



# Monte-Carlo Spectral Integration in COSMO Radiation Scheme



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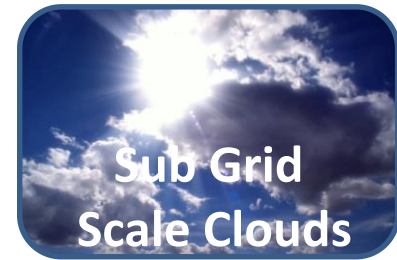
ICCARUS Seminar, DWD Offenbach, February 26, 2018

# Outline

- General description of radiative transfer calculation methods
- COSMO radiation scheme RG92
- What is Monte-Carlo Spectral Integration?
- The COSMO-MCSI setup
- Run time and errors comparisons
- Model performance verifications against observational data
- Conclusions

# Testing & Tuning of Revised Cloud Radiation Coupling - $T^2(RC)^2$

Revised  
Cloud/aerosols  
Radiation  
Coupling



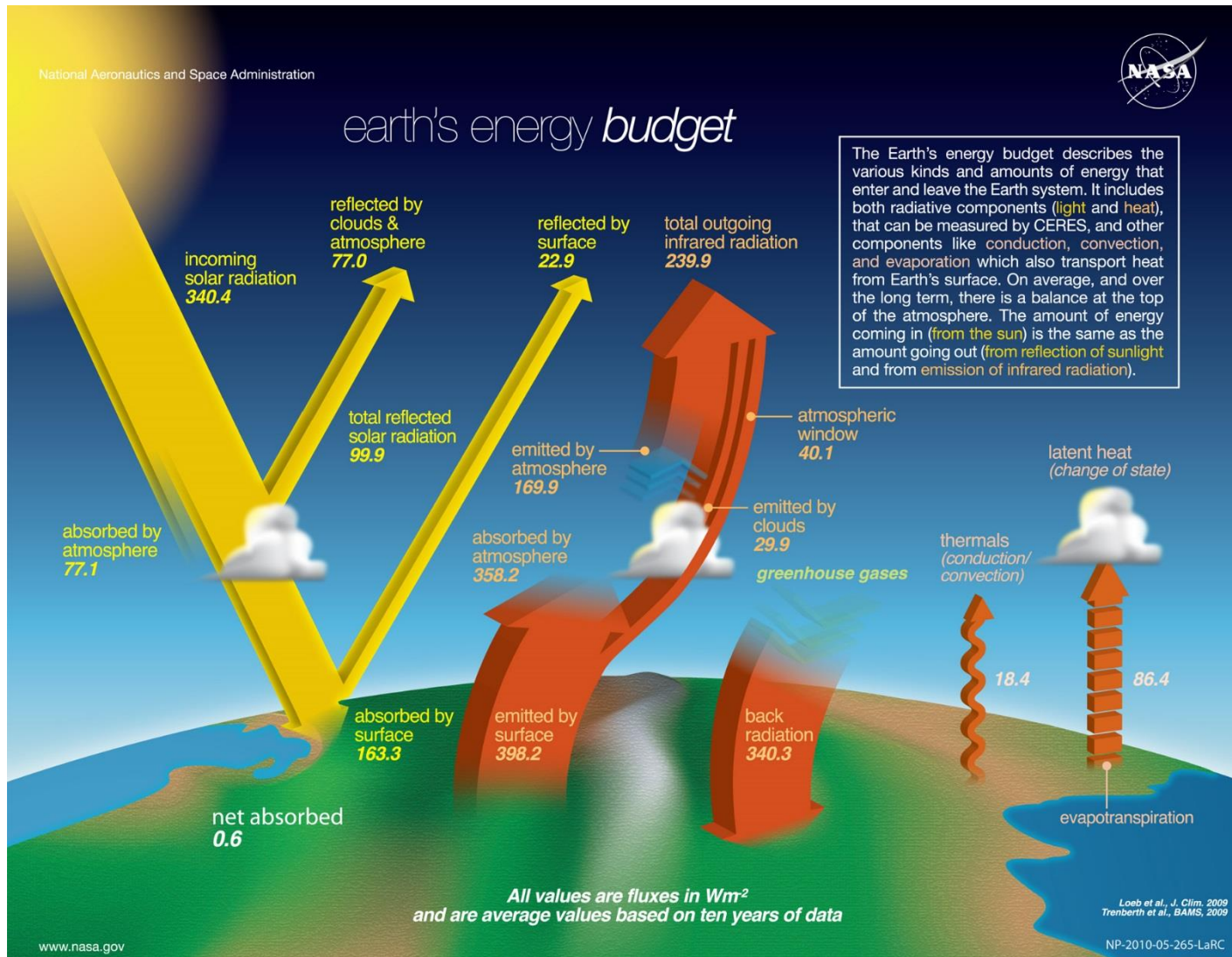
Testing &  
Tuning



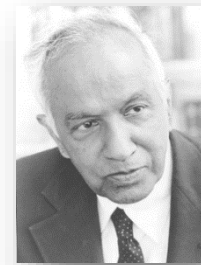
Run Time  
Optimization



# Earth's Energy Budget



# Chandrasekhar's General Radiative Transfer Equation (RTE)



$$\mu \frac{dI_v(\tau_v, \mu, \phi)}{d\tau_v} = I_v(\tau_v, \mu, \phi) - \varpi_v(\tau_v) \frac{j_v^{sca}(\tau_v, \mu, \phi)}{\sigma_{sca}(\nu)} - (1 - \varpi_v(\tau_v)) \frac{j_v^{thermal\ emi}(\tau_v)}{\sigma_{abs}(\nu)}$$

$I_v$  - radiance

$\tau_v$  - extinction optical thickness (abs. + sca.) – here treated as vertical coordinate

$\mu$  - cos(Solar zenith angle)

$\phi$  - azimuth angle

$\varpi_v$  - single scattering albedo, ( $\varpi_v=0$  “pure abs”,  $\varpi_v=1$  “pure sca”)

$$\varpi_v = \frac{n_{sca}\sigma_{sca}}{n_{sca}\sigma_{sca} + n_{abs}\sigma_{abs}}$$

$\sigma_{sca/abs}(\nu)$  - scattering/absorption cross section  $j_v^{sca/emi}(\tau_v, \mu, \phi)$  - sca./emission ability

$$\mu \frac{dI_v(\tau_v, \mu, \phi)}{d\tau_v} = I_v(\tau_v, \mu, \phi) - \varpi_v(\tau_v) \int_{4\pi} d\Omega' \frac{1}{4\pi} P_v(\mu', \phi', \mu, \phi) I_v(\tau_v, \mu', \phi') - (1 - \varpi_v(\tau_v)) B_v(T(\tau_v))$$

$P_v(\mu', \phi', \mu, \phi)$  - scattering phase function (probability that ray from  $\mu', \phi'$  will scatter to  $\mu, \phi$ )

$B_v(T(\tau_v))$  - Planck's Function

# Assumptions we make for simplicity

- Plane parallel atmosphere
- Local thermal equilibrium (LTE) – only thermal emissions considered
- Two-stream approximation (1D problem)  $\downarrow \uparrow$ 
  - Isotropic scattering for half sphere (many scatterings)
  - Rayleigh/Henyey-Greenstein phase functions

$$P_{Ray}(\theta) = \frac{3}{4}(1 + \cos^2 \theta)$$

$$P_{HG}(\theta) = \frac{1 - g^2}{(1 + g^2 - 2g\cos\theta)^{3/2}} \begin{cases} 0^\circ & \frac{1 + g}{(1 - g)^2} \\ 180^\circ & \frac{1 - g}{(1 + g)^2} \end{cases}$$

