

Status of Modtran/LibRadTran simulation and Comparison in december 2016

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For PCWG team of DESC-LSST collaboration

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Overview

- 1 Reminder on the specifications
- 2 The prescription
- 3 Practical file organization
- 4 Example of a comparison inside LibRadTran

Data points of the grid

Altitude

$h = 2.75$ km.

Wavelength range

$\lambda \in [250, 1200]$ nm, $\Delta\lambda = 1$ nm

Airmass range

z computed for 31 points in steps of 0.1 : $z \in [1, 3]$

The Atmospheric Profile

Compare predictions for 2 realistic atmospheric models.

- US standard atmosphere → probably good for LSST, but perhaps a little too wet
- Subarctic winter → very dry air, probably well suited for LSST site

The above atmospheric models are both available in Modtran(Data Card : Card1.MODEL) and LibRadTran (Data Card : atmosphere_file.)

Main absorption models used in LibRadTran

- Representative wavelengths parameterization (REPTRAN),
- Pseudo-spectral calculation adapted from LOWTRAN.
- CRS: switching off spectral parameterization.

Absorption models in LibRadTran :

1) REPTRAN, 2) LOWTRAN, 3) CRS,

The absorption models in Modtran

- **Band models methods**, The Modtran or Lowtran models,
- **The correlated-k method**, The Modtran and Lowtran either in slow or medium modes.

Absorption models in Modtran :

1) MODTRAN band Model, 2) LOWTRAN Band Model , 3) MODTRAN correlated-k and slow speed, 4) MODTRAN correlated-k and slow speed, 5) MODTRAN correlated-k and medium speed

The Data card for specifying the absorption model in Modtran is *CARD1 MODTRN* . One should provides the values *T,M,C,K,F,L* and *CARD1 SPEED* is *S* or blank.

Different modes : Selected interaction processes

- Simulate with only molecular scattering (Rayleigh scattering) (code name *sc*).
- **Simulate pure molecular absorption (code name *ab*).**

Variation of Precipitable Pressure Water PWV

absorption profiles for 31 points $p_{wv} \in [0 \text{ mm}, 15 \text{ mm}]$ in steps of 0.5 mm

Variation of Ozone O_3

absorption profiles for 21 points $O_3 \in [200 \text{ Dobson}, 600 \text{ Dobson}]$.

- **Simulate the combination of molecular scattering and molecular absorption (code name *sa*).**
 - perform similar variations in PWV and O_3

File naming : Part 1

filename

$P_O_ \{rte\} _ \{atm\} _ \{proc\} _ \{mod\} _ zXX _ wwXX _ ozXX . extension$

P : **RT** or **MT** for LibRadtran or Modtran,

O : Observatory site **LS** or **HP** or **GM** or **MK** for LSST, OHP, Gemini South, Mauna Kea,...

{rte} : **pp** or **ps**,

{atm} : **us** or **sw**,

{proc} : **sc** for pure molecular scattering, **ab** for pure molecular absorption, **sa** for the combination of molecular scattering and absorption,

{mod} : in LibRadTran : **rt** for Reptran model, **lt** for Lowtran model, **cr** for CRS model, **fu** for the Fu and Liou model, **k2** for Kato2 model, and **kt** for Kato model.

{mod} : in ModTran : **mt** for Modtran band model), **mk** for Modtran correlated-k model, **lt** for Lowtran model.

zXX : Airmass z , where XX is the value of the airmass on 2 digit $XX = 2 \times z$,

wwXX : Precipitable water vapour pwv , where XX is the value of the pwv on 3 digit $XXX = 10 \times pwv$, pwv in mm unit,

ozXX : Ozone oz , where XX is the value of the oz on 2 digit $XX = oz/10$, oz is Dobson unit.

For example a filename for LibRadtran, for LSST could be :

RT_LS_pp_us_sa_rt_z15_ww030_oz30.txt

where $pwv = 3$ mm and $oz = 300$ Dobson unit and $z = 1.5$.

Example of US standard atmosphere, with Reptran, for airmass $z = 1$ and $z = 2$, when varying PWV :

See repository at :

<https://github.com/LSSTDESC/PC5AtmosphericExtinction/tree/master/LibRadTran/simulations/RT/2.0/LS/pp/us>

```
ls
```

```
PC5AtmosphericExtinction/LibRadTran/simulations/RT/2.0/LS/pp/us
```

```
RT_LS_pp_us_ab_rt_z10_wv0.0OUT RT_LS_pp_us_ab_rt_z20_wv30.0OUT
RT_LS_pp_us_ab_rt_z10_wv10.0OUT RT_LS_pp_us_ab_rt_z20_wv35.0OUT
RT_LS_pp_us_ab_rt_z10_wv100.0OUT RT_LS_pp_us_ab_rt_z20_wv40.0OUT
RT_LS_pp_us_ab_rt_z10_wv105.0OUT RT_LS_pp_us_ab_rt_z20_wv45.0OUT
RT_LS_pp_us_ab_rt_z10_wv110.0OUT RT_LS_pp_us_ab_rt_z20_wv5.0OUT
RT_LS_pp_us_ab_rt_z10_wv115.0OUT RT_LS_pp_us_ab_rt_z20_wv50.0OUT
RT_LS_pp_us_ab_rt_z10_wv120.0OUT RT_LS_pp_us_ab_rt_z20_wv55.0OUT
RT_LS_pp_us_ab_rt_z10_wv125.0OUT RT_LS_pp_us_ab_rt_z20_wv60.0OUT
RT_LS_pp_us_ab_rt_z10_wv130.0OUT RT_LS_pp_us_ab_rt_z20_wv65.0OUT
RT_LS_pp_us_ab_rt_z10_wv135.0OUT RT_LS_pp_us_ab_rt_z20_wv70.0OUT
RT_LS_pp_us_ab_rt_z10_wv140.0OUT RT_LS_pp_us_ab_rt_z20_wv75.0OUT
```

