarily. It implies a star
c with luminosity 32300 L_sun, radius 6.4745 R_sun, and mass 3.745 M_sun.

```

```

table star mihalas 30400.0
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -3.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_mihalas.ovr"
punch results last "stars_mihalas.rlt"
punch transmitted continuum last "stars_mihalas.trn" no header
punch continuum last "stars_mihalas.con" units microns
c
c stars_mihalas.in
c class stars
c =====

```

This is a test that the code can correctly access the small Mihalas grid of NLTE model atmospheres Constant temperature is assumed since this tests shape of continuum not thermal physics.

Checks: table star mihalas command works.

---

## **stars\_optimize1 *optimizer on stellar grids***

```

title optimizer on stellar grids
c
c commands controlling continuum =====
table star tlusty ostar 3-dim 31200 3.6 -1.12 vary
ionization par -3
c
c commands for density & abundances =====
hden 4 vary
init file "hheonly.ini"
element oxygen on
c
c commands controlling geometry =====
case B
stop zone 1
c
c other commands for details =====
optimize tolerance 0.03
constant temper 4.3 vary
optimize phymir sequential
optimize iteration 500
c the following spectrum was derived by running this sim at hden 5, t=4,
c and T_star=51200
c resulting eden was 5.040 and T=4, which we shall assert that we find
optimize lines

```

```

o ii 3729 0.0416
o ii 3726 0.1179
o ii 7323 0.1535
o ii 7332 0.1237
o 3 5007 9.5265
totl 4363 0.0900
end of lines
c
c commands controlling output =====
print line faint -2
c
c
c optimize_phymir.in
c class optimizer
c =====
c

```

This checks whether the optimizer sets correct limits for Teff. The line spectrum was calculated at  $T = 1e4$  K,  $n_H = 1e5 \text{ cm}^{-3}$ ,  $T_{\text{star}} = 51200$  K.

---

## **stars\_optimize2 *optimizer on stellar grids***

```

title optimizer on stellar grids
c
c commands controlling continuum =====
table star costar 36200 vary
ionization par -3
c
c commands for density & abundances =====
hden 4 vary
init file "hheonly.ini"
element oxygen on
c
c commands controlling geometry =====
Case B
stop zone 1
c
c other commands for details =====
optimize tolerance 0.03
constant temper 4.3 vary
optimize phymir sequential
optimize iteration 500
c the following spectrum was derived by running this sim at hden 5, t=4,
c and T_star=51200
c resulting eden was 5.041 and T=4, which we shall assert that we find
optimize lines
o ii 3729 0.0335
o ii 3726 0.0950
o ii 7323 0.1238
o ii 7332 0.0998
o 3 5007 9.9714
totl 4363 0.0943
end of lines
c
c commands controlling output =====
print line faint -2
c
c
c optimize_phymir.in
c class optimizer

```



```
c =====
c
```

This checks whether the optimizer sets corrects limits for Teff. The line spectrum was calculated at  $T = 1e4$  K,  $n_H = 1e5 \text{ cm}^{-3}$ ,  $T_{\text{star}} = 51200$  K.

---

### **stars\_optimize3 *optimizer on stellar grids***

```
title optimizer on stellar grids
c
c commands controlling continuum =====
table star costar age 5.8 log 36 msol vary
ionization par -3
c
c commands for density & abundances =====
hden 4 vary
init file "hheonly.ini"
element oxygen on
c
c commands controlling geometry =====
Case B
stop zone 1
c
c other commands for details =====
optimize tolerance 0.03
constant temper 4.3 vary
optimize phymir sequential
optimize iteration 500
c the following spectrum was derived by running this sim at hden 5, t=4,
c and log(age)=6.3
c resulting eden was 5.040 and T=4, which we shall assert that we find
optimize lines
o ii 3729 0.0623
o ii 3726 0.1765
o ii 7323 0.2300
o ii 7332 0.1854
o 3 5007 8.2979
totl 4363 0.0784
end of lines
c
c commands controlling output =====
print line faint -2
c
c
c optimize_phymir.in
c class optimizer
c =====
c
```

This checks whether the optimizer sets corrects limits for  $\log(\text{age})$ . The line spectrum was calculated at  $T = 1e4$  K,  $n_H = 1e5 \text{ cm}^{-3}$ ,  $\log(\text{age}/\text{yr}) = 6.3$ .

---

### **stars\_ostar2002 *tlusty O star grid***