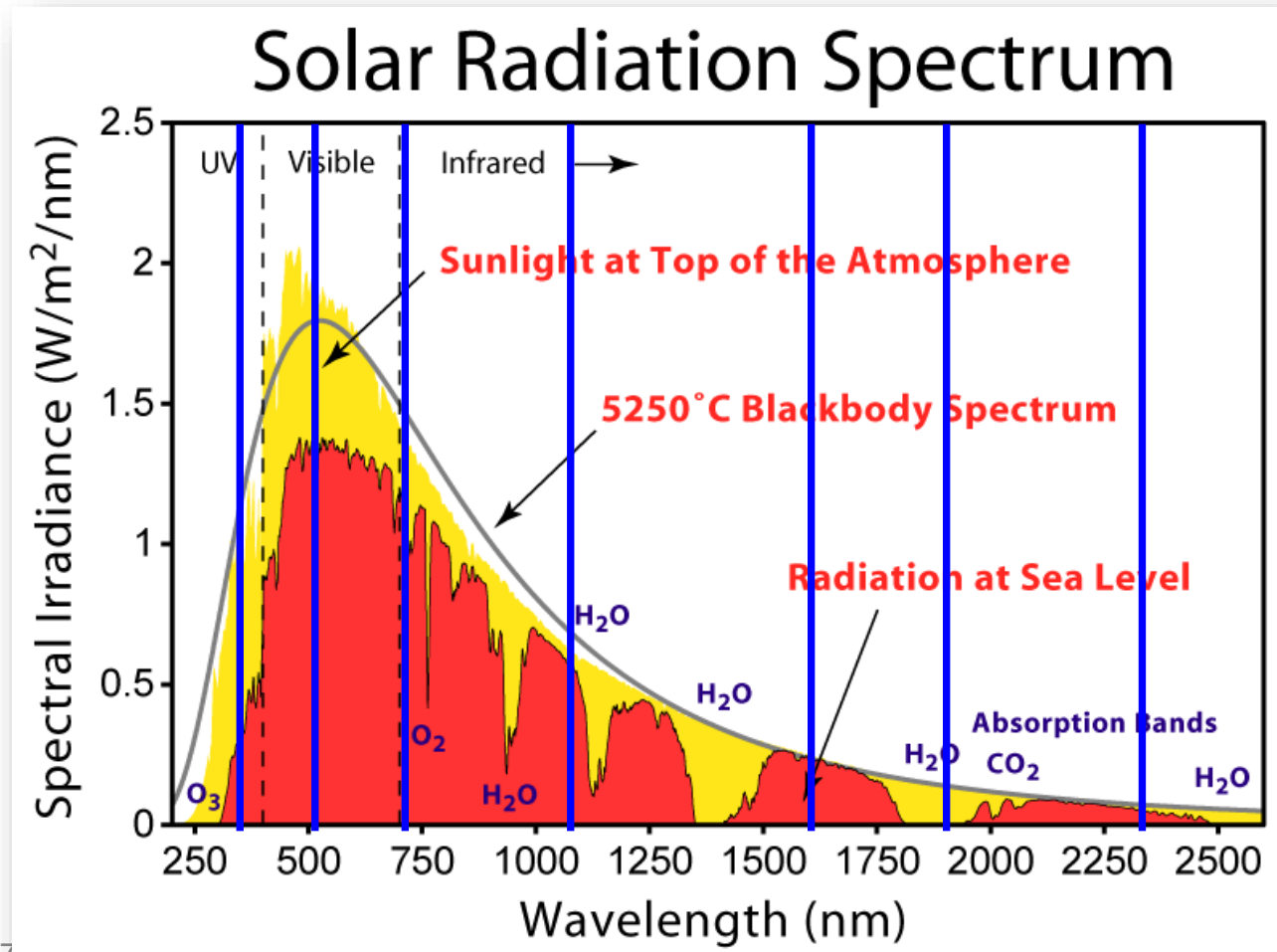
$$g = \frac{1}{2} \int_{-1}^1 P(\cos\theta) \cos\theta d\cos\theta$$

$$g = \begin{cases} 1 & \text{forward} \\ 0 & \\ -1 & \text{backwards} \end{cases}$$



# And that's not all...

The RTE needs to be computed for each  $(x, y, z, \nu)$  and separately for each of the gases, aerosols, hydrometeors



COSMO radiation:

3 Visible bands

5 Thermal bands

# The k-distribution Method

- For gases ( $H_2O$ ,  $CO_2$ ,  $O_3$ ...) the absorption is rapidly changing as a function wavelength. Line by line (LBL) methods are too expensive for NWP
- In the k-distribution method gases absorption spectra for each band is transformed from wavelength to cumulative probability space

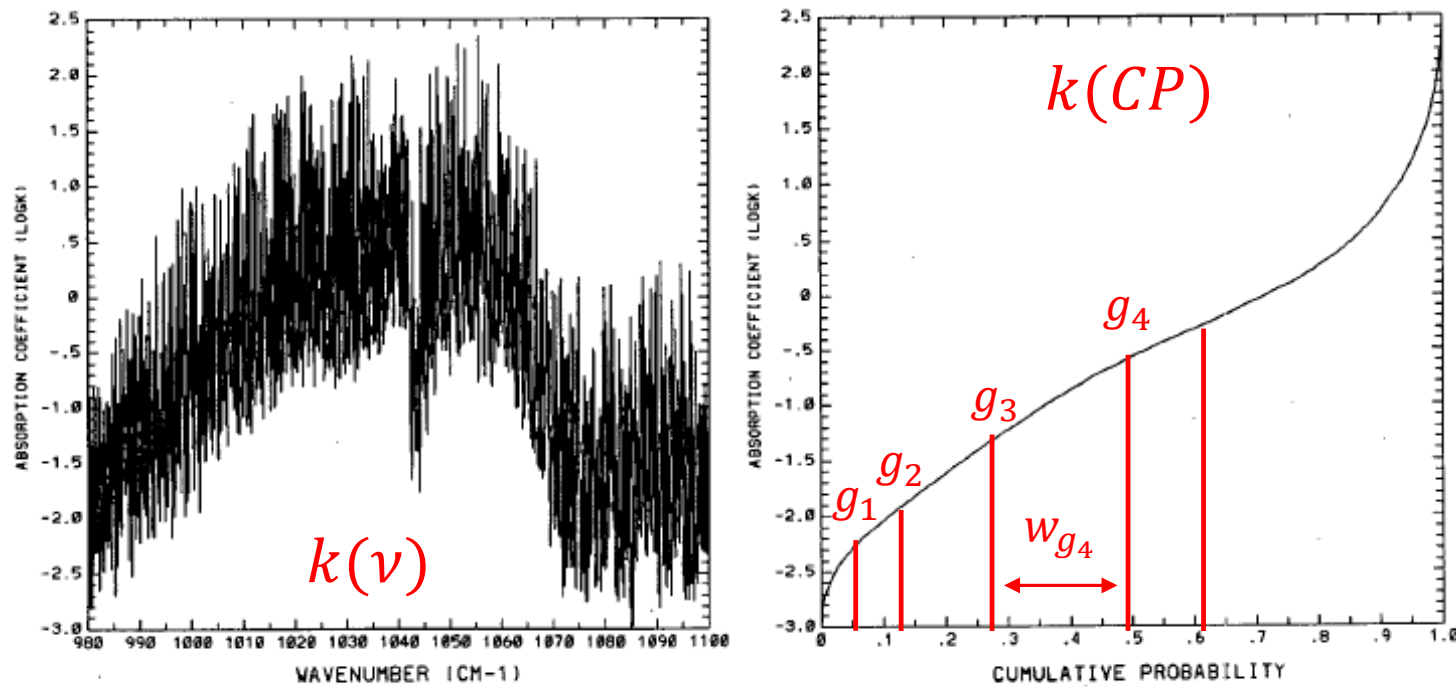


FIG. 1. Absorption coefficient  $k$  in (cm atm)<sup>-1</sup> as a function of (a) wavenumber and (b) cumulative probability for the  $O_3$  9.6- $\mu$ m band for a pressure of 25 mb and a temperature of 220 K.



# Exponential Sum Fitting Technique - ESFT

- One simple application of KDM is calculating the transmission function for a wide spectral interval and to fit it to series of exponentials as function of path length  $u$  :

$$T_{\bar{\lambda}}(u) \approx \frac{1}{\Delta\lambda} \int_{\Delta\lambda} e^{-k_{\lambda} u} d\lambda \approx \int_0^1 e^{-k(g)u} dg \approx \sum_g w_g e^{-k_g u}, \quad \sum_g w_g = 1$$

- The total flux is a some of pseudo-monochromatic fluxes for all three gasses, in all spectral intervals  $b$  and for all  $g$ -points

$$F(x, y, z, t) \approx \sum_b w_b \sum_{g_{b,1}} \sum_{g_{b,2}} \sum_{g_{b,3}} w_{g,b,1} w_{g,b,2} w_{g,b,3} F(\delta_0 + \delta_{g,b,1} + \delta_{g,b,2} + \delta_{g,b,3})$$

$\delta_0$  - optical thickness of gray constituents only (clouds, aerosols)

- But still - computationally very expensive!

