kCARTA : Fast pseudo line-by-line Radiative Transfer Algorithm for Allsky Infrared Sounding

sergio@umbc.edu

Sergio DeSouza-Machado¹

¹UMBC GESTAR2

June 16-17, 2025

Outline

- Motivation
- Line-by-Line codes (eg monoRTM, UMBC-LBL) many minutes/hours
- Pseudo Line-by-Line codes (eg kCARTA, LBLRTM, σ-IASI) seconds to minutes
- Fast Codes (eg SARTA, PCRTM, OSS) milliseconds
 - ► Flux (OLR) codes eg RRTM, ecRad milliseconds

Motivation

- New generation sounders : low noise, high spectral resolution sounders
 - ▶ NASA Atmospheric Infrared Sounder (AIRS) 09/2002-
 - ESA Infrared Atmospheric Sounding Interferometer (IASI) 06/2007-
 - NOAA Cross Track Infrared Sounder (CrIS) 03/2012-
- Designed to make observations for Numerical Weather Prediction
 - ► long/overlapping timespans allow for 40+ year record for climate studies
- Need fast, accurate RTA to fully exploit millions of daily observed radiances
 - accurate spectroscopy
 - \star latest line parameters eg HITRAN or GEISA databases
 - ★ accurate lineshapes : voigt or line mixing or
 - accurate radiative transfer : clear sky or allsky
 - esoteric needs (Non-Local Thermodynamic Equilibrium) NLTE
 - jacobians for retrievals
 - flux calculations for Outgoing Longwave Radiation (OLR) studies

Sun provides our energy

Earth Radiation budget balances incoming (shortwave) solar and outgoing terrestrial (longwave) radiation



$$T_{sun}, T_{earth}$$

$$P_{in} = \sigma A_{sun} T_{sun}^{4} = f 4 \pi r_{sun-earth}^{2}$$

$$P_{out} \rightarrow f \pi R_{earth}^{2} (1 - a) = 4 \pi R_{earth}^{2} \sigma T_{earth}^{4}$$

$$T_{sun} = 5600 \text{ K } T_{earth}^{radiating} = 255 \text{ K}$$

$$T_{earth}^{surface} = 285 \text{ K } f = 1360 \text{ W/m2 at}$$
TOA



Sergio DeSouza-Machado

Importance of the infrared to studying Earth Atmosphere

- Terrestrial Planck function peaks in infrared → OLR studies
- Greenhouse Gases (GHG) such as H₂O, CO₂, O₃, CH₄, CFCs have spectral signatures; also have temperature effects
- Clouds & large aerosols (\geq 4 μ m) impact infrared, 24/7
- Radiative Transfer Equation : height sensitivity



Spectroscopy

https://asl.umbc.edu/pub/packages/lbl.pdf

Atomic and molecular Spectroscopy - Basics 1



Literature

Introduction to Molecular Spectroscopy : Gordon Barrow First Course in Atmospheric Radiation : Grant Petty GENLN2 : https://opensky.ucar.edu/islandora/object/3A3407 UMBC LBL : https://asl.umbc.edu/pub/packages/lbl.pdf

Ideal Gas Law

Gas Cell

L =length of Gas cell

 $P, PS_g = \text{total pressure, gas partial pressure}$

$$T =$$
temperature (K)

R = molar gas constant (8.31 J/mol/K)

 $PS_g/P = mr_g = mixing ratio$

$$pV = nRT \rightarrow Q_g = \frac{n}{V}L = CF\frac{PS_g \times L}{R \times T}$$

 $\mathsf{CF}=\mathsf{conversion}\ \mathsf{to}\ \mathsf{eg}\ \mathsf{molecules}/\mathsf{cm2}$

Find optical depth of all gases $k(\nu, P, T) = \sum_{g} k_{g}(\nu, Q_{g}, P, PS_{g}, T_{g})$



Atomic and molecular Spectroscopy - Basics 3A

 $\mathsf{OD} = \mathsf{linestrength}(\mathsf{T}) \times (\mathsf{gas} \; \mathsf{amount} \times \mathsf{pathlength}) \times \mathsf{lineshape}$

$$[cm - 1/(molecules/cm2)]$$
{molecules/cm3 × cm}{ $1/cm - 1$ }

$$k_g^l(\nu, \nu_0) = S(T) \times Q_g \times f_{lineshape}(\nu, \nu_0^l)$$
$$k(\nu, P, T) = \sum_g k_g(\nu, Q_g, P, PS_g, T_g)$$

Master Equation for Spectroscopy

$$k(\nu, P, T) = \sum_{g}^{\text{all gases all lines}} \sum_{l}^{\text{all gases all lines}} k_g((\nu, Q_g, P, PS_g, T_g), (\text{lineshape}^{l}(\nu_0^{l}, \gamma^{l})))$$

