

## Monte Carlo RT

The traditional transfer equation is

$$\hat{n} \cdot \vec{\nabla} I_{\nu} = \underbrace{-\chi_{\nu} p I_{\nu}}_{\text{absorption}} + \underbrace{j_{\nu}}_{\text{emission}} + \underbrace{\sum_i \eta_i \sigma_{\nu}^i \int \left( \frac{1}{\sigma_{\nu}^i} \frac{d\sigma_{\nu}^i}{d\Omega} \right) I_{\nu}(\hat{n}') d\Omega'}_{\text{scattering}}$$

" $g(\hat{n}', \hat{n})$ " = "phase function"

This is actually the same as the Boltzmann transport equation for particles; i.e., photons.

$$\frac{\partial f}{\partial t} + \vec{v} \cdot \vec{\nabla} f = \left( \frac{\partial f}{\partial t} \right)_{\text{interaction}} \quad f = \text{phase space \# density of photons}$$

since the specific intensity

$$I_{\nu}(\vec{x}, \hat{n}) = \frac{dE}{dA_1 dt d\nu d\Omega}$$
$$= c h \nu f$$

photons travel with velocity

$$\vec{v} = c \hat{n}$$

and in steady-state  $\frac{\partial f}{\partial t} \rightarrow 0$

So

Transport eqn = Boltzmann equation

Physically, the transfer equation says

- i) In absence of interactions the number of photons along any ray is constant; photons travel in straight lines
- ii) Absorption removes photons from the beam = destruction
- iii) emission adds photons to beam = creation
- iv) scattering removes photons from one direction, changes them to another = scattering conserve photon #

## Traditional Solution to RT equation:

- integrate along path (ray tracing)

$$I_{\nu} = e^{-\tau} I_{\nu}^0 + \int S e^{-\tau'} d\tau'$$

- Problem:  $S = \frac{j_{\nu} + \int n_{\nu} g I_{\nu}}{\chi_{\nu}}$

depends on solution for  $I_{\nu}$  from all other directions at each point along line of sight

⇒ must integrate every direction, at every point for every frequency + iterate! (lambda-iteration)

- Problem:
  - 1) 6-D integration is computationally expensive & memory intensive
  - 2) Temperature & level populations depend on radiation field ⇒ opacity and emissivity not known a priori.
  - 3) lambda-iteration fails without special techniques that enforce energy (flux) conservation

## Monte Carlo RT

- probabilistic simulation of the transport of photons (flow of energy) in the system
- It is an exact soln to the transfer eqn (using a statistical sampling of the paths)

## Photon Packets

- Divide the total system luminosity  $L$  into  $N_p$  packets of energy. The packet energy

$$E_p = \frac{L \Delta t}{N_p}$$

$\Delta t =$  arbitrary simulation time

- Each packet is an ensemble of physical photons with the same position, direction, and frequency. The number of physical photons in the packet is

$$n = E_{\gamma} / h\nu$$

- Since each packet is an ensemble of particles it may be partially polarized. Normalizing the Stokes parameters to the packet energy

$$I = 1$$

$$Q = (E_{\uparrow} - E_{\downarrow}) / E_{\gamma}$$

$$U = (E_{\rightarrow} - E_{\leftarrow}) / E_{\gamma}$$

$$V = (E_{\odot} - E_{\ominus}) / E_{\gamma}$$

Photon Weights

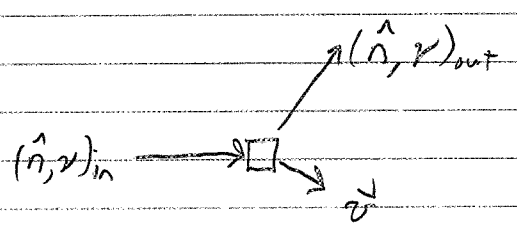
- Sometimes we need to use photon packets with unequal energies
- Define the "weight" of the packet

$$W = \frac{\text{packet energy}}{E_{\gamma}}$$

- Example: Coherent scattering in moving media

In the rest frame of the gas, photon energy is conserved; in lab frame number of photons is unchanged

However photons are Doppler shifted, so new packet energy is



$$E_{\gamma}^{out} = \left( \frac{\nu_{out}}{\nu_{in}} \right) E_{\gamma}^{in}$$

so the packet weight is updated

$$W_{out} = \left( \frac{\nu_{out}}{\nu_{in}} \right) W_{in}$$

After a number of scatterings

$$W = \Pi \left( \frac{v_{out}}{v_{in}} \right)_{\text{scatterings}}$$

- Note: 1) for non-relativistic motion this effect is often ignored  
 2) the total change in packet energy is the work done by the radiation force on the gas

### MC Probabilities

- To simulate the transport of photons, we require the probabilities of photon emission and interaction
- For equal energy packets, number of packets is proportional to the energy

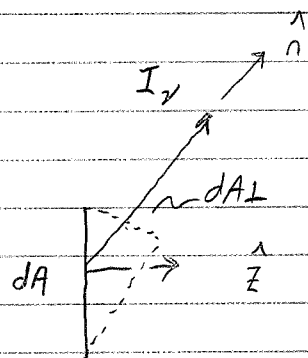
⇒ energy distributions = probability distributions

