
High Spectral Resolution Infrared Radiance Modeling Using Optimal Spectral Sampling (OSS) Method

J.-L. Moncet and G. Uymin

Background

- **Optimal Spectral Sampling (OSS) method is a fast and accurate monochromatic RT modeling technique applicable to a wide range of remote-sensing platforms**
- **Originally developed to support instrument/sounding algorithm trades, validation studies**
 - Same method is applicable across the spectrum (from microwave to UV)
 - Monochromatic radiative transfer makes it directly applicable to multiple scattering, non-positive ILS (interferometers)
 - Physically based approach makes the method robust and ultimately alleviates need for extensive fast model validation
 - Trade-off between accuracy and speed depends on specifics of problem at hand
- **Computationally efficient calculation of radiances and Jacobians makes OSS attractive for operational environment**
- **Initial research-grade version of OSS model:**
 - NPOESS CrIS, CMIS, ATMS EDR algorithms
 - AMSU/MHS processing at AER
 - NAST-I (W. Smith)
- **Parallel R&D effort (under Navy SBIR) led to improved training and new faster and more accurate RT model**
- **Used for AIRS instrument**

ESFT and k -distribution methods

Correlated- k assumption breaks down for single absorber in the presence of lines of different strengths or non-regularly spaced lines

No satisfactory treatment of gas mixture when relative abundance of individual species changes with altitude

- The exponential sum fitting (ESFT) and k -distribution methods approximate band transmittances in homogenous atmospheres as,

$$\bar{\tau}(u) = \sum_{i=1}^N w_i e^{-k_i u}$$

- Weights w_i can be interpreted in terms of the probability distribution of the absorption coefficient over the spectral interval

$$w_i = \Delta g_i = \int_{k'_{i-1}}^{k'_i} p(k) dk \quad (k'_{i-1} < k'_i)$$

and k_i is a representative k -value for the interval $[k'_{i-1}, k'_i]$

- Problems with extending ESFT or k -distribution concept to inhomogeneous atmospheres lies in difficulty of ensuring physical consistency between the k -values in each layers
 - Extension of the k -distribution method to non-homogeneous atmospheres is based on observation that minima and maxima of absorption in different layers coincide spectrally
 - Correlated- k method vertically integrates RT equation in g -space
 - Equivalent to assuming a correspondence between k 's of same ranking in different layers

OSS solution for inhomogeneous atmospheres

- Proper treatment of overlapping absorbers requires accurate characterization of the multivariate probability distribution of absorption coefficients for all layers and molecules
- High dimensionality of the problem makes it impractical to attempt to solve directly for the k 's without use of appropriate constraints
- **OSS solution:**
 - Reduce the problem to a one-dimensional frequency search
 - Impose the constraint that the k 's correspond to actual values of absorption coefficient for all molecules and layers at the selected frequencies

OSS parameters generation

- **Parameter generation starts from a set of M uniformly spaced monochromatic transmittances (or radiances)**
 - **Compute with a line-by-line model (currently LBLRTM)**
 - **Use a globally representative ensemble (S) of atmospheres**
- **Search for the smallest subset of frequencies (nodes) and associated weights**