Key idea

$$k_g(\nu, Q_g, P, PS_g, T_g) = Q_g \times S(T) \times f_{\textit{lineshape}}(\nu, P, PS_g, T) = Q_g \sigma(\nu, T, P, PS_g)$$

Atomic and molecular Spectroscopy - Basics 3B

Line Strength S(T)

Partition fraction Z(T) gives population stats for a given line line center shift due to pressure $\nu_0(P) = \nu_0(0) + \lambda P$ line strength $S(T) = S(296K) \frac{Z(296K)}{Z(T)} \times f(exp^{-hcE/kT})$

Line Broadening : Lorentz vs Doppler

$$\begin{split} \gamma_{L} &= (PS \times sbroad + (P - PS) \times fbroad) \times (296/T)^{n} \\ \gamma_{L}^{air}(P, T) &\sim \gamma_{0}^{air} \frac{P}{P_{ref}} (\frac{T_{ref}}{T})^{n} \sim 0.1 \frac{P}{P_{ref}} \text{ cm}^{-1} \\ \gamma_{D}(T) &= \nu_{0} (\frac{kT}{m})^{1/2} \sim \sqrt{T/m} \end{split}$$

Lineshape

Typically use Voigt lineshape for IR Note : H₂O, CO₂, CH₄ more complicated $f_{lorentz} = \frac{1}{\pi} \{ \frac{\gamma_L}{(\nu - \nu_0)^2 + \gamma_L^2} \}$ $g_{doppler} = \frac{1}{\gamma_D(T)} (\frac{ln2}{\pi})^{1/2} e^{-(ln2(\frac{\nu - \nu_0}{\gamma_D(T)})^2)}$ $f_{voigt} = f_{lorentz} \bigotimes g_{doppler}; f_{voigt} \rightarrow f_{lorentz} \text{ as pressure increases}$ $k(\nu) = Qf_{voigt}(\nu)S(T)$

Atomic and molecular Spectroscopy - Basics 2



Spectroscopic databases

- Publicly available spectroscopic databases include HITRAN and GEISA
- Line strengths, Line centers, Line widths
- Lineshape (typically voigt, though water/CO2 more complicated)

Sergio DeSouza-Machado

HITRAN

	NOM	e	\frown	\sim	Logged	n as Sergio Des	iouza-Mao	hado <u>Logo</u>
Home	Data Acce	ss D	ocumentation	Conferences	Links	About		
Line-by-Line S	Search						\$	0
4. Select or Crea	te Output For	nat						
Select an output for	mat or create a ne	v one, then	5. Start Data Search					
		_						
Available Output Fo	ormats	Ou	tput Format Descr	ription				
.par (160 chars)	000	.pa	r (160 chars)					
test2	(4 0 0	The	e 160-byte fixed-width	h format used since HITRA	N 2004 - see Table	1 in <u>Rothman</u>	et al., J	OSRT 96, 13
IESI/		(20	05)					
		(20	<u>05</u>).					
Create New Outpu	t Format	(20 Fie Thi	<u>05)</u> . ld separator: [no sepa s output format has v	arator], Line endings: Wine rariable-width fields and r	dows (CR LF) no header line.			
Create New Outpu	t Format	<u>(20</u> Fie Thi	0 <u>5</u>). ld separator: [no sepa s output format has v	arator], Line endings: Wine variable-width fields and r	dows (CR LF) no header line.			
Create New Outpu	t Format	<u>(20</u> Fie Thi	05). Id separator: [no sepa s output format has v	arator], Line endings: Wine variable-width fields and r	dows (CR LF) no header line.	<u>C</u> Fortran	Err	Pof
Create New Outpu	t Format	<u>(20</u> Fie Thi	05). Id separator: [no sepa s output format has v Parame Molecule ID	arator], Line endings: Wina araiable-width fields and r eter	dows (CR LF) 10 header line. Units	<u>C</u> Fortran Format I2	Err	Ref
Create New Outpu	t Format	(20 Fie Thi	05). Id separator: [no sepa s output format has v Parame 3 Molecule ID 3 Isotopologue	arator], Line endings: Win rariable-width fields and r eter ID	dows (CR LF) no header line. Units	<u>C</u> Fortran Format I2 I1	Err	Ref
Create New Outpu	t Format	(<u>20</u> Fie Thi	 dseparator: [no separator: soutput format has v Parame Molecule ID Isotopologue ν 	arator), Line endings: Win arariable-width fields and r eter ID	dows (CR LF) to header line.	C Fortran Format I2 I1 F12.6	Err	Ref
Create New Outpu	t Format	(<u>20</u> Fie Thi	 (1) S (2) S 	arator), Line endings: Wini rariable-width fields and r eter ID	dows (CR LF) to header line. Units cm ⁻¹ :/(molec.cm ⁻²)	C Fortran Format I2 I1 F12.6 E10.3	Err ✓	Ref ✓
Create New Outpu	tFormat	(<u>20</u> Fie Thi	 (55). kd separator: [no separato	arator), Line endings: Wini rariable-width fields and r eter ID	dows (CR LF) to header line. Units cm ⁻¹ /(molec.cm ⁻²) s ⁻¹	C Fortran Format I2 I1 F12.6 E10.3 E10.3	Err ✓	Ref ✓ ✓
Create New Outpu	t Format	(<u>20</u> Fie Thi	(d separator: [no separator:]no separator:]no separator:]no separator Parame Molecule ID Isotopologue ν S A γ _{air}	arator), Line endings: Win rariable-width fields and r eter ID cm ⁻¹	dows (CR LF) to header line.	 Si Fortran Format I2 I1 F12.6 E10.3 E10.3 F5.4 	Err ~ ~	Ref ~ ~ ~