```
title tlusty O star grid
c
```

```

c commands controlling continuum =====
c Approximate model of a hot star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the new Tlusty model option.
c I choose the T_eff and log(g) values arbitrarily. It implies a star
c with luminosity 32300 L_sun, radius 6.4745 R_sun, and mass 3.745 M_sun.
table star tlusty ostar 30400.0 4.2
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -3.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
c
c in this model the Lyman lines are pumped significantly by the stellar continuum;
c disable this process to avoid critical dependence on the precise shape of the
continuum.
no induced processes
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_ostar2002.ovr"
punch results last "stars_ostar2002.rlt"
punch transmitted continuum last "stars_ostar2002.trn" no header
c
c stars_ostar2002.in
c class stars
c =====

```

This is a test that the code can correctly access the large block of Tlusty O-star model atmosphere continua.

Checks: table star tlusty ostar command works.

## **stars\_ostar2002\_3d *tlusty Ostar grid***

```

title tlusty Ostar grid
c
c commands controlling continuum =====
c Approximate model of a hot star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the new Tlusty model option.
c I choose the T_eff, log(g) and log(Z) values arbitrarily.
table star tlusty ostar 3-dim 30400.0 4.2 -1.573
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -3.0 parsec

```

```

sphere
stop zone 1
set dr 0
c
c other commands for details =====
c
c in this model the Lyman lines are pumped significantly by the stellar continuum;
c disable this process to avoid critical dependence on the precise shape of the
continuum.
no induced processes
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_ostar2002_3d.ovr"
punch results last "stars_ostar2002_3d.rlt"
punch transmitted continuum last "stars_ostar2002_3d.trn" no header
c
c stars_ostar2002_3d.in
c class stars
c =====

```

This is a test that the code can correctly access the large block of Tlusty O-star model atmosphere continua.

Checks: table star tlusty ostar 3-dim command works.

---

## **stars\_ostar2002\_all *tlusty Ostar grid***

```

title tlusty Ostar grid
c
c commands controlling continuum =====
c Approximate model of 10 hot stars ionizing their surroundings.
c Assume H density of 1e+6 per cc. This run tests all the various
c metallicities contained in the OSTAR 2002 grids. I choose the T_eff
c and log(g) values such that the selection algorithm gets a good workout.
table star tlusty ostar Z+0.3 30000.0 4.5
luminosity 4.509202522 solar
table star tlusty ostar Z+0.0 31400.0 4.5
luminosity 4.509202522 solar
table star tlusty ostar Z-0.3 45000.0 4.2
luminosity 4.509202522 solar
table star tlusty ostar Z-0.7 30400.0 4.2
luminosity 4.509202522 solar
table star tlusty ostar Z-1.0 27500.0 3.0
luminosity 4.509202522 solar
table star tlusty ostar Z-1.5 55000.0 4.75
luminosity 4.509202522 solar
table star tlusty ostar Z-1.7 27500.0 4.75
luminosity 4.509202522 solar
table star tlusty ostar Z-2.0 55000.0 3.00
luminosity 4.509202522 solar
table star tlusty ostar Z-3.0 44400.0 3.2
luminosity 4.509202522 solar
table star tlusty ostar Z-INF 53000.0 4.75
luminosity 4.509202522 solar
table star tlusty ostar Z+0.0 48400.0 3.87
luminosity 4.509202522 solar
table star tlusty ostar Z-0.7 48400.0 3.55
luminosity 4.509202522 solar

```