# Fast ESFT - FESFT

- First guess would be neglecting overlapping absorption bands of different gases in other words considering only the dominant gas in each band – causes systematic errors we cannot afford
- **FESFT** - Calculate each gas + gray constituents separately and then combine:

$$\bar{T}_{1,\lambda} = \frac{F^1}{F^0} = \frac{\sum_g w_g F(\delta_0 + \delta_1)}{F^0}$$

$$F \approx \prod_{i=1}^{N_{gas}} \bar{T}_{i,\lambda} F^0$$

- CPU run time gain is factor of ~3
- Reasonable accuracy
- COSMO's default scheme

Still not fast enough →  
compromise on spatial/temporal resolution

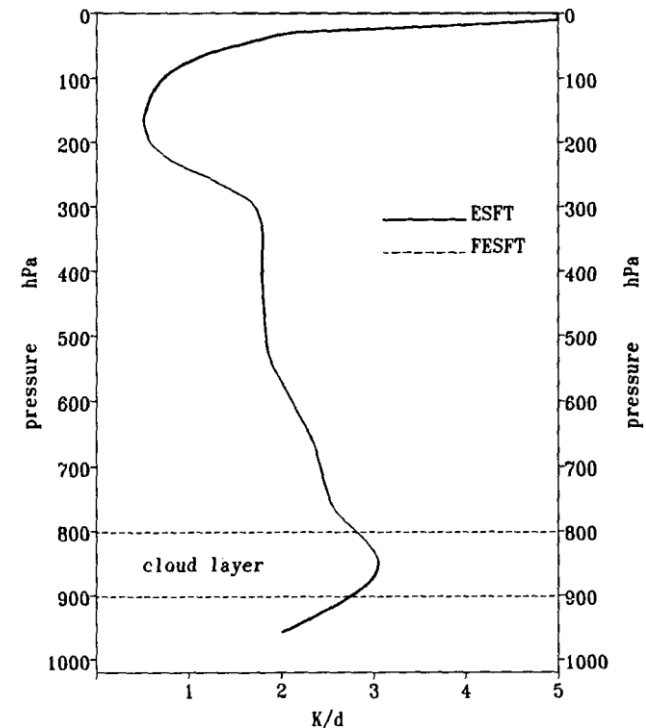
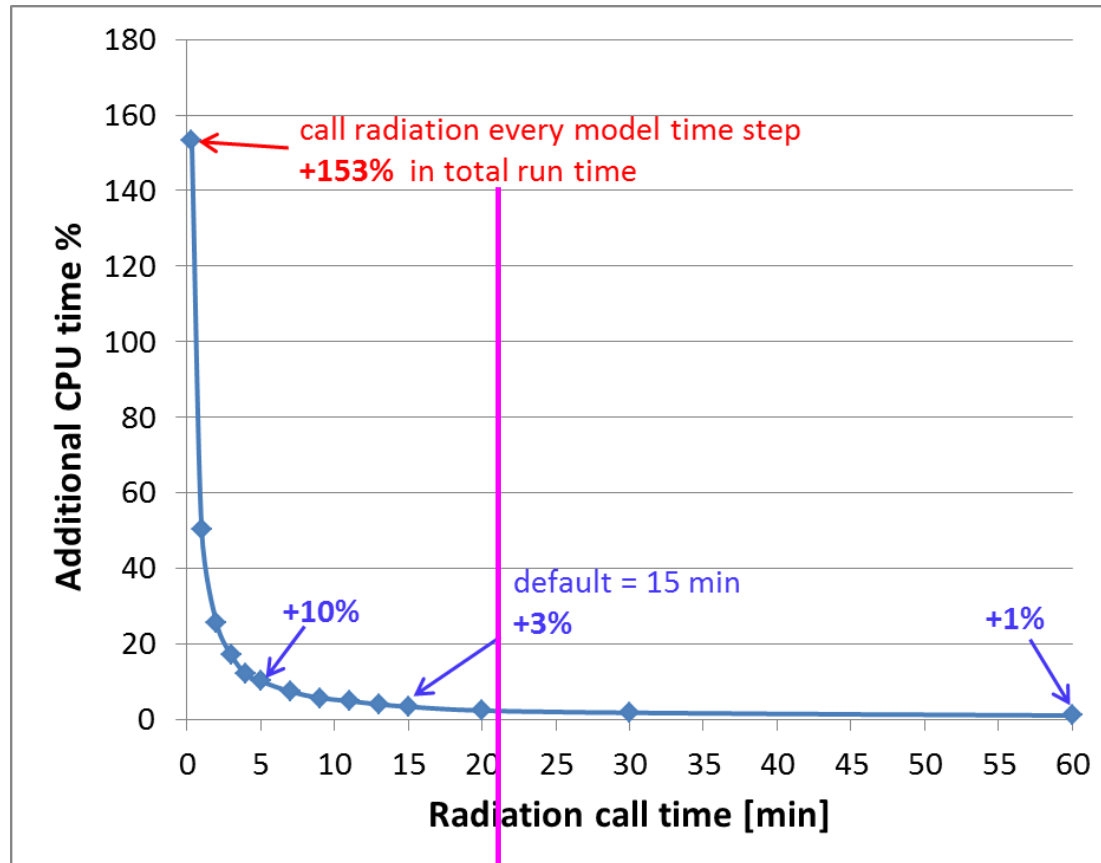


FIG. 1. Comparison of solar heating rates in a midlatitude summer atmosphere using the exponential sum-fitting technique in its original (ESFT) and its approximate fast (FESFT) version. A cloud with  $10 \text{ g m}^{-2}$  liquid water content is located between the 1000- and 2000-m heights. Solar zenith angle is  $30^\circ$  and a surface albedo of 0.20 is assumed.