$$\left\{ (v_i, w_i) \quad i = 1, \dots, N \right\}$$

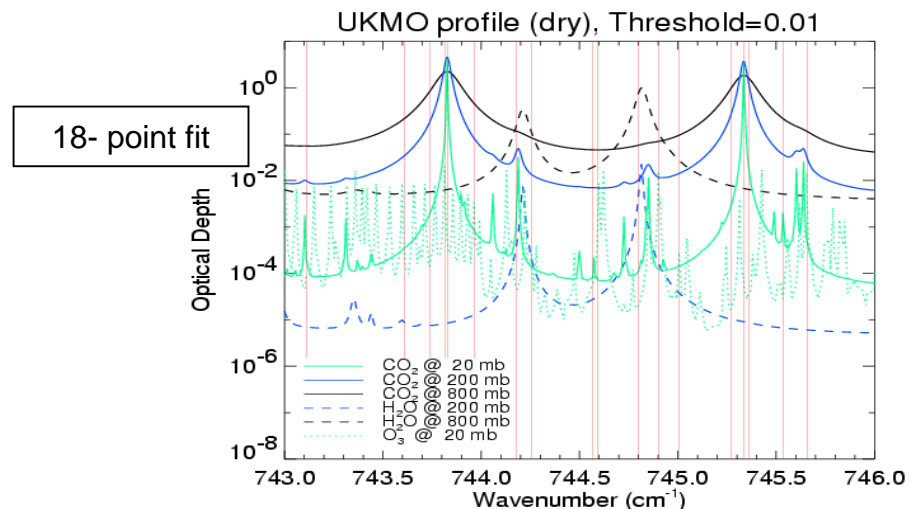
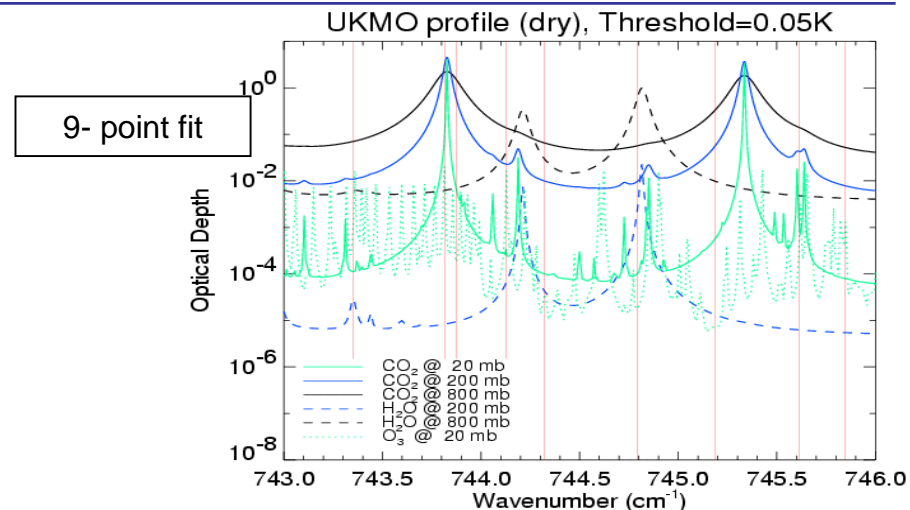
for which the *rms* error computed over the ensemble S

$$\mathcal{E}_N(p_l) = \left[\sum_s \left(\bar{\tau}^s(p_l) - \sum_{i=1}^N w_i \tau_{v_i}^s(p_l) \right)^2 \right]^{\frac{1}{2}}$$

is less than a prescribed tolerance for all levels

OSS Search Procedure

- Procedure consists of starting with number of nodes $N=1$ and searching for the spectral location ν_1 that produces the smallest error among the M possible locations
- N is then incremented by one and search for ν_N proceed in the same fashion
- Procedure stop when prescribed error tolerance for training set is reached
- For each trial combination, weights are obtained by linear regression
- Details of search method are provided in companion poster



Radiance training

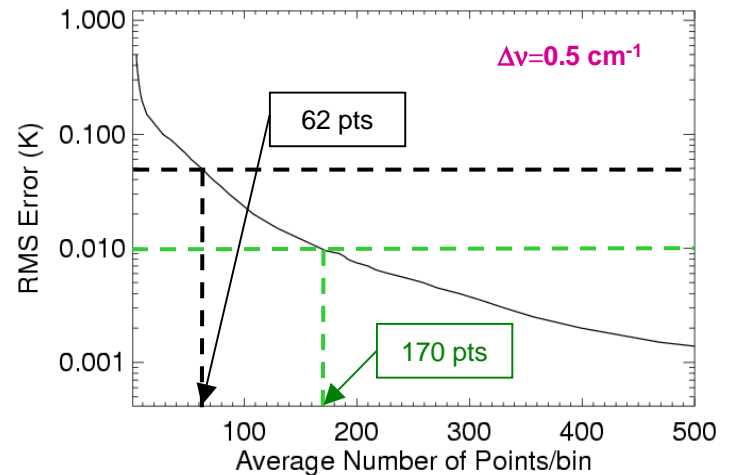
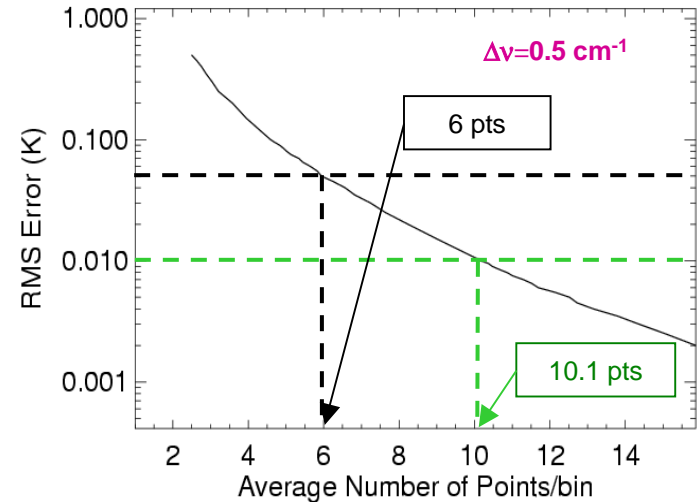
- For infrared remote sensing, fit is done in radiances (or brightness temperatures), which naturally emphasizes levels near the peak of channel weighting function