```

table star tlusty ostar Z+0.3 43400.0 3.65
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -2.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
c
c in this model the Lyman lines are pumped significantly by the stellar continuum;
c disable this process to avoid critical dependence on the precise shape of the
continuum.
no induced processes
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_ostar2002_all.ovr"
punch results last "stars_ostar2002_all.rlt"
punch transmitted continuum last "stars_ostar2002_all.trn" no header
c
c stars_ostar2002_all.in
c class stars
c =====

```

This is a test that the code can correctly access the large block of Tlusty O-star model atmosphere continua.

Checks: table star tlusty ostar metallicity option works, also test selection algorithm.

---

## **stars\_rauch\_3d** *Rauch 3-dimensional grid*

```

title Rauch 3-dimensional grid
c
c commands controlling continuum =====
table star rauch 3-dim T = 150000 K, g = 6.5 , log(Z) = -0.6
luminosity 4.5 solar
table star rauch old 3-dim T = 233000 K, g = 7.5 , log(Z) = -0.32
luminosity 4.5 solar
c
c commands controlling geometry =====
sphere
radius 18
stop zone 1
set dr 0
c
c commands for density & abundances =====
abundances planetary
hden 3.0
c
c other commands for details =====
no level2
c

```

```

c commands controlling output =====
plot continuum range .1
punch overview last "stars_rauch_3d.ovr"
punch results last "stars_rauch_3d.rlt"
punch continuum last "stars_rauch_3d.con" units microns
c
c stars_rauch_3d.in
c class stars
c =====

```

This model checks that the code is able to access the 3-dim versions of the Rauch grids of stellar atmosphere models. Constant temperature is set since this test is intended to test shape of ionizing continuum, not thermal physics.

---

## **stars\_rauch\_h+he *Rauch PN central star***

```

title Rauch PN central star
c
c commands controlling continuum =====
c this tests the various interpolation modes of the 3d grid
table star rauch h+he T = 150000 K, g = 6.5 , f(He) = 0.2
luminosity 4.5 solar
table star rauch h+he T = 155000 K, g = 6.5 , f(He) = 0.2
luminosity 4.5 solar
table star rauch h+he T = 150000 K, g = 6.25 , f(He) = 0.2
luminosity 4.5 solar
table star rauch h+he T = 150000 K, g = 6.5 , f(He) = 0.25
luminosity 4.5 solar
table star rauch h+he T = 155000 K, g = 6.25 , f(He) = 0.2
luminosity 4.5 solar
table star rauch h+he T = 155000 K, g = 6.5 , f(He) = 0.25
luminosity 4.5 solar
table star rauch h+he T = 150000 K, g = 6.25 , f(He) = 0.25
luminosity 4.5 solar
table star rauch h+he T = 155000 K, g = 6.25 , f(He) = 0.25
luminosity 4.5 solar
c
c commands controlling geometry =====
sphere
radius 18
stop zone 1
set dr 0
c
c commands for density & abundances =====
abundances planetary
hden 3.0
c
c other commands for details =====
no level2
c
c commands controlling output =====
plot continuum range .1
punch overview last "stars_rauch_h+he.ovr"
punch results last "stars_rauch_h+he.rlt"
punch continuum last "stars_rauch_h+he.con" units microns
c
c stars_rauch_h+he.in
c class stars
c =====

```

This is a model of a very hot planetary nebula, and checks that the code is able to access the H+He version of the Rauch grid of stellar atmosphere models.

---

## **stars\_rauch\_h-ca** *Rauch central star hot PN*

```

title Rauch central star hot PN
c NB
c this uses the H-Ca rauch atmospheres
c
c commands controlling continuum =====
table star old rauch T = 250000 K, g = 7.5
luminosity 4.5 solar
c
c commands controlling geometry =====
sphere
radius 18
stop zone 1
set dr 0
c
c other commands for details =====
no level2
c
c commands for density & abundances =====
hden 3.0
abundances planetary
c
c commands controlling output =====
plot continuum range .1
punch overview last "stars_rauch_h-ca.ovr"
punch results last "stars_rauch_h-ca.rlt"
punch continuum last "stars_rauch_h-ca.con" units microns
c
c stars_rauch_h-ca.in
c class stars
c =====

```

This is a model of a very hot planetary nebula, and checks that the code is able to access the H-Ca Rauch grid of stellar atmosphere models.

---

## **stars\_rauch\_h-ni** *Rauch hot PN star*