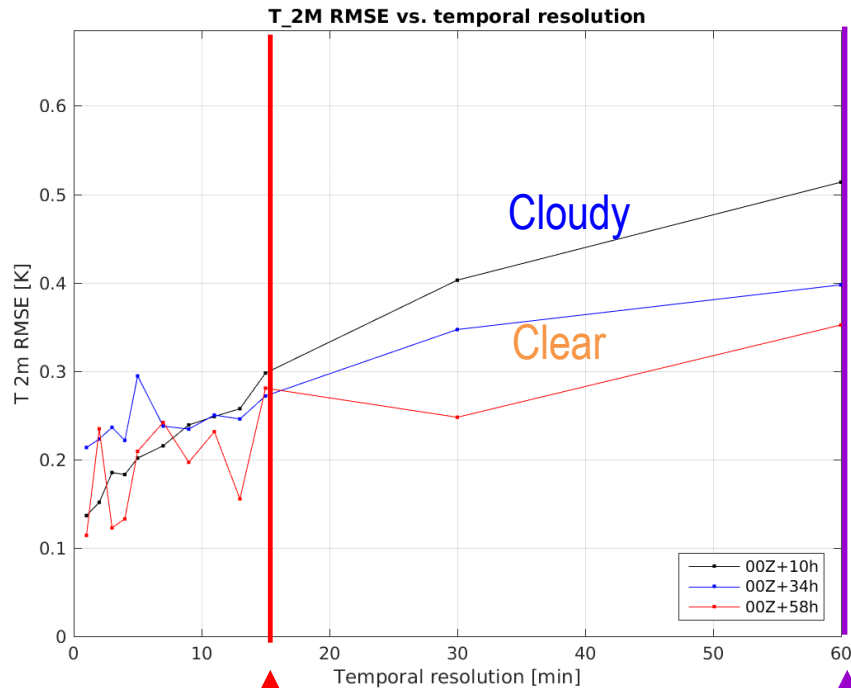
# Radiation Temporal Resolution



COSMO 2.8  
operational setup  
using FESFT

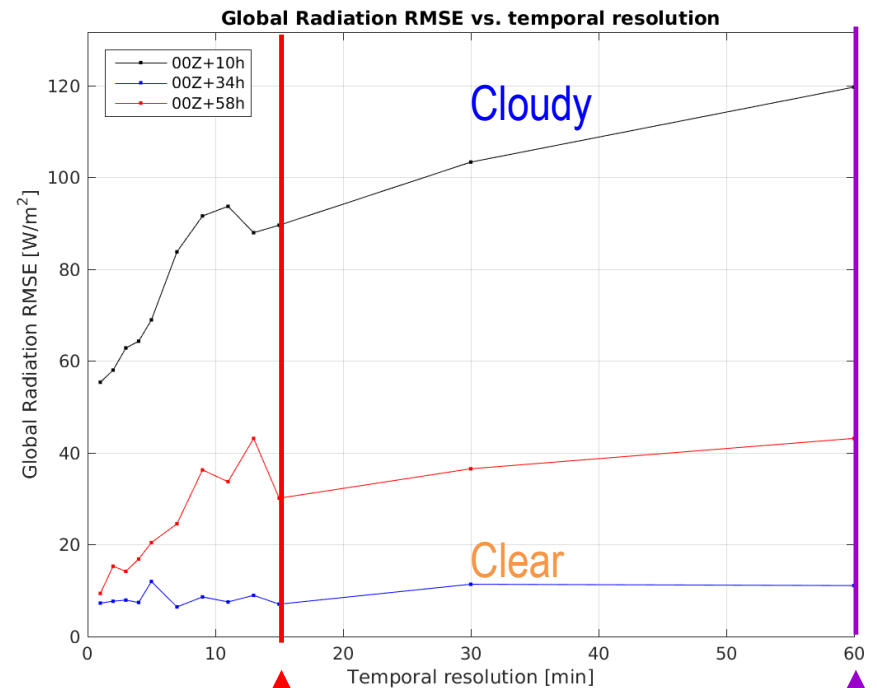
COSMO\_radiation = 1.5 X COSMO model

# Radiation Temporal Resolution



15 min

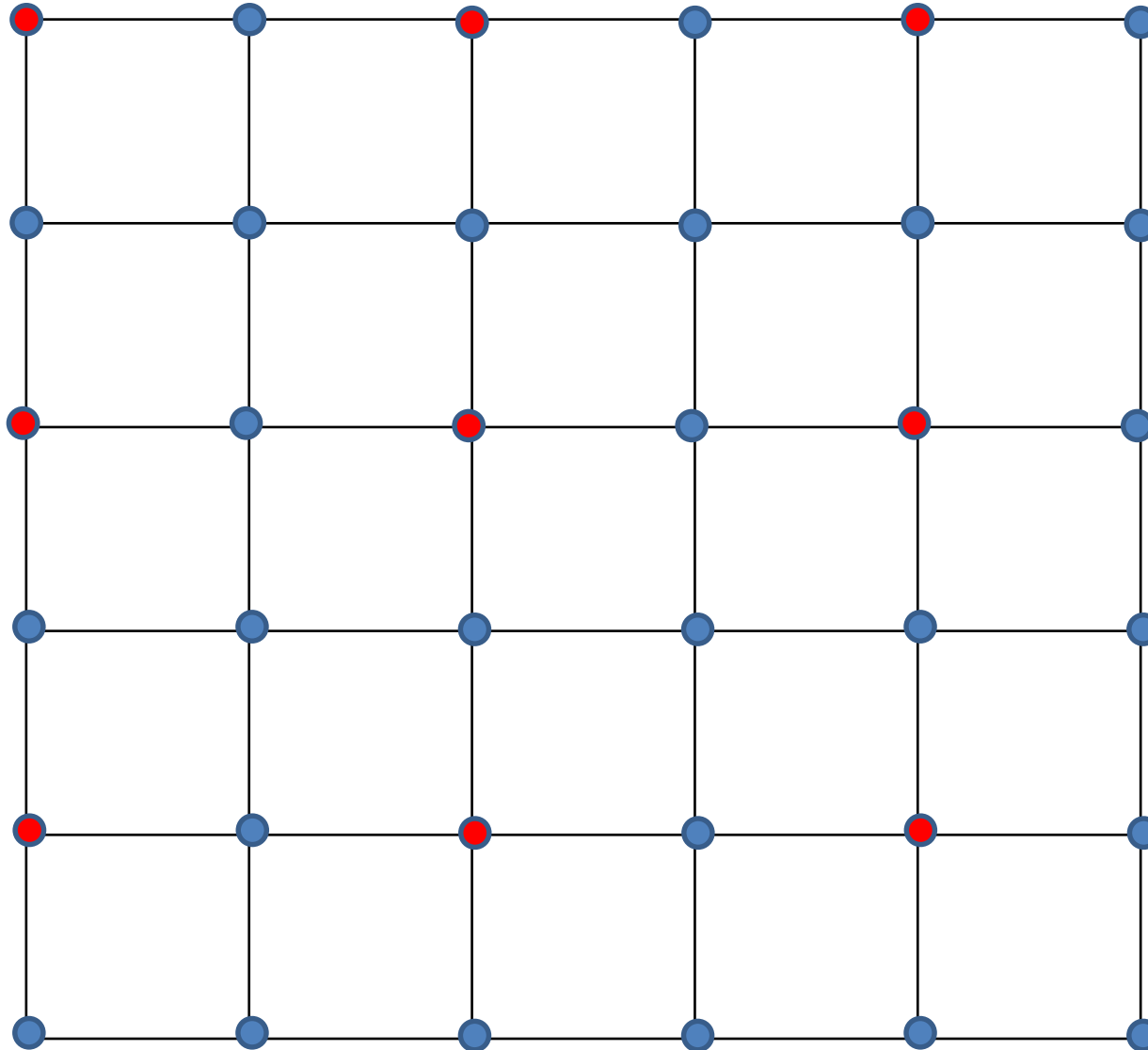
60 min



15 min

60 min

# Radiation Spatial Resolution



nradcoarse = 2

# Monte-Carlo Method

Manhattan Project



Stanislaw Ulam



John von Neumann



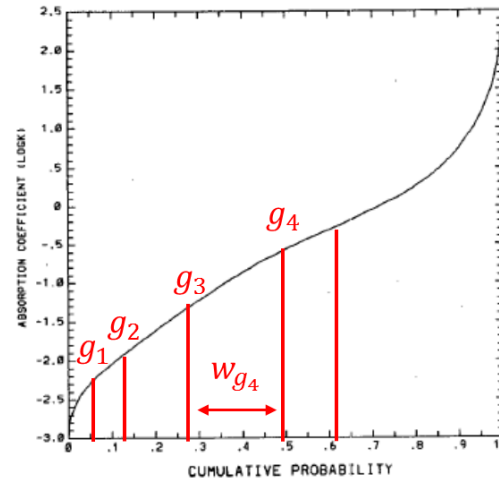


# Monte-Carlo Spectral Integration - MCSI

J. Adv. Model. Earth Syst., Vol. 1, Art. #1, 9 pp.

## Monte Carlo Spectral Integration: a Consistent Approximation for Radiative Transfer in Large Eddy Simulations

Robert Pincus<sup>1</sup> and Bjorn Stevens<sup>2</sup>



$$P(g) = \frac{1}{w_{g(b)}}$$

Back to **ESFT** but instead of doing this every 15 min (45 time steps):

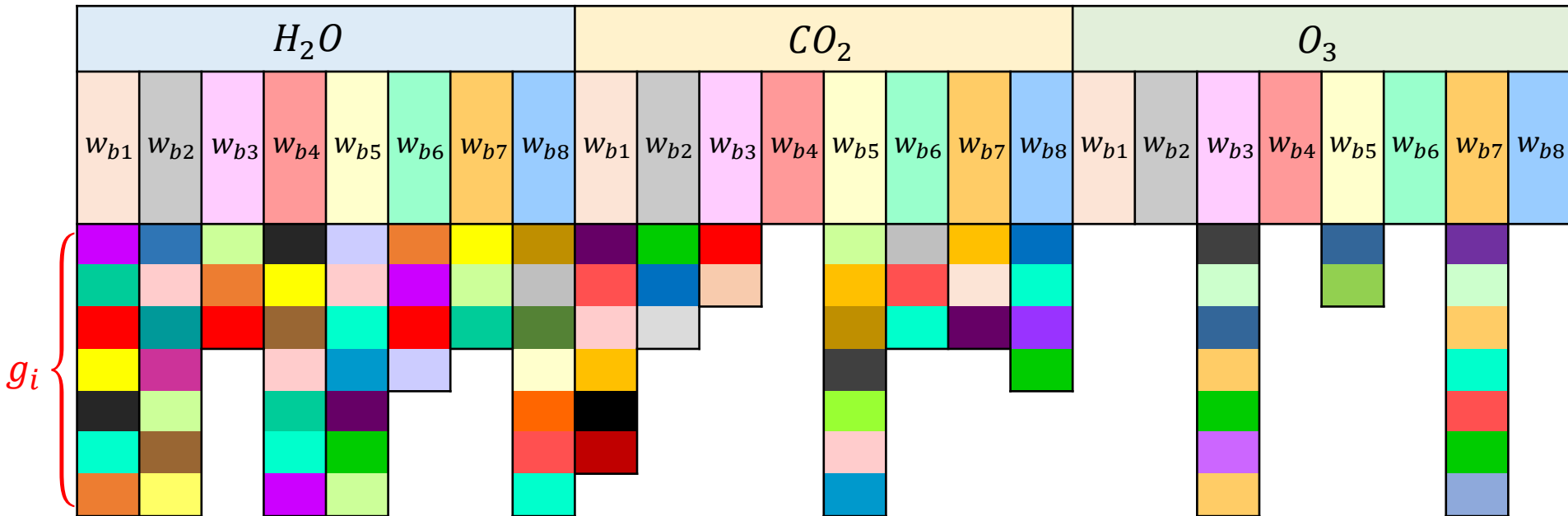
$$F(x, y, z, t) \approx \sum_b w_b \sum_{g_{b,1}} \sum_{g_{b,2}} \sum_{g_{b,3}} w_{g,b,1} w_{g,b,2} w_{g,b,3} F(\delta_0 + \delta_{g,b,1} + \delta_{g,b,2} + \delta_{g,b,3})$$

Pick only one *g* point according to its probability weight for each gas & band more frequently (i.e. every time step):

$$F(x, y, z, t) \approx \sum_b w_b F(\delta_0 + \delta_{g',b,1} + \delta_{g',b,2} + \delta_{g',b,3})$$

Locally temporal big errors that averages fast to an accurate solution!