$$\mathcal{E}_N = \left[\sum_s \left(r^s - \sum_{i=1}^N w_i r_{v_i}^s \right)^2 \right]^{\frac{1}{2}}$$

- Radiance training takes into account functions that vary slowly across channel passband (Planck function, surface emissivity/reflectivity,...etc)
- Set of training scenes includes appropriate variability
 - Viewing angle
 - Surface emissivity and reflectivity
 - Observer altitude
 - Solar angles
- Stratification (in viewing angle, surface emissivity,...etc) may be used to ensure uniform level of accuracy (threshold *rms*) across the range of conditions

OSS vs. Frequency (or Radiance) Sampling Method

- Comparison of number of points used by OSS and RSM provide some quantitative assessment of ability of OSS approach to exploit spectral redundancies
- RSM treats radiance as random variable and relies on fact that estimates of mean improves as number of samples increases (no weighting)
- For a same level of accuracy, reduction in number of samples with OSS, in 700-750 cm^{-1} region, is greater than 90% (in this example line-by-line calculations are sampled on 10^{-4} cm^{-1} grid, i.e. 5000 pts per bin)



OSS Radiative Transfer Model (1)

- Radiative transfer is performed monochromatically (one node at a time) from pre-computed absorption coefficients for the fixed and variable constituents

$$R \cong \sum_{l=1}^N (T_{l-1}^{\uparrow} - T_l^{\uparrow}) B(\Theta_l) + \varepsilon_s T_N B_s + (1 - \varepsilon_s) T_N \sum_{l=1}^N (T_l^{\downarrow} - T_{l-1}^{\downarrow}) B(\Theta_l)$$

- Planck function and surface emissivity are evaluated at the precise “node” spectral location (as opposed to using effective values for the channel)
- Several channels may share the same node: computed radiance is added (after appropriate weighting is applied) to partial results for the channels that use that node
- Number of variable molecules varies from node to node
- Molecules currently included are: H₂O, CO₂, O₃, N₂O, CO, CH₄
- Grouping between fixed and variable gases is decided at run time

OSS Radiative Transfer Model (2)

- Total monochromatic optical depth for a layer

$$\tau = \tau_{fix} + \tau_w + \sum_{m=2}^{Nvar} \tau_m$$

- Fixed gases optical depth

$$\tau_{fix} = k_{fix} u_{dry} = k_{fix} \left(\frac{\Delta P}{g} - u_w \right)$$

where k_{fix} is the effective absorption coefficient for the fixed gases mixture