₿ E"

6 n_{air}

 $\boldsymbol{6} \delta_{\mathrm{air}}$

F10.4

F4.2 \checkmark ~

<u>F8.6</u> ~ \checkmark

cm⁻¹

cm⁻¹·atm⁻¹

Line-by-Line codes

https://asl.umbc.edu/pub/packages/lbl.pdf



UMBC-LBL

Basics

- Matlab based
- input : gasID,(T,P,PS,Q),($\nu_{start}, \nu_{end}, d\nu$),(preset parameters)
- output : $\nu = (\nu_{start} : d\nu : \nu_{end} d\nu), k(\nu); d\nu = 0.0025 \text{ cm}^{-1};$
- Reads in HITRAN/GEISA spectral database line parameters
- Loops over lines, applying voigt lineshape to each line (Fortran Mex)
- \bullet Does H_2O (without basement, HDO, continuum) and CO_2 (line mixing)

Size/Timings

- 40 molecular gases H₂O,CO₂,O₃,CH₄ plus HDO, self+foreign continuum
- 40 cross section gases CFCs
- 100 layers : 1100 mb \geq Pressure \geq 0.005 mb (0 km \leq z \leq 82 km); thickness 250 m at GND, 4 km at TOA); span 605-2830 cm^{-1} in 25 cm^{-1} chunks
- Takes a long time to compute ODs for one atmosphere

How to speed up parts of the code

The "preset parameters" in the previous slide include nbox = number of points to boxcar integrate, and width of "near", "medium", "far" intervals

Algorithm

- 1 cm⁻¹ chunks, $d\nu/nbox$ resolution
- bin lines : "near" "medium" "far"
- "near" within ± 1 cm⁻¹ of the bin
 OD at fine resolution dν/nbox (red), box car integrate to dν
- "medium" within ± 1-2 cm⁻¹ of the bin : OD at resolution 0.1 cm⁻¹, interpolate to dν (green)
- "far" within \pm 2-25 cm⁻¹ of the bin : OD at resolution 0.5 cm⁻¹, interpolate to d ν (not shown)
- repeat at next 1 cm⁻¹ interval



Example : Ozone lines from HITRAN 2020

Some numbers

46985 lines In the 980-1080 cm $^{-1}$ interval; choose the 1020-1021 cm $^{-1}$ bin

- 1629 lines in the fine bin [1019,1022] cm^{-1}
- \bullet 1099 lines in the medium bins [1018,1019] $\rm cm^{-1}$ and [1022,1023] $\rm cm^{-1}$
- \bullet 23194 lines in the far bins [0995,1018] $\rm cm^{-1}$ and [1023,1046] $\rm cm^{-1}$



Pseudo Line-by-Line codes : kCARTA

https://asl.umbc.edu/pub/packages/kcarta1.22.pdf



How to speed this up

Some challenges

- About 8 seconds for 1 cm⁻¹, about 36 seconds to do 25 cm⁻¹, for one layer
- Need 100 layers so about 60 minutes to do 25 cm⁻¹; to span 605-2830 cm⁻¹ would take 89 chunks = 89 hours
- We have 80 gases so about
- It's not so bad, there are about 6 main HITRAN gases; not all gases have so many lines, in every spectral region
- Very Accurate! Too slow to be practical for remote sensing applications!

T and Q profiles

49 AFGL regression profiles, including Tropical, Mid Lat Summer, Mid Lat Winter, Sub Arctic Summer, Sub Arctic Winter (cold/hot and dry/wet)



How to speed this up A

Idea A : Make a database

٠

$$k_g(\nu, T) = Q_g \sum_{l}^{all \ lines} S_l(T) f_{voigt}(l, \gamma(P, PS_g, T)) = Q_g k_g(T)$$

- Make database in term of $T_{USStandard}(z)$, $Q_{USStandard}(g, z)$ for all 80 gases
- Use 10 K temperature offsets from US Standard (reference) profiles : $T_{ref}(z) - 50, T_{ref}(z) - 40, ... T_{ref}(z), ... T_{ref}(z) + 40, T_{ref}(z) + 50$
- \bullet Will take time to generate, but when saved can find optical depths for arbitrary profile $(\mathsf{T}(z), Q(g, z))$
 - n = -5 : 1 : +5 (11 temperature offsets)
 - interpolate in temperature k^{ref}_g(v, z, T), Q^{ref}_g(z)) = interp(T_{ref}(z) + n * 10, k^{ref}_g(v, T_{ref}(z) + n * 10, Q^{ref}_g, z), T(z)).
 - ► scale with gas amount $k_g(\nu, z, T) = k_g^{ref}(\nu, z, T), Q_g^{ref}(z)) \frac{Q(z)}{Q_r^{ref}(z)}$

Idea B : Reduce Database Size!