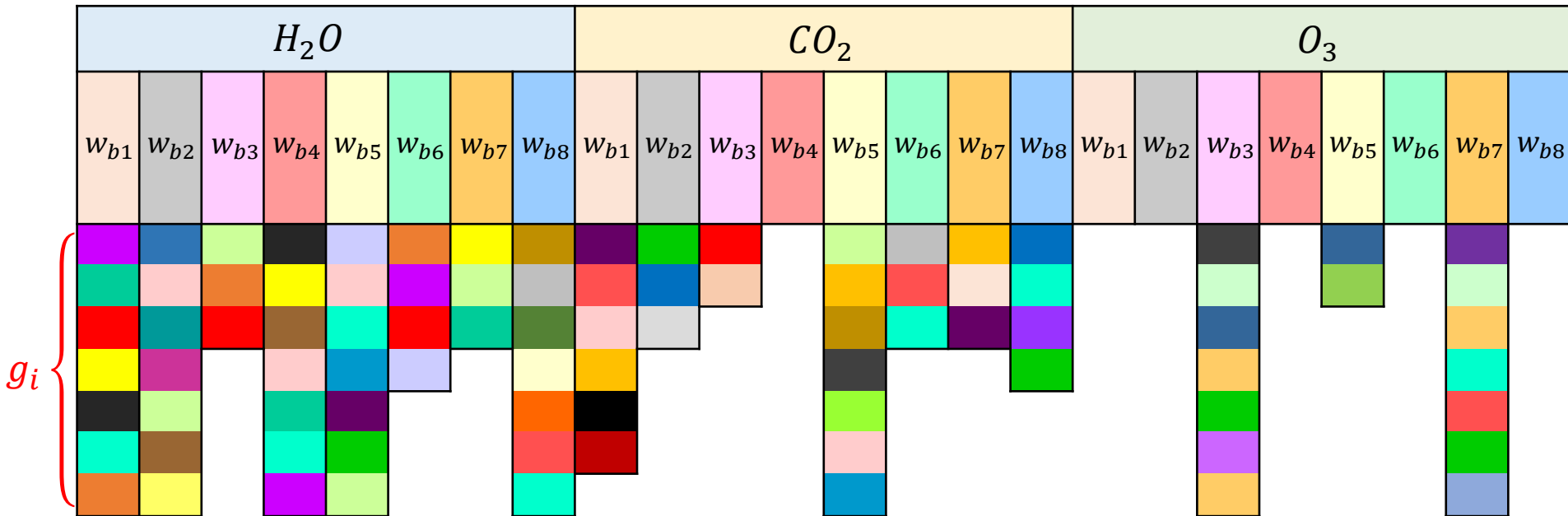
# COSMO ESFT Diagram



Example: for spectral interval  $b=7$  we have  $3 \times 3 \times 7 = 63$  calls `inv_th/inv_so` subroutines which calculate the fluxes

→ Total of **301** calls to `inv_th/inv_so` subroutines

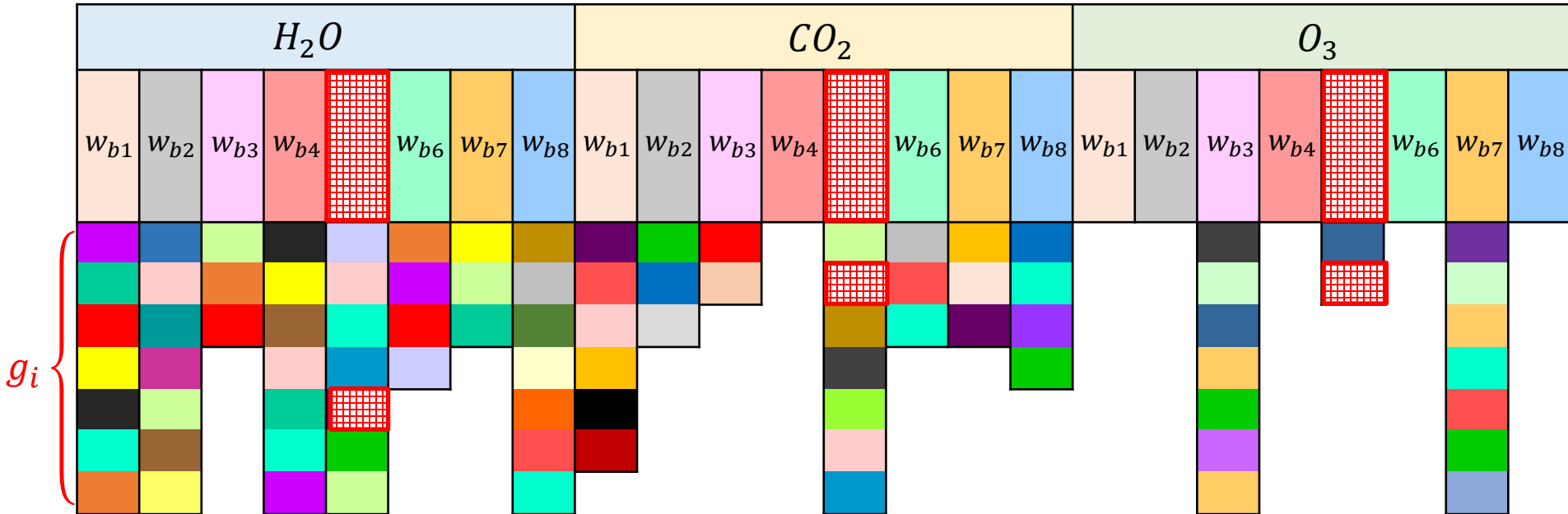
# COSMO FESFT Diagram



→ Here we calculate each b, g only once (all small boxes)  
total of **87** calls to inv\_th/inv\_so subrutines

$$\text{CPU gain} \approx \frac{\text{calls decreases}}{\text{frequency increase}} = \frac{301/87}{1} = 3.46$$

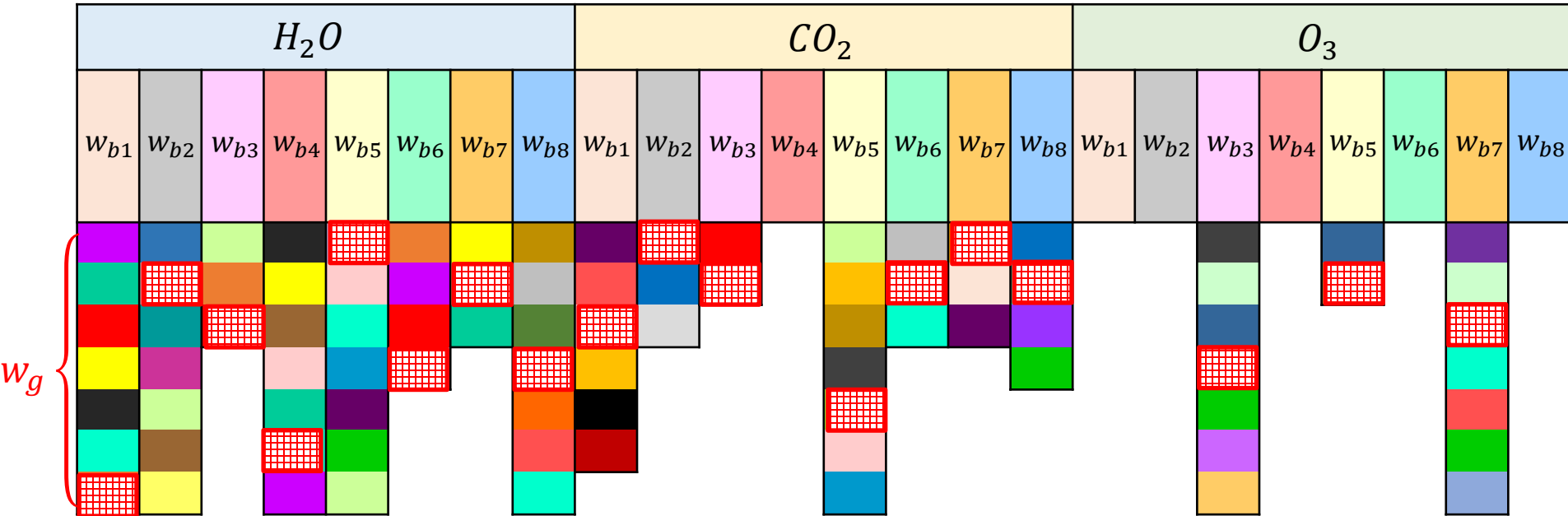
# COSMO MCSI Diagram – Classic Version



→ Only 1 call to inv\_th/inv\_so subroutines  
instead of 301 calls in ESFT

$$\text{CPU gain} \approx \frac{\text{calls decreases}}{\text{frequency increase}} = \frac{301/1}{45} = 6.7$$