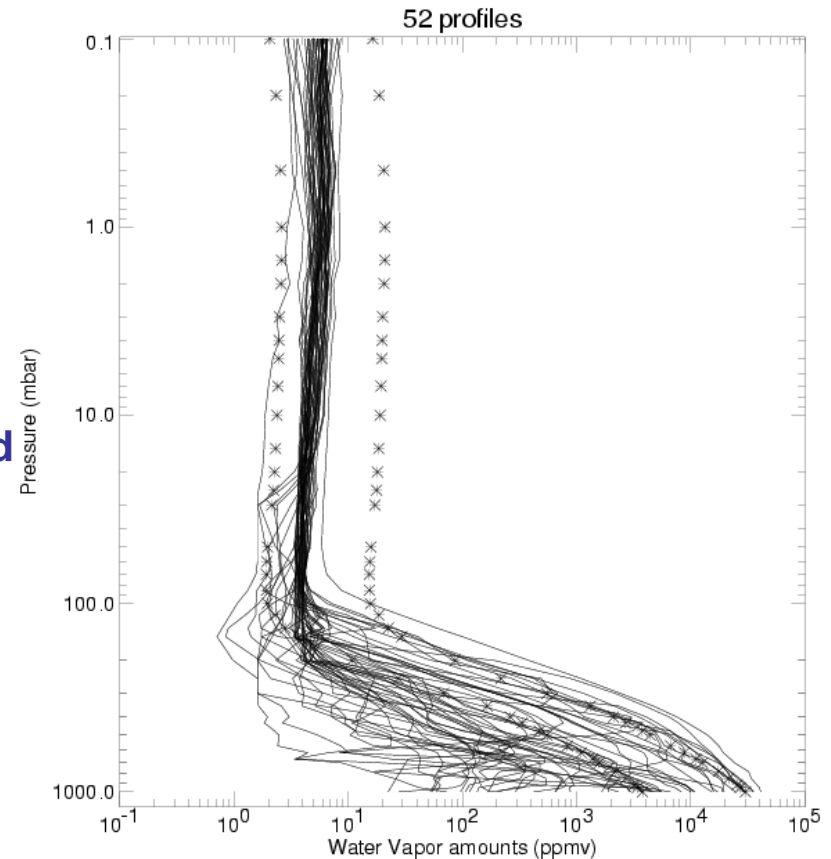
- Dry variable gases

$$\tau_m = k_m u_m$$

- Water vapor self-broadening effect handled by assuming that absorption coefficient depends linearly on specific humidity

$$\tau_w = \left[k'_w + \bar{q} \times \frac{\partial k_w}{\partial \bar{q}} \right] u_w$$

- Approximation does not hold in the near wing (non-linear dependence of absorption of line half-width)



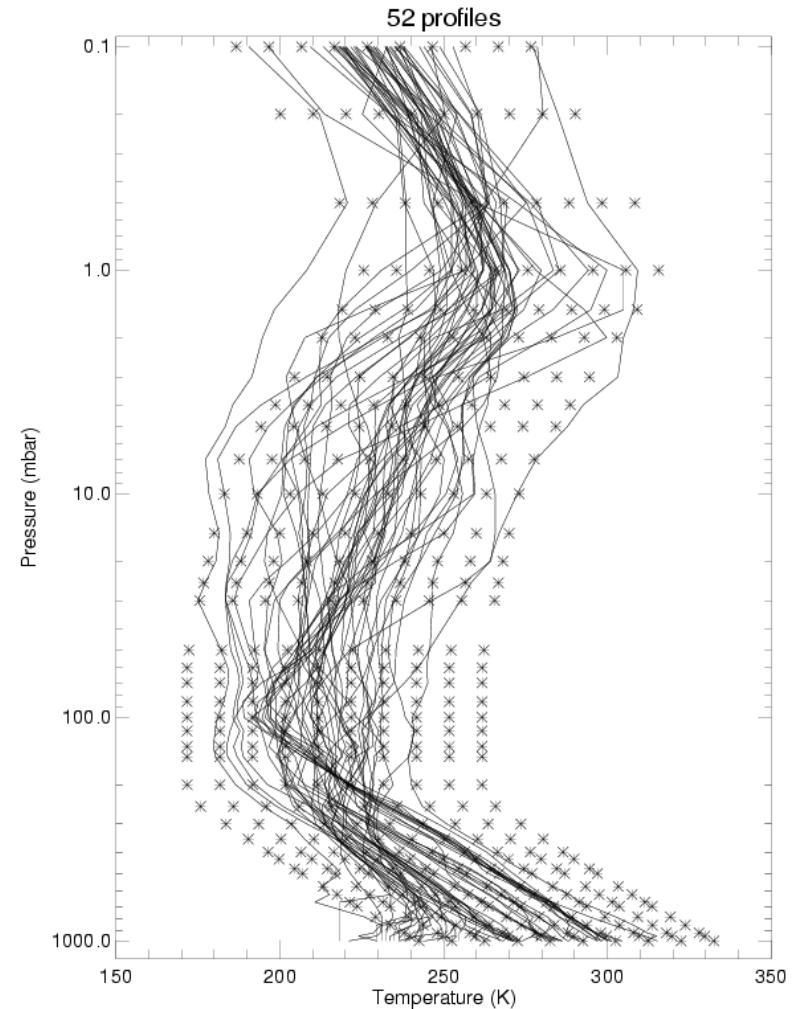
OSS Radiative Transfer Model (3)