```

title Rauch hot PN star
c
c commands controlling continuum =====
table star rauch T = 150000 K, g = 6.5
luminosity 4.5 solar
c
c commands controlling geometry =====
sphere
radius 18
stop zone 1
set dr 0
c
c commands for density & abundances =====
abundances planetary
hden 3.0

```

```

c
c other commands for details =====
no level2
c
c commands controlling output =====
plot continuum range .1
punch overview last "stars_rauch_h-ni.ovr"
punch results last "stars_rauch_h-ni.rlt"
punch continuum last "stars_rauch_h-ni.con" units microns
c
c stars_rauch_h-ni.in
c class stars
c =====

```

This is a model of a very hot planetary nebula, and checks that the code is able to access the H-Ni version of the Rauch grid of stellar atmosphere models. Constant temperature is set since this test is intended to test shape of ionizing continuum, not thermal physics.

---

### **stars\_rauch\_helium** *Rauch hot PN star*

```

title Rauch hot PN star
c
c commands controlling continuum =====
table star rauch helium T = 240000 K, g = 8.5
luminosity 4.5 solar
c
c commands controlling geometry =====
sphere
radius 18
stop zone 1
set dr 0
c
c commands for density & abundances =====
abundances planetary
hden 3.0
c
c other commands for details =====
no level2
c
c commands controlling output =====
plot continuum range .1
punch overview last "stars_rauch_helium.ovr"
punch results last "stars_rauch_helium.rlt"
c
c stars_rauch_helium.in
c class stars
c =====

```

This is a model of a very hot planetary nebula, and checks that the code is able to access the pure helium version of the Rauch grid of stellar atmosphere models.

---

### **stars\_rauch\_hydr** *Rauch hot PN star*

```

title Rauch hot PN star
c
c commands controlling continuum =====

```

```

table star rauch hydr T = 240000 K, g = 8.5
luminosity 4.5 solar
c
c commands controlling geometry =====
sphere
radius 18
stop zone 1
set dr 0
c
c commands for density & abundances =====
abundances planetary
hden 3.0
c
c other commands for details =====
no level2
c
c commands controlling output =====
plot continuum range .1
punch overview last "stars_rauch_hydr.ovr"
punch results last "stars_rauch_hydr.rlt"
c
c stars_rauch_hydr.in
c class stars
c =====

```

This is a model of a very hot planetary nebula, and checks that the code is able to access the pure hydrogen version of the Rauch grid of stellar atmosphere models.

---

## **stars\_rauch\_pg1159 *Rauch hot [WR] PN***

```

title Rauch hot [WR] PN
c
c commands controlling continuum =====
table star rauch pg1159 T=150000K g=6.5
luminosity 4.5 solar
c
c commands controlling geometry =====
sphere
radius 18
stop zone 1
set dr 0
c
c commands for density & abundances =====
abundances planetary
hden 3.0
c
c other commands for details =====
no level2
c
c commands controlling output =====
plot continuum range .1
punch overview last "stars_rauch_pg1159.ovr"
punch results last "stars_rauch_pg1159.rlt"
c
c stars_rauch.in
c class stars
c =====

```

This is a model of a very hot planetary nebula, and checks that the code is able to access the PG1159 version of the Rauch grid of stellar atmosphere models.



---

## **stars\_starburst99** *Starburst 99 SED*

```

title Starburst 99 SED
c
c commands controlling continuum =====
c read the table starburst and interploate to an age of 1e8 years
table star log age=8 "starburst99.mod"
luminosity 43
c add background continuum - starburs 99 does not extend
c into fir so code would complain about zero continuum
c adding the cosmic background stops this complaint
background
c
c commands for density & abundances =====
hden 5
c
c commands controlling geometry =====
stop zone 1
radius 20
c
c other commands for details =====
constant temper 4
c
c commands controlling output =====
punch continuum "stars_starburst99.con" last
punch continuum last "stars_starburst99.con" units microns
c
c
c stars_starburst99.in
c class stars
c =====
c

```

this is a demonstration of the use of a Starburst 99 spectrum. It was provided to me by anand Srianand, and we used it in Srianand et al. 2003. A constant temperature is set since this is to test shape of continuum not thermal physics. It also tests access to user-supplied grids as the mechanism is the same.

---

## **stars\_starburst99\_2d** *Starburst 99 2D SED*