# COSOMO MCSI Diagram – Soft Version



→ Only 8 calls to inv\_th/inv\_so subrutines  
instead of 301 calls in ESFT!

$$\text{CPU gain} \approx \frac{\text{calls decreases}}{\text{frequency increase}} = \frac{301/8}{45} = 0.83$$

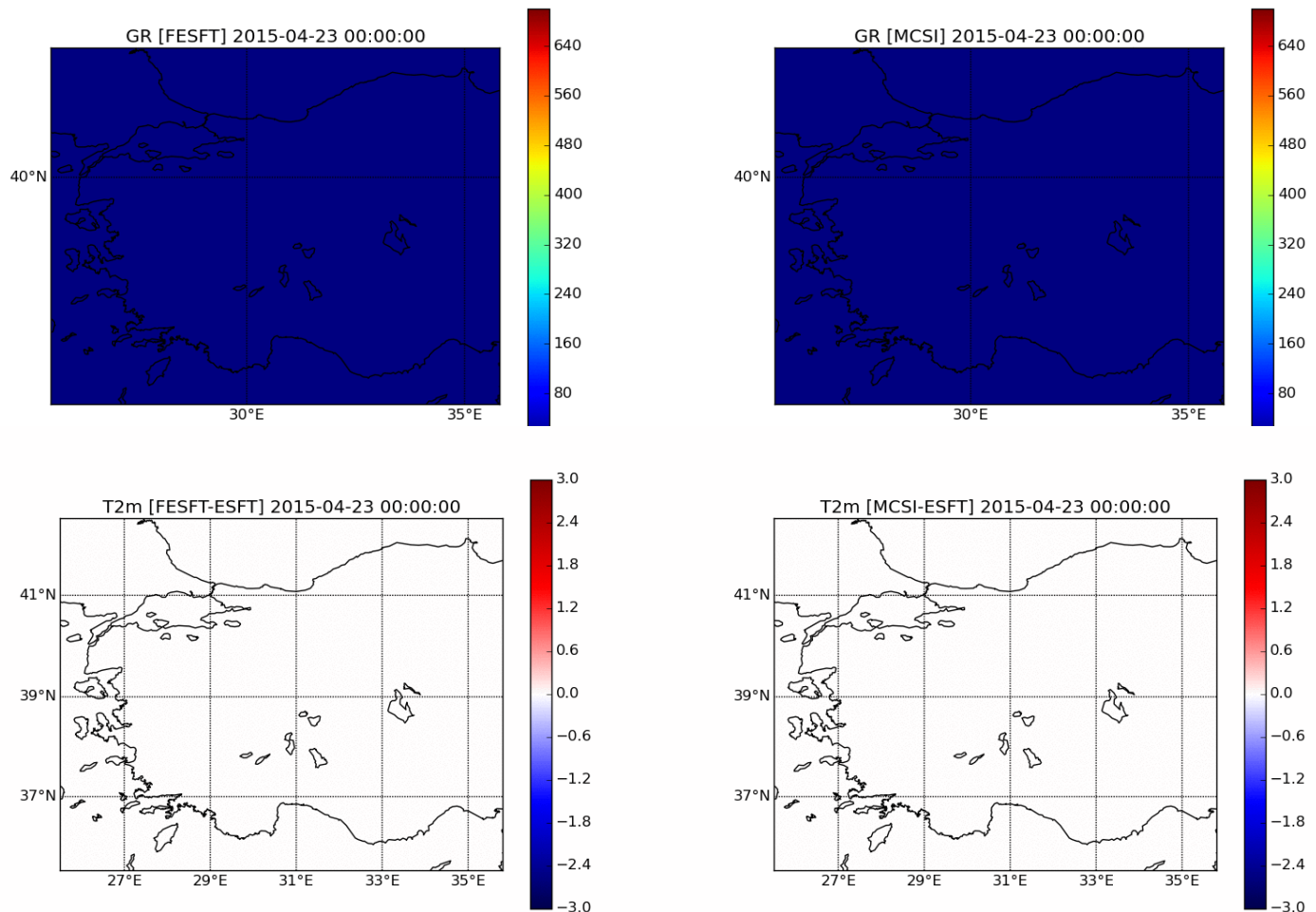
# COSMO Radiation Module

```
MODULE src_radiation
...
SUBROUTINE organize_radiation
...
SUBROUTINE fesft      ! ESFT & FESFT
....
    DO jspec= 1, nspec    ! Spectral loop
        ...
        DO jh2o = 1, ih2o    ! Loop over H2O coefficients
            DO jco2 = 1, ico2    ! Loop over CO2 coefficients
                DO jo3 = 1, io3    ! Loop over O3 coefficients
                    ...
                    CALL    inv_th/so
                    ...
            ...
        ...
    ...
...
...
```