- Absorption coefficients are stored in each layer as a function of temperature
- Use quadratic interpolation in temperature
- 3-points Lagrange interpolation provides continuous 1st derivatives

$$k(\theta) = \frac{(\theta - \theta_i)(\theta - \theta_{i+1})}{(\theta_{i-1} - \theta_i)(\theta_{i-1} - \theta_{i+1})} (k_{i-1} - k_{i+1})$$

$$+ \frac{(\theta - \theta_{i-1})(\theta - \theta_{i+1})}{(\theta_i - \theta_{i-1})(\theta_i - \theta_{i+1})} (k_i - k_{i+1}) + k_{i+1}$$

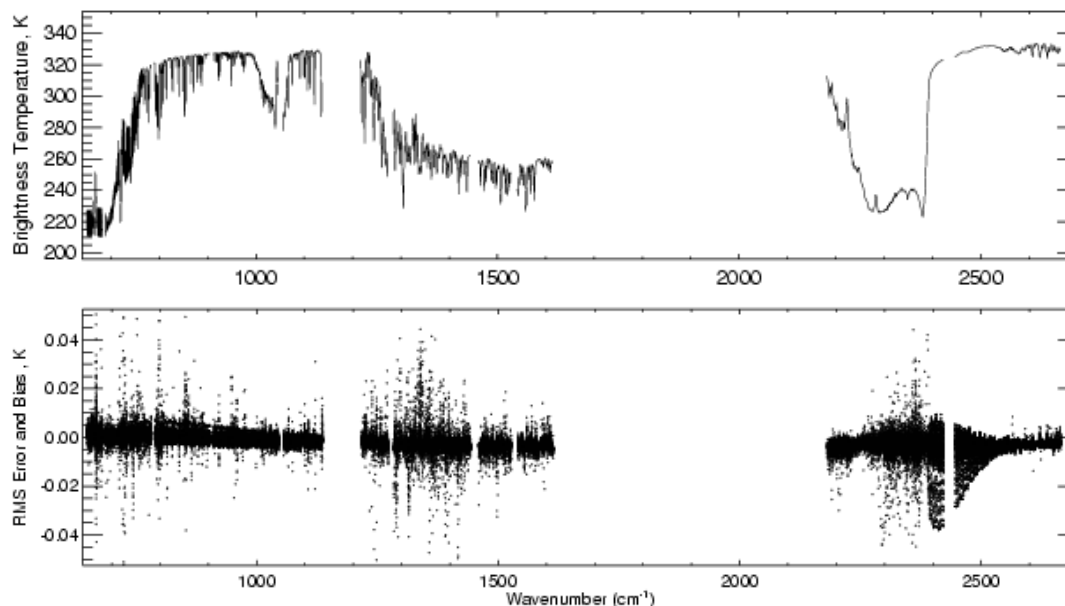
Depend only on layer temperature; computed only once prior to performing the RT calculations