- The raw database is huge : (80 gases) \times (89 chunks of 25 cm⁻¹ at 0.0025 cm⁻¹ spacing) \times (100 layers) \times (11 T offsets)
- There are spectral features : use Singular Value Decomposition to compress
- $k_g(\nu, T, P) \rightarrow k_g^{compressed}(\nu, T, P)$
- We call this the kCompressed database
 - check compression by running radiative transfer and ensuring errors in TOA brightness temperature are less than 0.1 K
- Generated using H1998, H200,H2004, H2008, H2012, H2016, H2020, ... (and also GEISA 2015)
- Also have done (FIR) 10-605 cm⁻¹ and (NIR,Vis,UV) 2830-44000 cm⁻¹ (making sure point spacing d ν changes with central wavenumber); but older H2004,H2008



Raw database

- \bullet Use UMBC-LBL to compute ODs for \sim molecular gases and \sim cross section gases, using US Standard Temperature and gas profiles
 - ▶ for H₂O we use 5 different variable water widths to account for self broadening of water
 - ▶ We use LBLRTM (v12.8) to generate CO₂ and CH₄ since they have more recent line mixing coefficients
- 25 cm⁻¹ chunks at 0.0025 cm⁻¹ resolution (10000 points per chunk) spanning 605-2830 cm⁻¹ (89 chunks)
- 100 Pressure layers spanning 1100 mb (surface) to 0.005 mb (TOA)
- $\bullet\,$ 11 Temperature offsets from the US Standard Temperature profile : 0 K, $\pm\,$ 10 K, \pm 20 K, ..., \pm 50 K
- For each gas, each 25 cm-1 chunk $k_g(\nu, T, P) = (100 \times 11) \times 10000$
- Takes about a week to grind out on a HPC
- $\bullet\,$ Database size $\sim\,60\,$ Gb

SVD compressed database

• $m \times n$ Matrix of absorption coefficients K ($m = 100 \times 11, n = 10000$) $K_{ij} = U\Sigma V^T = \sum_{l=1}^{L} U_{il} \Sigma_l V_{jl} + \sum_{l=L+1}^{n} U_{il} \Sigma_l V_{jl}$

• Many of the singular values are small ... drop the second term. $U = m \times n, \Sigma = n \times n, V = n \times n \rightarrow U = m \times L, \Sigma = L \times L, V = n \times L.$ $L \simeq 30 - 40$ for H_2O, CO_2 $L_{avg} \simeq 5$ for rest of gases $K = U\Sigma V^T \simeq \hat{U}\hat{K}$ $\hat{K}_{ij} = \sum_{l=1}^{L} \Sigma_l V_{jl}$

• Reconstructing the absorption coefficients for the *m*-th layer

$$K_{m(g)}(\nu) = \frac{q_{m(g)}}{q_{m(g)}^{ref}} \sum_{l=1}^{L} \hat{k}_{l(g)}(T_m, m) \ \hat{u}_l$$

 \hat{u}_l are basis vectors, $\hat{k}_{l(g)}(T_m)$ are interpolated coeffs from $\hat{K}(T)$ • Database ~ 776 Mb (200 Mb (H₂O), 76 Mb (CO₂), 500 Mb (other gases))

https://asl.umbc.edu/pub/packages/kcarta1.22.pdf https://amt.copernicus.org/articles/13/323/2020/

Sergio DeSouza-Machado



Add on a Radiative Transfer Algorithm

Clear Sky

- After you have the optical depths, add in : solar source term, emissivity, reflectivity, ray tracing and do Radiative Transfer
- kCARTA : the k-Compressed Atmospheric Radiative Transfer Algorithm
- Analytic jacobians, can do flux calculations
- Takes about 30-45 seconds to uncompress and calculate TOA (or ground) infrared radiances $R(\nu)$ (605-2830 cm⁻¹) for an arbitrary profile
 - ► convolve with appropriate instrument Spectral Response Function
 - $R_i^{calc}(\nu) = \int d(\nu) R(\nu) SRF_i(\nu)$
- Takes about x3 longer to do 100 layer analytic jacobians dBT/dT(z),dBT/dQ(z),dBT/dST
- WARNING Since we use 0.0025 cm^{-1} resolution throughout the (605-2830 cm^{-1}) spectral range
 - noticeable differences when you compare with LBLRTM since that uses variable (higher) spectral resolution
 - ► but after you convolve with an instrument SRF, these differences are very small