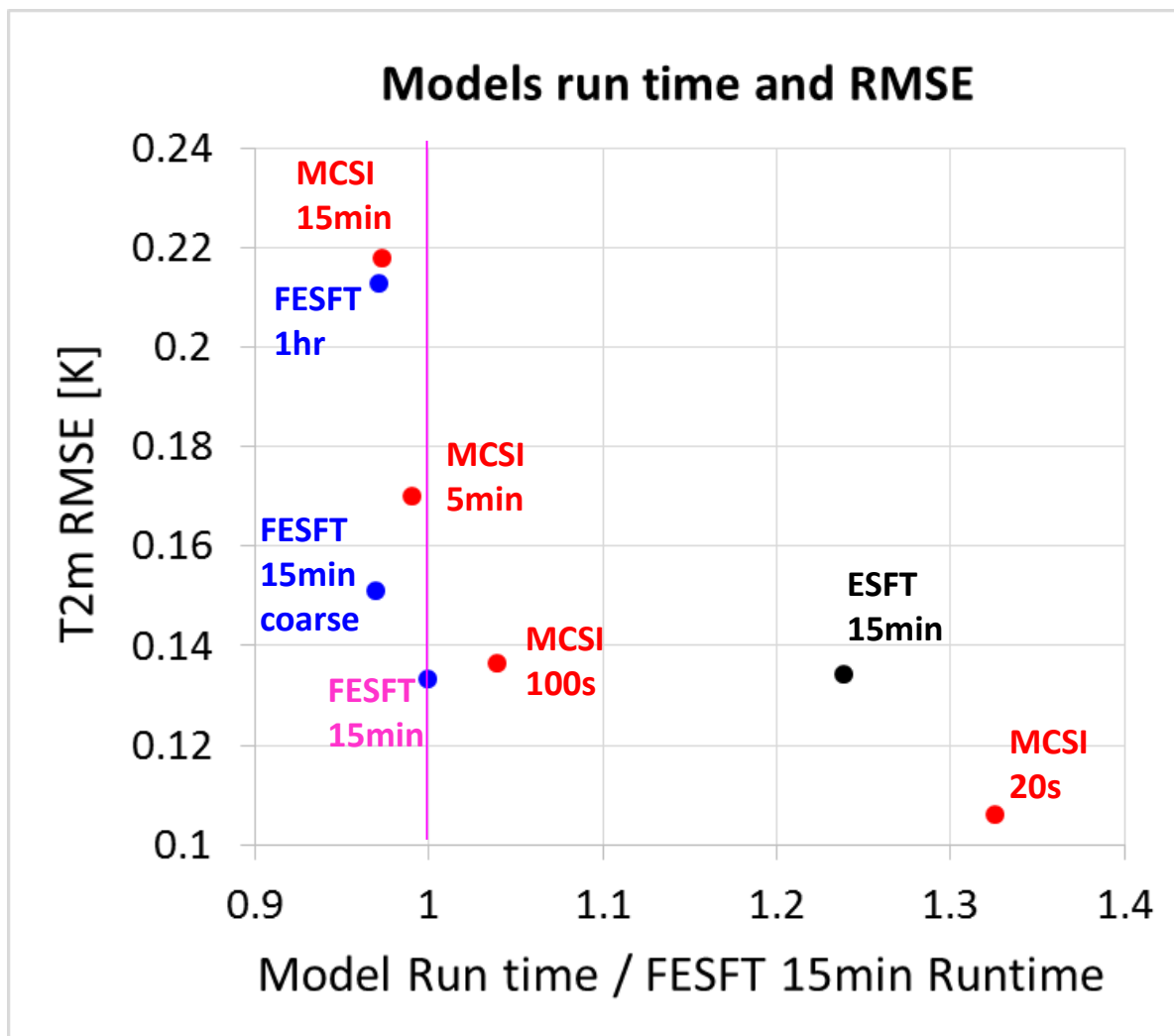


# Run Time & Errors Comparisons

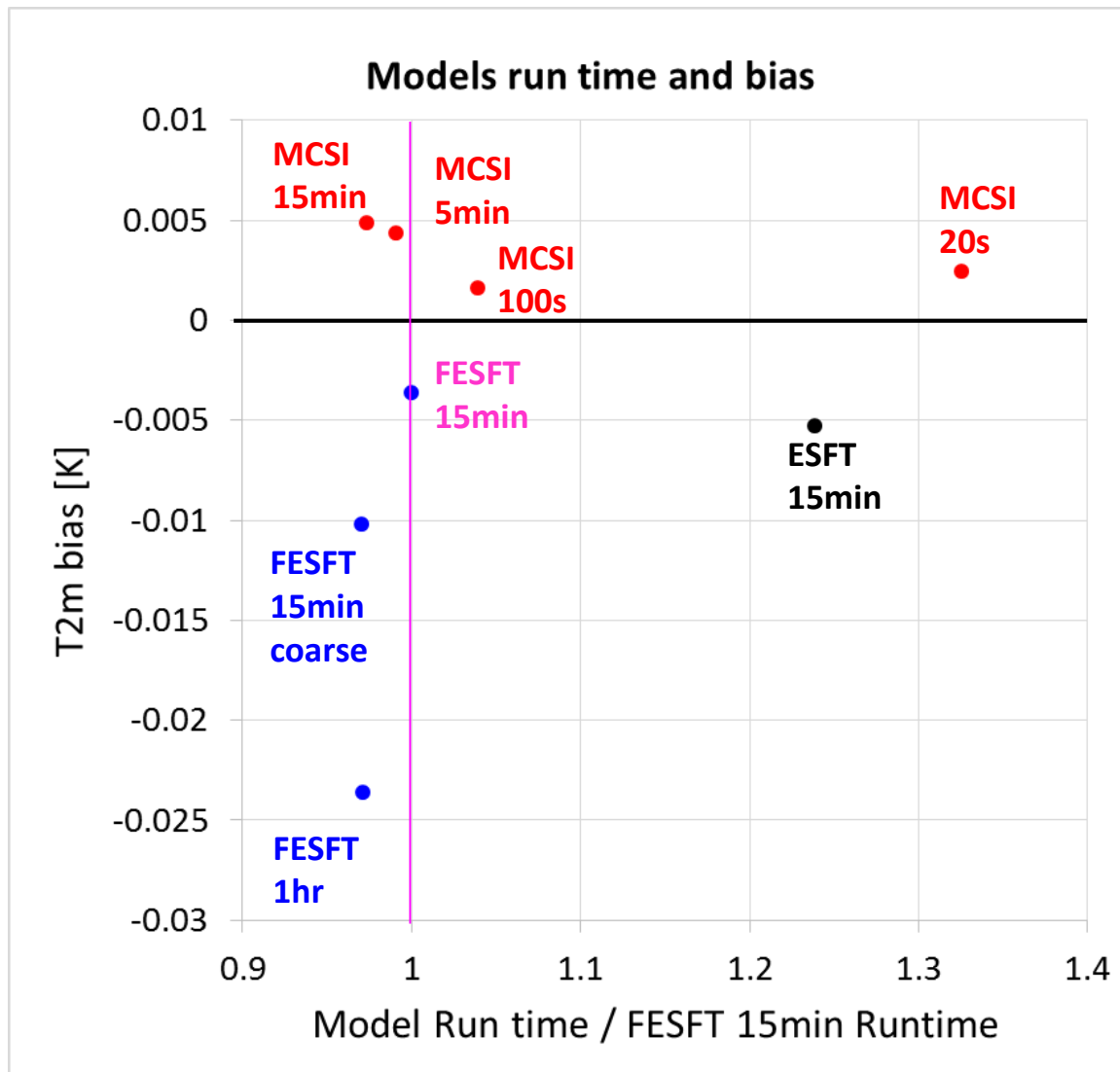
- COSMO-2.8km
- Test case: 23-25/04/2015 – Turkey
- Partial cloudiness + High wind speeds
- Stand alone computer 1-node 4 CPUs



# Run Time & Errors Comparisons

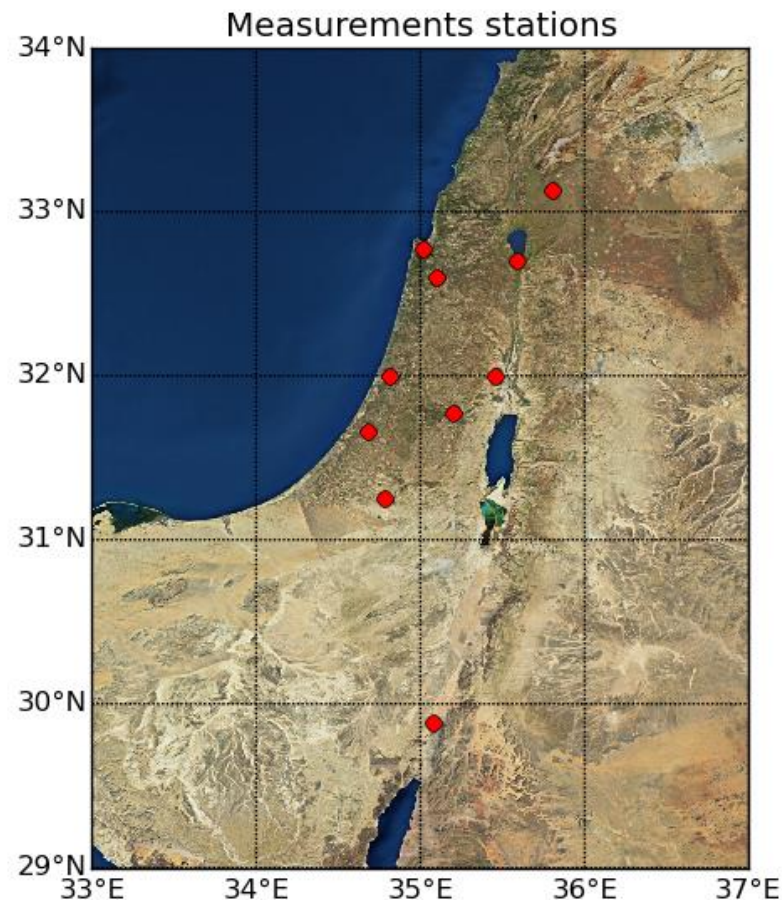
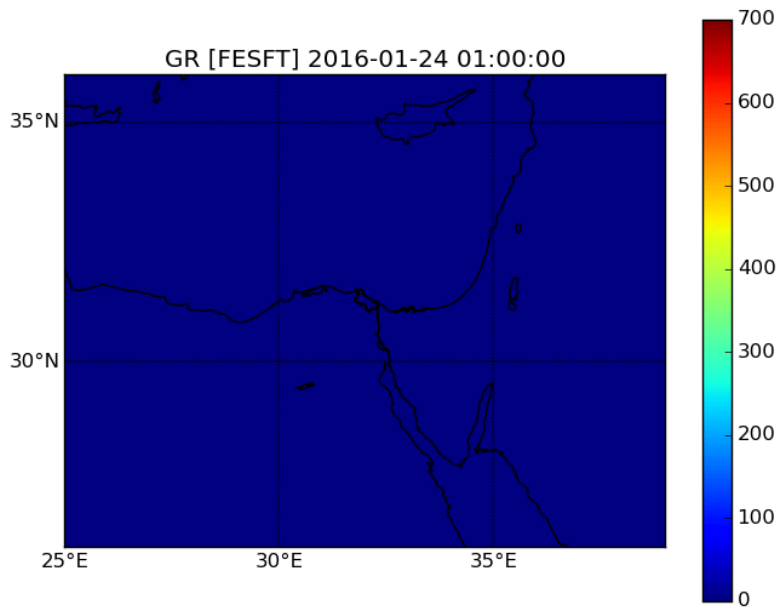