Impact of absorption coefficient errors

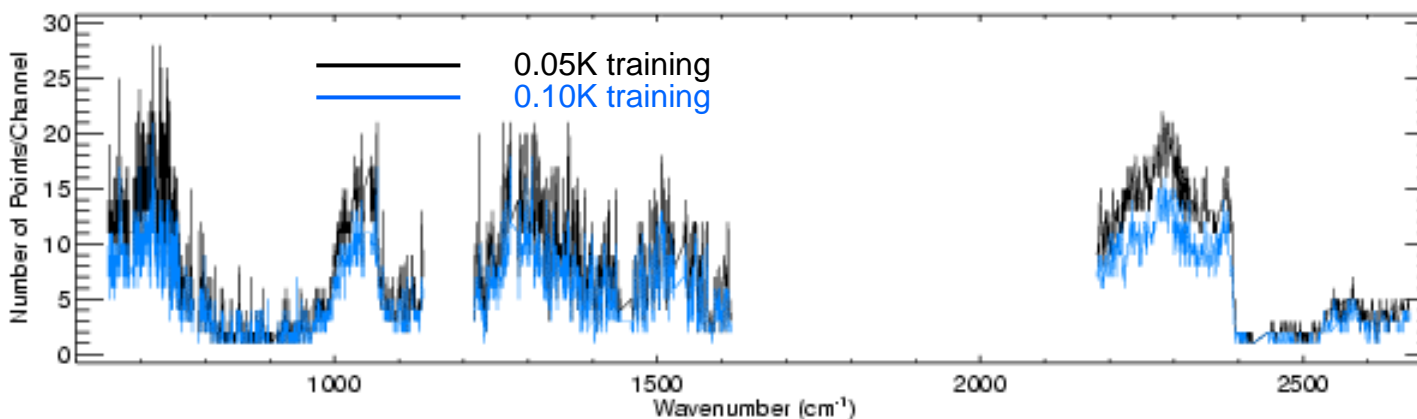
- Impact of errors in treatment of absorption coefficients on computed brightness temperatures is small
- Maximum errors for AIRS reach 0.05 K (*rms* error < ~0.02K)
- Could be reduced further in CO₂ bands and 2400-2500 cm⁻¹ region (N₂ continuum) by extending temperature range of the tables (and by reducing temperature step size - no impact on computational efficiency)
- Not a high priority

Example:
AIRS instrument -
nadir viewing



AIRS OSS model

- Selection accuracy of 0.05K and 0.1 K in brightness temperature
- EIA: 0 to 60 °
- Random surface emissivity in range 0.9-1.
- Variable H₂O and O₃



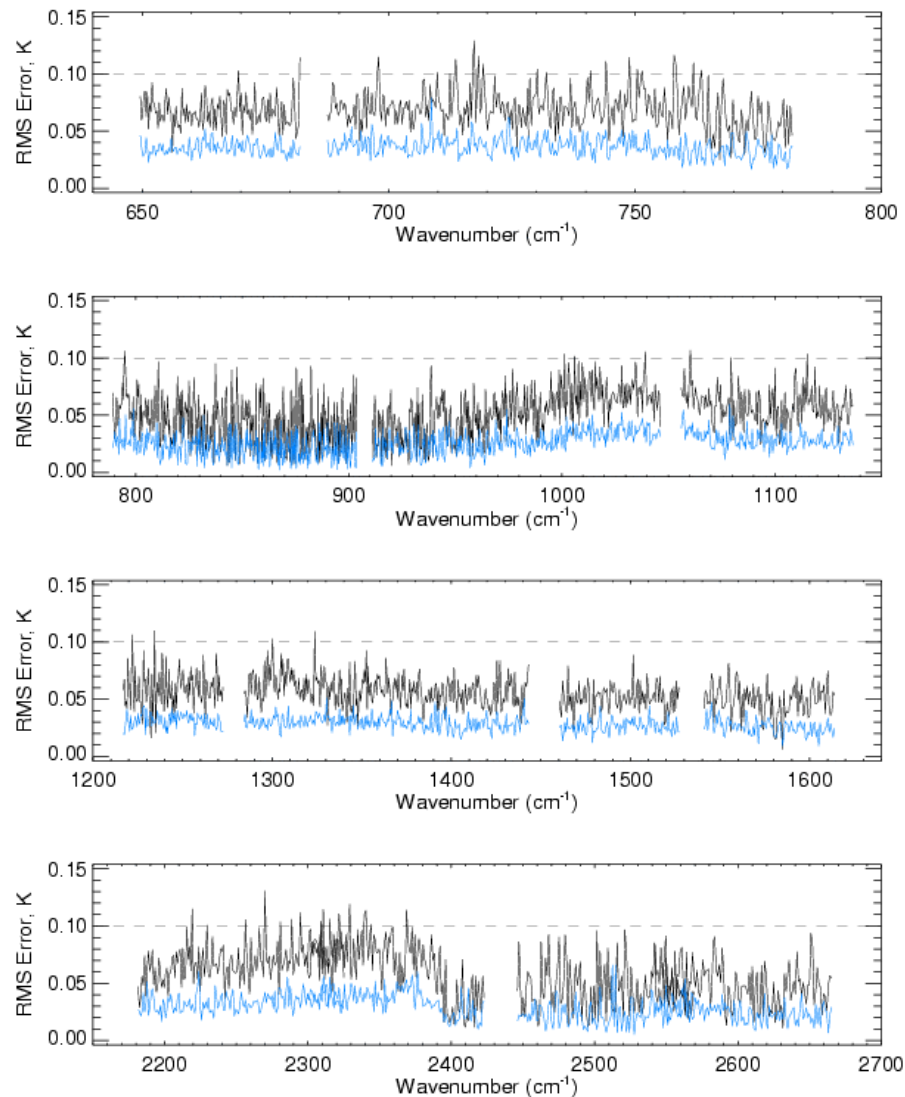
- 73-74% of nodes are common to at least two channels
- Average number of points per channel is 2.1 and 1.36 for 0.05K and 0.1K training, respectively