Details

 $R(\nu, \theta) = R_{surface} + R_{atmospheric emission} + R_{background thermal} + R_{solar}$

- $R_{surface}(\nu, \theta)$ = surface to TOA at angle θ
- $R_{atmospheric\ emission}(\nu, \theta)$ = thermal emission from each angle at angle θ
- $R_{background \ thermal}(\nu, \theta) =$ downward thermal radiation from all angles, reflected at surface at angle θ
- $R_{solar}(\nu, \theta, \theta_{sun}) =$ downward solar radiation at θ_{sun} reflected at surface at angle θ

Clear Sky

At steady state, 1D radiative transfer (RT) described by Schwartzchild equation (ignoring solar and background contributions)

$$\mu \frac{dR(\nu,\theta)}{\kappa_e dz} = -R(\nu,\theta) + J(\nu)$$

- $\mu = cos(\theta)$, dz is the vertical coordinate
- κ_e is the total extinction (due to gases, clouds etc)
- $\kappa_e dz = dk$ is the optical depth
- $R(\nu, \theta)$ is the radiance intensity we want to solve for
- *J* is the source function
 - Clear Sky, in LTE : $J = B(\nu, T)$
 - Clear Sky, in NLTE : $J = rB(\nu, T)$



Soln to Schwartzchild Eqn

$$R(\nu,\theta) = \frac{\epsilon_{s}(\nu)B(\nu,T_{s})\tau_{atm}(\nu,\theta) + \int_{surface}^{TOA} B(\nu,T(z))\frac{\partial\tau(\nu,\theta)}{\partial s}ds + \frac{1-\epsilon_{s}(\nu)}{\pi}\tau_{atm}(\nu)\int d\Omega^{+}\int_{TOA}^{surface} B(\nu,T(s))\frac{\partial\tau(\Omega)}{\partial s}\cos(\theta)ds + \rho_{s}(\nu)B_{\odot}(\nu)\cos(\theta_{\odot})\tau_{atm}(\nu,\theta_{\odot})\tau_{atm}(\nu,\theta)$$

Layer by Layer Iterative Solution

- Surface Emission : $R_s(\nu) = \epsilon_s(\nu)B(\nu, T_s)\tau_{GND \to TOA}(\nu, \theta)$
- Layer Emission : $R_{\text{lay}} = \sum_{i=1}^{i=N} B(\nu, T_i)(1.0 \tau_i(\nu))\tau_{i+1 \to TOA}(\nu, \theta)$
- Background thermal : $R_{th}^{surface}(\nu) = \pi \rho_s \sum_{i=N}^{i=1} B(T_i) [\tau_{i-1 \rightarrow ground}(\nu, \theta_{diff}) - \tau_{i \rightarrow ground}(\nu, \theta_{diff})]$ We speed up the integral over hemispheres by using layer by layer/spectrally varying effective diffusive angles; see AMT paper
- Solar reflected : $R_{\bigcirc}(\nu) = \rho_s(\nu, \theta, \phi) B_{\bigcirc}(\nu) cos(\theta_{\bigcirc}) \times \tau_{N \to ground}(\nu, \theta_{\bigcirc})) \tau_{ground \to TOA}(\nu, \theta)) \Omega_{\bigcirc}$

Typical AIRS spectrum (KCARTA convolved with AIRS SRFs)





Comparisons to LBLRTM

- kCARTA uses fixed resolution 0.0025 $\rm cm^{-1}$
- LBLRTM varies spectral resolution with wavenumber
- (L) kCARTA less accurate on line centers at larger wavelengths (3K errors at 15 um); improves as wavelengths decrease (0.1 K errors at 10 um)
- (R) After convolution with instrument SRFs, on average kCARTA-LBLRTM within NeDT. Plots averaged over 49 profiles.



kCARTA 1D RTA Extras : (A) Jacobians

Jacobians

$$R(\nu) = \epsilon_s B(T_s, \nu) \tau_{1 \to TOA}(\nu) + \sum_{i=1}^{i=N} B(T_i, \nu) (1.0 - \tau_i(\nu)) \tau_{i+1 \to TOA}(\nu)$$

Differentiation with respect to the *m*-layer variable s_m , (gas amount or layer temperature $s_m = q_{m(g)}, T_m$) yields

$$\frac{\partial R(\nu)}{\partial s_m} = \left[\epsilon_s B(T_s)\tau_{1\to TOA}\right](-1)\frac{\partial k_m(\nu)}{\partial s_m} + \left[\sum_{i=1}^{m-1}(1.0-\tau_i(\nu))B_i(\nu)\tau(\nu)_{i+1\to TOA}\right](-1)\frac{\partial k_m(\nu)}{\partial s_m} + \left[(1.0-\tau_m(\nu))\frac{\partial B_m(\nu)}{\partial s_m} - B(T_m,\nu)\frac{\partial \tau_m(\nu)}{\partial s_m}\right]\tau_{m+1\to TOA}(\nu)$$