# Run Time & Errors Comparisons



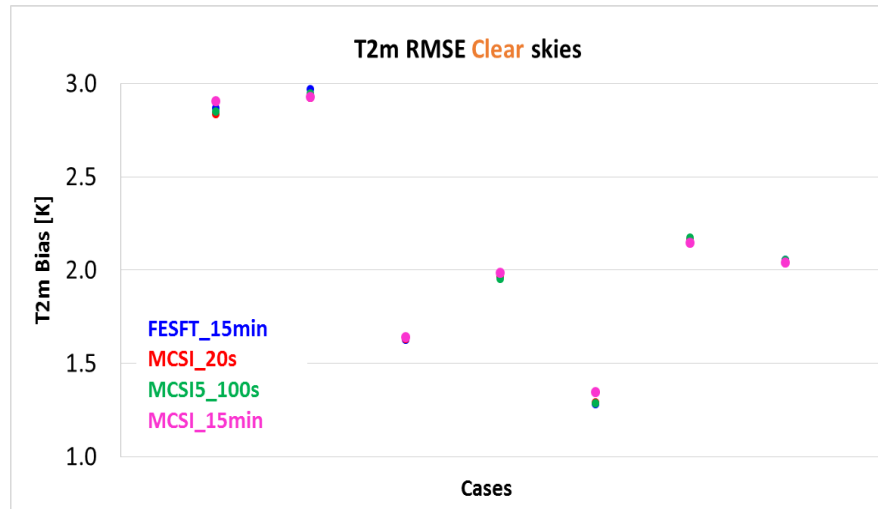
# Testing MCSI Scheme vs. Ground Based Measurements

- **29 test cases** in different weather situations, lead time of 30h/42h
- **10** measurement stations – T2m validation
- Compare 3 models:
  - **FESFT** – 15 min / 45 steps
  - **MCSI** – 20 s / 1 step
  - **MCSI** – 100 s / 5 step
  - **MCSI** – 15 min / 45 steps



# Testing vs. Ground Based Measurements

## Clear skies



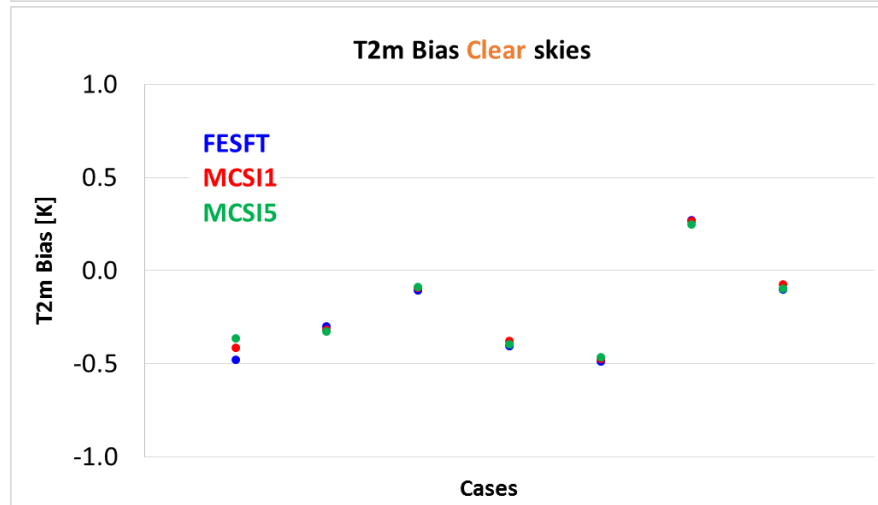
T2m RMSE:

FESFT\_15min 2.13 [K]

MCSI\_20s 2.12 [K]

MCSI\_100s 2.13 [K]

MCSI\_15min 2.14 [K]



T2m bias:

FESFT\_15min -0.23 [K]

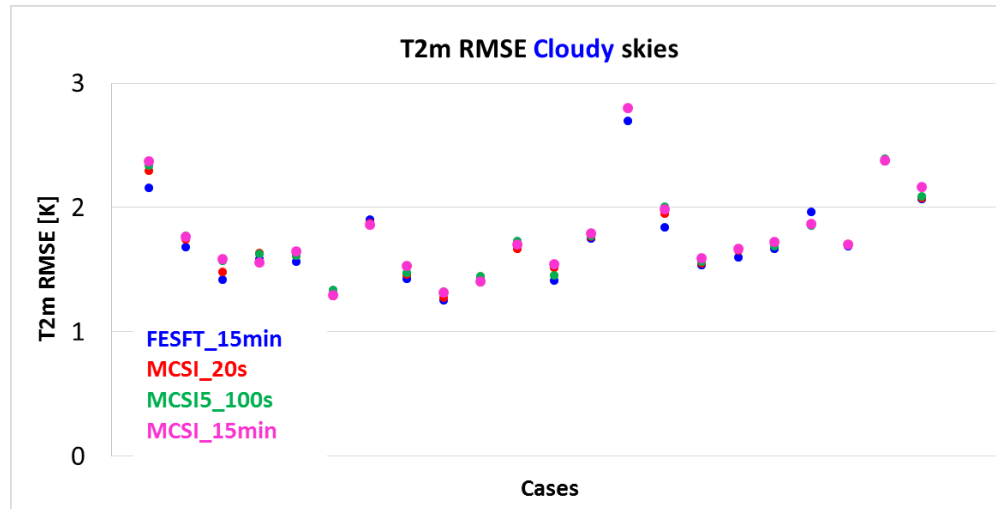
MCSI\_20s -0.21 [K]

MCSI\_100s -0.21 [K]

MCSI\_15min -0.21 [K]

# Testing vs. Ground Based Measurements

## Cloudy skies



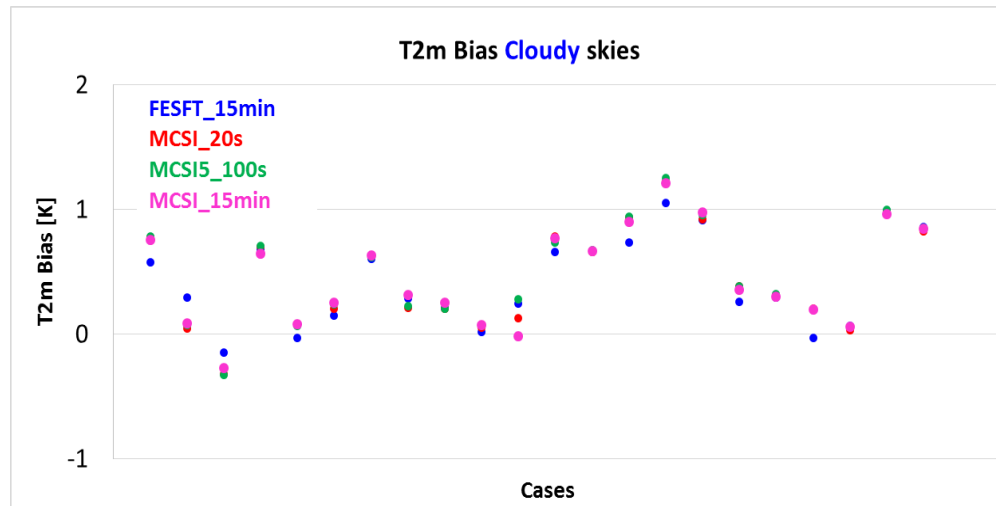
T2m RMSE:

FESFT\_15min 1.73 [K]

MCSI\_20s 1.76 [K]

MCSI\_100s 1.78 [K]

MCSI\_15min 1.79 [K]



T2m bias:

FESFT\_15min 0.42 [K]

MCSI\_20s 0.45 [K]

MCSI\_100s 0.47 [K]

MCSI\_15min 0.46 [K]



# Summary

- “Full” radiation scheme calculations is impractical in NWP applications
- We can compromise on the : **spatial**, **temporal** or **spectral** resolutions
- Each has it own advantages and disadvantages. The MCSI greatest strength is that the “dilution” of computations is wise and based on statistical reasoning
- The MCSI is now implemented in COSMO (**itype\_mcsi = 1**) gives a reasonable and comparable results in both CPU and performance to the default FESFT scheme
- MCSI did not show a significant advantage to the FESFT which deserves a change in the default scheme choice
- Nevertheless, the tests shown here were done on 2.8 km / 20 seconds model resolution. It is possible that MCSI can be preferable when using different model uses (climate, LES) and model resolutions.

thank you

tusind tak  
謝謝 dakujem vám  
ありがとう  
ngiyabonga

dziękuję  
merci  
baie dankie  
धन्यवाद molte grazie

suksema  
danke  
gracias  
obrigada  
obrigado  
teşekkür ederim  
شكرا  
tack så mycket

takk  
gràcies  
tānan  
תודה

dank u  
mahalo  
teşekkür edire