File naming : Part 2

Naming absorption models for Modtran

filename

*P_O-**{rte}**-**{atm}**-**{proc}**-**{mod}**_zXX_wvXX_ozXX.extension*

- {mod}** :
- **mt** for Modtran band model (CARD1.MODTRN='T'),
 - **mm** for Modtran band model (CARD1.MODTRN='M'),,
 - **lt** for Lowtran band model (CARD1.MODTRN='L' or 'F'),
 - **mks** for Modtran correlate-k in slow mode (speed), (CARD1.MODTRN='K' or 'C' and CARD1.SEED='S'),
 - **mkm** for Modtran correlate-k in medium mode (speed), (CARD1.MODTRN='K' or 'C' and CARD1.SEED='M'),

For example a filename for Modtran, for LSST could be :

MT_LS_pp_us_sa_mt_z15_wv030_oz30.txt

where *pwv* = 3 mm and *oz* = 300 Dobson unit.

Example of Hierarchy of directories

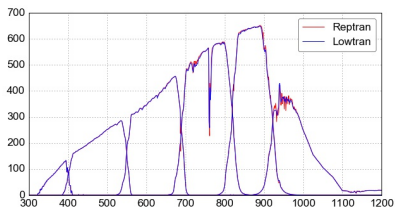
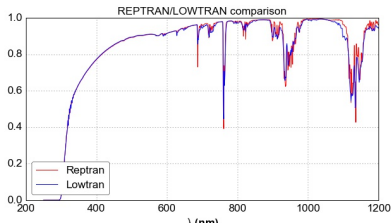
rootdir	top directory
rootdir/RT/VXX/LS/pp/us/sc/	pure molecular scattering
rootdir/RT/VXX/LS/pp/us/ab/rt/wv/	pure molecular absorption for varying <i>pwv</i>
rootdir/RT/VXX/LS/pp/us/ab/rt/oz/	pure molecular absorption for varying <i>oz</i>
rootdir/RT/VXX/LS/pp/us/sa/rt/wv/	molecular absorption and scattering for varying <i>pwv</i>
rootdir/RT/VXX/LS/pp/us/sa/rt/oz/	molecular absorption scattering for varying <i>oz</i>
rootdir/RT/VXX/LS/pp/us/sc/ae/	pure molecular scattering and default aerosols scattering
rootdir/RT/VXX/LS/pp/us/sa/rt/ae/	molecular absorption scattering and default aerosols scattering
rootdir/RT/VXX/LS/pp/us/sc/as/	pure molecular scattering and special aerosols scattering
rootdir/RT/VXX/LS/pp/us/sa/rt/as/	molecular absorption scattering and special aerosols scattering

- *RT* means LibRadTran
- *VXX* = 2.0 is the version number of Radtran or Modtran.
- *LS* means LSST site,
- *pp* means geometry of parallel planes for the atmosphere

Example of comparison inside LibRadTran : REPTRAN vs LOWTRAN

Magnitude calculation

$$F_{\Delta\lambda}^{ADU} = \frac{\pi D^2}{4g_{el}hc} \int_{\Delta\lambda} T^{atm}(\lambda) T^{filt}(\lambda) \epsilon_{CCD}(\lambda) S_{\lambda}^E(\lambda) \lambda d\lambda$$



Example

For a flat SED $S_{\lambda}^E(\lambda) = cte$, with $z = 1$:

filter	mag-shift (mmag)	filter	mag-shift (mmag)	filter	mag-shift (mmag)
U	- 3.0	I	+ 8.9	G	-2.6
Z	7.3	R	+ 1.5	Y4	4.5

Status of the work

- GitHub Repository created by Nicolas at <https://github.com/DarkEnergyScienceCollaboration/PC5AtmosphericExtinction>
 - Prescription note written to proceed with air transparency simulations and posted at <https://github.com/DarkEnergyScienceCollaboration/PC5AtmosphericExtinction/tree/master/doc/Prescriptions>
 - Simulation with LibRadTran **done** and at posted at (see README) <https://github.com/DarkEnergyScienceCollaboration/PC5AtmosphericExtinction/tree/master/LibRadTran/simulations>
 - Simulation with Modtran **soon performed** and at posted at (see README) <https://github.com/DarkEnergyScienceCollaboration/PC5AtmosphericExtinction/tree/master/ModTran>
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- Ready for Atmospheric properties Analysis,
 - Almost ready for comparison, at least comparing the different models inside LibRadTran.

The End