Individual jacobian terms $\frac{\partial k_m}{\partial s_{m(g)}}$ are as follows

- gas amount derivative is $\frac{\partial k_m}{\partial q_{m(g)}} = \frac{k_m}{q_{m(g)}}$ (with added complexity for water, to account for self broadening),
- temperature derivative $\frac{\partial k_m}{\partial T}$ cumulatively obtained *while* performing the temperature interpolations during individual gas database uncompression.



Tropical Jacobians (kCARTA convolved with AIRS SRFs)

Gas Jacobians are dBT/d(log(Q(p))) in K Temperature Jacobian is dBT/dT(p) in K/K



Sergio DeSouza-Machado

Column Tropical Jacobians (kCARTA convolved with AIRS SRFs)

Gas Jacobians are dBT/d(log(Q(p))) in K Temperature Jacobian is dBT/dT(p) in K/K



Sergio DeSouza-Machado

Variation in local zenith angle

- Can use plane parallel atmosphere
- Default is to slowly vary angle as ray propagates through spherical layers



satellite, while *I* is the sun

Variation in layer temperature with optical depth

- Default : constant layer temperature
- eg LBLRTM use linear-in-tau variation of temperature; let $\tau(\nu) = e^{-\kappa(\nu)}$

$$I(\nu) = I_0(\nu)\tau + (1-\tau)\left\{B_{a\nu}(\nu) + (B_u(\nu) - B_{a\nu}(\nu))\left(1 - 2\left(\frac{1}{\kappa} - \frac{\tau}{1-\tau}\right)\right)\right\}$$

kCARTA 1D RTA Extras : (C) NLTE

NLTE Computation

- $\bullet\,$ Daytime Solar radiation is preferentially absorbed by some CO_2 and O_3 bands
- This gives higher kinetic energies and hence vibrational temperatures than other bands/other molecules for $z \ge 40$ km
- $\bullet\,$ Nadir sounders (AIRS/CrIS/IASI) : 10 K effect at 4.3 um T(z) sounding region
- $\bullet\,$ Due to the Planck function \ll in 15 um (limb sounders see effect)
- Manuel Lopez Puertas (U. of Granada, Spain) has given AIRS team CO2 vibrational temperatures as function of solar angle and temperature profile
- Some bands remain in LTE, others go into NLTE
- UMBC has developed a line-by-line model for kCARTA to make a fast model for SARTA
- Lorenzo Cassini (UniBas) currently visiting UMBC to develop a NLTE model using LBLRTM for $\sigma\text{-}\mathsf{IASI}$
- GENLN2, KOPRA and RFM also have NLTE models

Tvib vs Tkin

TRP profile; Solzen = 0 and Solzen = 90 Tvib \rightarrow Tk as day \rightarrow night



Modelling for SARTA

- 3 models since 2005
- Manuel has tweaked Tvib collisional model
- We tweaked our fitting parameters
- Have not fitted latest 2025 data
- O-C Bias (AIRS) has oscillated between about ± 1 K

Sergio DeSouza-Machado

NLTE : Basic Idea

- LTE : compute ODs as usual (contribution "in-between" lines, using local kinetic temperature T(z))
- NLTE : 'laser'' effect (enhanced OD and Planck emission) mostly ''on top of lines/line centers'' using vibrational temperature $T^i_{vib}(z)$

OD Modification

- $\tau_{\textit{NLTE}}(\nu)$: needs additional factor β
- As solar angle θ_{\odot} increases beyond 90°the factor $\alpha_{\textit{NLTE}}(\nu, \theta_{\odot}) \rightarrow 1$
- In reality there is a ''dusk'' effect (θ_{\odot} between 90° and 120°) which we have difficulty modeling accurately
- Nadir low earth orbit sounders such as AIRS, CrIS, IASI have
 - $\blacktriangleright~\theta_{\bigodot}$ between 10-60 $^\circ {\rm in}$ tropics and mid latitude regions
 - θ_{\odot} above 60 °in polar regions
 - Plus there can be aurora events in the polar regions

$$\eta(\nu,\theta_{\odot}) = \sum_{LTE} \tau_{LTE}(\nu) + \sum_{NLTE} \alpha_{NLTE}(\nu,\theta_{\odot}) \tau_{NLTE}(\nu)$$

NLTE : Planck Modification

As solar angle θ_{\odot} increases beyond 90°the factor $\beta_{\textit{NLTE}}(\nu, \theta_{\odot}) \rightarrow 1$

$$B_{NLTE}(\nu, \theta_{\odot}) = B_{Planck}(\nu, T)\beta_{NLTE}(\nu, \theta_{\odot})$$