AIRS model accuracy

Validate: 52
diverse
ECMWF profile

Errors are due
to OSS
sampling and
optical depth
interpolation
(no radiative
transfer
errors)

- Independent set: 52 ECMWF profiles (see Saunders, 2003)
- Nadir viewing angle
- Surface emissivity =0.99
- Same 101-level RT scheme used for both LBLRTM and fast OSS model



Analytical Jacobians (1)

- With monochromatic RT, analytical Jacobian computation is trivial
- For example, in the case of a non reflective surface computation of Jacobians wrt molecular amounts reduces to

$$\partial R / \partial u_{ml} = D_l \partial \tau_l / \partial u_{ml}$$

where

$$D_l = \bar{B}_l T_l - \Sigma_{l+1}$$

Contribution to TOA
radiance of upwelling
radiation incident at bottom of
layer *l*

is independent of *X* and is derived in the process of merging layers successively (from bottom up) for radiance computation

- Temperature Jacobians:

$$\partial R / \partial \Theta_l = \frac{\partial B_l}{\partial \Theta_l} (T_{l-1} - T_l) + D_l \partial \tau_l / \partial \Theta_l$$

Analytical Jacobians (2)

- Derivative wrt amount for variable molecules

- Dry variable gases

$$\frac{\partial \tau_l}{\partial u_{ml}} = k_{ml}(\theta_l)$$

- Water vapor

$$\frac{\partial \tau(\bar{\theta}_l, \bar{q}_l)}{\partial u_w} = \frac{\partial \tau_w(\bar{\theta}_l, \bar{q}_l)}{\partial u_w} + \bar{k}_{fix} \frac{\partial u_{fix}}{\partial u_w}$$

$$+ \sum_{m \in dry} k_m \left(\frac{\partial u_m}{\partial u_w} \right)$$

$$\frac{\partial \tau_w(\bar{\theta}_l, \bar{q}_l)}{\partial u_w} = k_w(\bar{\theta}_l, \bar{q}_l) + \frac{\partial k_w(\bar{\theta}_l)}{\partial \bar{q}_l} \bar{q}_l$$

Dry gas amounts derived from relative concentration with respect to moist air (not mixing ratios)