```

title Starburst 99 2D SED
c
c commands controlling continuum =====
c read the table starburst and interpolate to an age of 2.09e6 years
c and a metallicity of 12.
table star log age=6.32 Z=12. "starburst99_2d.mod"
luminosity 43
c add background continuum - starburs 99 does not extend
c into fir so code would complain about zero continuum
c adding the cosmic background stops this complaint
background
c
c commands for density & abundances =====
hden 5
c
c commands controlling geometry =====
stop zone 1

```

```

radius 20
c
c other commands for details =====
constant temper 4
c
c commands controlling output =====
punch continuum "stars_starburst99_2d.con" last
punch continuum last "stars_starburst99_2d.con" units microns
c
c
c stars_starburst99_2d.in
c class stars
c =====
c

```

this is a demonstration of the use of a manually crafted Starburst 99 2D grid, allowing for interpolation in both age and metallicity. It was provided by Christophe Morisset. A constant temperature is set since this is to test shape of continuum not thermal physics. It also tests access to user-supplied grids as the mechanism is the same.

## **stars\_werner** *Werner stars grid*

```

title Werner stars grid
c
c commands controlling continuum =====
table star werner 190000 g=7.5
luminosity total 38
c
c commands for density & abundances =====
hden 4
abundances planetary no grains
grains agm no qheat single
c
c commands controlling geometry =====
radius 17
sphere
set dr 0
stop zone 1
c
c other commands for details =====
* >>chng 06 jul 01, add this to stars set since want to test shape
* of continuum and resulting ion dist, not thermal equilibrium of nebula
constant temperature 4.373
c
c commands controlling output =====
c
punch overview last "stars_werner.ovr"
punch results last "stars_werner.rlt"
punch transmitted continuumlast "stars_werner.trn"
c
c stars_werner.in
c class stars
c =====
c

```

This checks that the code can access Kevin Volk39s Werner atmospheres.

Checks:

- table star Werner atmosphere works. constant temperature is set to that this tests shape of stellar continuum

---

## stars\_wmbasic *wmbasic stellar SED*

```

title wmbasic stellar SED
c
c commands controlling continuum =====
c Approximate model of a hot star ionizing its surroundings.
c Assume H density of 1e+6 per cc. This run tests the WMBASIC model option.
c I choose the T_eff, log(g), and log(Z) values arbitrarily.
table star wmbasic 30000.0 4.0 -0.3
luminosity 4.509202522 solar
table star wmbasic 32000.0 4.0 0.0
luminosity 4.509202522 solar
table star wmbasic 40000.0 3.6 -0.3
luminosity 4.509202522 solar
table star wmbasic 35000.0 4.0 -0.15
luminosity 4.509202522 solar
table star wmbasic 57000.0 4.0 0.0
luminosity 4.509202522 solar
table star wmbasic 37000.0 3.4 -0.3
luminosity 4.509202522 solar
table star wmbasic 32000.0 4.0 -0.1
luminosity 4.509202522 solar
table star wmbasic 36000.0 3.5 -0.12
luminosity 4.509202522 solar
c
c commands for density & abundances =====
abundances old solar 84
hden 6.0
grains ism
c
c commands controlling geometry =====
radius -2.0 parsec
sphere
stop zone 1
set dr 0
c
c other commands for details =====
c
c in this model the Lyman lines are pumped significantly
c by the stellar continuum; disable this process to avoid critical
c dependence on the precise shape of the continuum.
no induced processes
constant temperature 4
c
c commands controlling output =====
c
punch overview last "stars_wmbasic.ovr"
punch results last "stars_wmbasic.rlt"
punch transmitted continuum last "stars_wmbasic.trn" no header
c
c stars_wmbasic.in
c class stars
c =====

```

This is a test that the code can correctly access the WMBASIC O-star model atmosphere continua.

Checks: table star wmbasic command works.