 $-I(\nu, \theta, \theta_{\odot}) + \beta(\nu, \theta_{\odot})B_{Planck}(\nu, T)$

NLTE : Rad Transfer Modification

dz _{NLTE}	$= dz \eta(\nu, \theta_{\odot})$
$dI(\nu,\theta,\theta_{\odot})$	_
dzNITE	—

NLTE : Quick summary

- Daytime solar : some bands of CO₂ can be in NLTE above 40 km
- Need to modify the optical depth and the planck function
- Between 0-30 km, do usual LTE calc for radiative transfer
- Above 30 km, do NLTE calc for radiative transfer
- But you need the vibrational temperatures to be able to do this

Work done by Lorenzo Cassini (for future use in σ -IASI)

NLTE - LTE : (LBLRTM)

- function of solar zenith angle
- 415 ppmv for CO₂. Very small changes as CO₂ changes (0.25 K/40 ppmv)
- Largest effect is at solzen = 0°, \rightarrow 0 as solzen becomes 90°



NLTE - LTE : LBLRTM - KCARTA

- KCARTA : NLTE line mixing coeffs made in 200
- LBLRTM : Niro et. al (2012)
- US Standard Profile, 415 ppmv
- Other profiles show similar results



June 2019 Random observations, 2350 cm⁻¹ channel Observations



kCARTA 1D RTA Extras : (D) Flux

Flux Computation

- Default : 0.0025 cm^{-1} resolution is acceptable for integration with SRFs
- Limitations become evident when you integrate over wavenumber to get fluxes and/or heating rates
- When we use 0.0005 cm⁻¹ reolution, we approach LBLRTM heating rates and fluxes (but this slows down our code)



Sergio DeSouza-Machado

kCARTA 1D RTA Extras : (E) Scattering

All Sky

$$u \frac{d(\nu)}{dk_{e}} = -I(\nu) + B(\nu, T)(1 - \omega_{0}) + \frac{\omega_{0}}{2} \int_{-1}^{+1} I(\nu, k_{e}, \mu) P(\mu, \mu) d(\mu) + \frac{\omega_{0}}{4\pi} \pi I_{sun} P(\mu, -\mu_{sun}) e^{-k_{e}/\mu_{sun}}$$

- $\omega = k_s/k_e = 1 k_a/k_e$ is the single scattering albedo (0 for no scatter)
- $P(\mu,\mu')$ is probability of scattering from μ' into μ
- $P(\mu, -\mu_{sun})$ is probability of scattering from μ_{sun} into μ
- Can interface to DISORT or to PCLSAM to do scattering calculations
- This also has analytic jacobians for the PCLSAM version
 - ► Have also put in a correction (by G.Tang and P.Yang)
- Mie scattering for water clouds and aerosols, have Ping Yang/Bryan Baum scattering database for cirrus (different habits)

DISORT

- Solves integro-differential equation by using a number of streams (2,4,8, ...)
- Can be "slow" for line-by-line applications, especially if you also try to use it to do the RT for the clear sky part
- Stamnes, K., Tsay, S.-C., Wiscombe, W., and Jayaweera, K.: Numerically Stable Algorithm for discrete ordinate method Radiative Transfer in multiple scattering and emitting layered media, Appl. Opt., 27, 2502–2509, 1988

PCLSAM

- $\bullet\,$ More accurate in cases where single scattering albedo much smaller than 1
- Effective scattering of cloud and aerosol becomes very similar to an effective cloud optical depth $k_{total}(\nu) = k_{gas}(\nu) + k_{cloud}^{eff}(\nu, \omega)$
- Radiative Transfer now like earlier slides (fast/analytic jacobians)
- Chou, M.-D., Lee, K.-T., Tsay, S.-C., and Fu, Q.: Parameterization for Cloud Longwave Scattering for use in Atmospheric Models, J. Climate, 12, 1999
- G. Tang and P.Yang (2018) : Improvement of the simulation of cloud longwave scattering in broadband radiative transfer models, JAS (75)

Two Slab Clouds

- NWP models have ice/water cloud profiles (91 levels); but there are typically very few deg of information in IR radiances
- UMBC has developed a Two Slab Cloud approximation to reduce these profiles to two randomly overlapping clouds, using NWP information
 - Cloud/Aerosol 1 : type, fraction, amount, effective particle size, cloud top/bottom
 - Cloud/Aerosol 2 : type, fraction, amount, effective particle size, cloud top/bottom
 - Cloud Overlap fraction (random)
- Slab placement can be adjusted "as high as possible" "median of distribution" to get better agreement with eg DISORT based RTAs
- DeSouza-Machado, Strow, Tangborn, X. Huang, X. Chen, X. Liu, W. Wu, Q. Yang, "Single-footprint retrievals for AIRS using a fast TwoSlab cloud-representation model and the SARTA all-sky infrared radiative transfer algorithm", AMT 2018

Example

Cloud vertical profiles, reduced to one or two slabs. The red and blue curves come from the NWP model, while the cyan and magenta are the resulting locations (and loadings) for the slabs.