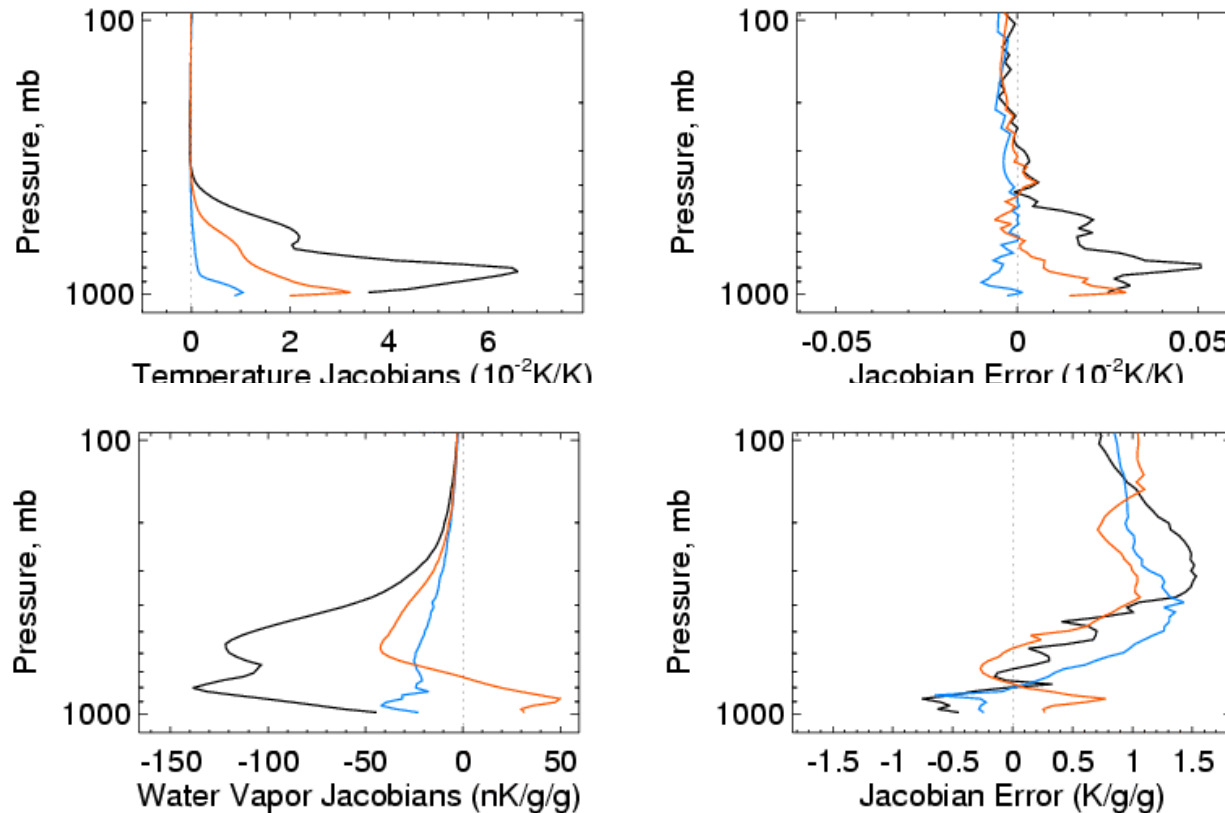
- Derivatives of optical depth wrt to temperature

$$\begin{aligned} \frac{\partial \tau(\bar{\theta}_l, \bar{q}_l)}{\partial \bar{\theta}} &= \frac{\partial k_w(\bar{\theta}_l, \bar{q}_l)}{\partial \bar{\theta}} u_w + \frac{\partial k_{fix}(\bar{\theta}_l)}{\partial \bar{\theta}} u_{fix} \\ &+ \sum_{m \in dry} \frac{\partial k_m(\bar{\theta}_l)}{\partial \bar{\theta}} u_m \end{aligned}$$

$$\begin{aligned} \frac{\partial k(\theta)}{\partial \theta} &= \frac{(2\theta - \theta_i - \theta_{i+1})}{(\theta_{i-1} - \theta_i)(\theta_{i-1} - \theta_{i+1})} (k_{i-1} - k_{i+1}) \\ &+ \frac{(2\theta - \theta_{i-1} - \theta_{i+1})}{(\theta_i - \theta_{i-1})(\theta_i - \theta_{i+1})} (k_i - k_{i+1}) \end{aligned}$$

AIRS Jacobians accuracy

- OSS Jacobians for temperature and water vapor compared to LBLRTM finite difference results
- Differences generally within 5 % (profile average) with 0.05 K model, reaching 8% (water vapor and ozone) for a few profiles



Computational Efficiency

- **Number of operation/layer for clear sky transmittance calculations at a single node:**
 - 3 multiplications per constituent (6 for water vapor) (see Slide 10 and 11)
 - 1 exponential
- **OSS and existing transmittance parameterizations (e.g. OPTRAN, RTTOV or SARTA) should become similar in terms of number of operations when number of nodes approaches ~2.5**
- **Average number of points per channel for AIRS model is 2.1 and 1.36 with 0.05K and 0.1K threshold**
- **Adding Jacobians wrt all state vector elements (atmosphere and surface) to radiance computation only doubles execution time (in clear sky)**

Future work

- Tune training for “non-localized” ILS (e.g. weakly apodized interferometer functions or narrowband imagers) for handling spectral variation of emissivity within band pass
- Tune training for clouds (with and w/o scattering in daytime and nighttime)
- Improve handling of variable viewing angles
- Trade improved schemes for treatment of layer emission (e.g. linear-in-tau) against number of levels