Fast RTA codes : SARTA

https://asl.umbc.edu/pub/packages/sarta.html

Fast Models

Need more speed

- Pseudo LBL codes are fast eg kCARTA (30 seconds), σ -IASI (0.8 seconds) per 0.0025 cm⁻¹ spectrum
- New generation sounders (AIRS, CrIS, IASI) 2-3 million spectra daily
- The sounders make observations with finite number of channels (eg AIRS has 2378 channels, IASI has about 8400 channels)

Fast Models

Basic Idea

- We use 49 regression profiles spanning Earth's typical temperature/H₂O/O₃ and other trace gases (eg CO₂,CH₄) to produce convolved layer-to-space transmittances for each instrument channel $\tau_{L2S}(i, l) = \int_{d\nu} e^{-\sum_{j=l+1}^{N} (K(\nu, j)/\cos(\theta))} SRF_i(\nu)$
- We then fit these convolved transmittances using linear combinations of gas amounts, temperatures, view angles such as
 τ_{L2S}(i, l) = a_lT(l) + a_{l+1}T(l + 1) + ... + a_NT(N) + q_lQ_g(l) + s_lsec(θ)
 τ_{L2S}(i, l) = Predictors(T(l), T(l + 1), ...T(N), Q_g(l), θ) × Coefficients
- Then $k_{eff}(i, l) = -log(\tau_{L2S}(i, l))$ and we can do clear sky/all sky radiative transfer exactly as for kCARTA, also with analytic jacobians
- SARTA = Stand Alone Radiative Transfer Algorithm takes about 0.02 seconds/AIRS spectrum; also has Two Slab Cloud Scattering
- Can then use this SARTA Fast Model in AIRS L2 operational retrievals at NASA or NOAA

Fast Models

SARTA TwoSlab Examples

The *i*-th channel all-sky radiance $r_i(\nu)$ is computed using four weighted radiance streams : f_{clr} is clear fraction, $c_{overlap}$ is cloud overlap, cx1, cx2 are exclusive cloud fractions $c_i - c_{overlap}$

$$r_i(\nu) = f_{clr}r_i^{clr}(\nu) + c_{overlap}r_i^{(12)}(\nu) + cx_1r_i^{(1)}(\nu) + cx_2r_i^{(2)}(\nu)$$







observations

Conclusions



Conclusions

- Accurate Radiative Transfer needed for weather and climate
 - number of layers (more layers, more accurate, slower)
 - spectroscopy (most gases : voigt; CO₂, CH₄ needs line mixing)
 - new line shapes being used (speed dependent voigt)
- Monochromatic line-by-line codes are very accurate
 - Accuracy does depend on the lineshape used
 - Can take minutes/gas for a 25 cm⁻¹ interval
- Pseudo line-by-line codes are quite accurate
 - kCARTA optimized to be accurate enough for current generation high spectral sounders R_i^{calc}(ν) = ∫ d(ν)R(ν)SRF_i(ν)
 - Monochromatic inaccuracies versus eg LBLRTM (0.0025 cm⁻¹ spacing) can easily be reduced by using eg 0.0005cm⁻¹ database (but slows down code)
 - ▶ takes about 30-45 seconds to run 605-2830 cm-1
 - ► can do analytic jacobians, fluxes, all sky scattering
- Sounders provide millions of daily radiance spectra
 - ► SARTA Fast Model takes 0.02 second/2378 channel AIRS spectrum
 - ► can do analytic jacobians, all sky scattering, fluxes (eg RRTM, ecRad)

Public code repositories

UMBC

Matlab UMBC LBL : https://github.com/sergio66/UMBC_LBL Matlab kCARTA : https://github.com/sergio66/kcarta F90 kCARTA : https://github.com/sergio66/kcarta_gen F90 SARTA : https://github.com/strow/sarta

AER

LBLRTM : https://github.com/AER-RC/LBLRTM RRTM : https://github.com/AER-RC/RRTMG_LW

ECMWF

RTTOVS

ecRad : https://github.com/ecmwf-ifs/ecrad

NOAA GFDL

pyLBL : https://github.com/GRIPS-code/pyLBL GRTcode : https://github.com/NOAA-GFDL/GRTCODE

and many others (eg PCRTM, OSS etc) \ldots

Sergio DeSouza-Machado

Thanks

L. Larrabee Strow and Scott Hannon and Howard Motteler

AIRS Science Team Chris Hepplewhite, Steve Buczkowski NLTE : Manuel Lopez Puertas, David Edwards, Lorenzo Cassini