#### i) Emission from a planar surface

the specific intensity

$$I_\nu = \frac{dE/dt}{dA d\nu d\Omega}$$

but  $dA_L = \mu dA$      $\mu = \cos\theta = \hat{n} \cdot \hat{z}$



$$\therefore \mu I_\nu = \frac{dE/dt}{dA d\nu d\Omega}$$

So the probability distribution

$$\frac{dP}{dA d\nu d\Omega} \propto \mu I_\nu$$

integrating over all outward directions we find

$$\frac{dP}{dA d\nu} \propto \int_{2\pi} \mu I_\nu d\Omega = F_\nu \quad (\text{or } H_\nu)$$

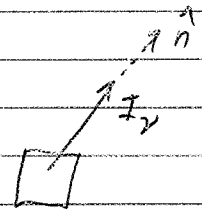
and then integrating over all frequencies

$$\frac{dP}{dA} \propto \int_0^\infty F_\nu d\nu = F \quad (\text{or } H)$$

ii) Emission from a volume element  $dV$

The emissivity

$$j_\nu = \frac{dE/dt}{dV d\nu d\Omega}$$



so the probability distribution

$$\frac{dP}{d\nu dV d\Omega} \propto j_\nu$$

integrating over direction followed by frequency gives

$$\frac{dP}{d\nu dV} \propto 4\pi j_\nu$$

$$\frac{dP}{d\nu} \propto 4\pi \int_0^{\infty} j_\nu dV$$

iii) Star + Envelope

The total system luminosity

$$L = L_* + L_{env}$$

where

$$L_* = 4\pi \int H dA$$

$$L_{env} = 4\pi \int j_\nu dV$$

The number of stellar and envelope photons is

$$N_* = (L_* / h\nu) N_\gamma$$

$$N_{env} = (L_{env} / h\nu) N_\gamma$$

so the probability of stellar and envelope photons is

$$P_* = L_* / (L_* + L_{env})$$

$$P_{env} = L_{env} / (L_* + L_{env}) = 1 - P_*$$

## MCRT Algorithm

1) pick stellar vs envelope photon

$\xi = \text{rand}(0, 1)$     if  $\xi < P_*$  then stellar photon  
else envelope photon

2) pick starting position, frequency, and direction

a) Star

i) position,  $\vec{x}_0$

$$\frac{dP}{dA} = 4\pi H / L_*$$

note:  $H$  might be position-dependent  
(e.g., gravity darkening)

ii) Frequency,  $\nu_0$

$$\frac{dP}{d\nu} = H_\nu / H$$

iii) Direction,  $\hat{n}$

$$\frac{dP}{d\Omega} = \mu I_\nu / (4\pi H_\nu)$$

b) Envelope

i) position

$$\frac{dP}{dV} = 4\pi \int_0^\infty J_\nu d\nu / L_{\text{env}}$$

ii) frequency

$$\frac{dP}{d\nu} = J_\nu / \int_0^\infty J_\nu d\nu$$

iii) direction

$$\frac{dP}{d\Omega} = \frac{1}{4\pi}$$

3) Doppler Shift photon frequency for source motion

$$\nu \rightarrow (1 + \hat{n} \cdot \vec{v}/c) \nu_0$$

4) Transport photon along straight line to interaction location

since  $\frac{dI_\nu}{d\tau} = -\chi_\nu \rho I_\nu$ , the probability for interaction is

$$dP = d\tau = \chi_\nu \rho ds$$

the number of photons interacting at a point is

$$N \xrightarrow{ds} N + dN$$

$$dN = -N d\tau$$

where  $N =$  number passing that point

$$\Rightarrow N = N_0 e^{-\tau}$$

the cumulative probability of interaction is

$$P = \frac{N - N_0}{N_0} = 1 - e^{-\tau} \quad (\text{Poisson distribution})$$

generate uniform random number

$$\xi = \text{rand}(0, 1) = 1 - e^{-\tau}$$

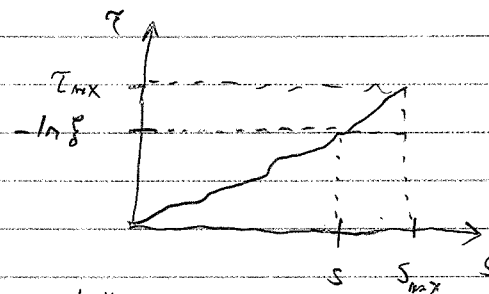
$$\tau = -\ln(1 - \xi)$$

but  $1 - \xi$  is also uniformly distributed  $(0, 1)$

so equivalently

$$\tau = -\ln \xi$$

$$\int_0^s \chi_\nu \rho(\vec{x}_0 + s' \hat{n}) ds' = -\ln \xi$$



solve for  $s$ . This is computationally challenging.  
Analytic soln is faster than numeric.

maximum value for  $s$ :

- i) photon escape  $\Rightarrow$  place photon in observation bin
- ii) photon hits star  $\Rightarrow$  reemit photon (locally) from stellar surface

otherwise, move photon to interaction location

$$\vec{x} = \vec{x}_0 + s\hat{n}$$

(Note: cartesian coords are usually used for this)

5) Find interaction process

The albedo is the scattering probability

$$a = \frac{\sigma_s}{\sigma_s + \kappa_{\nu}} = P_{\text{scatter}}$$

generate random number  $\xi = \text{rand}(0, 1)$

$\xi > a \Rightarrow$  absorb (i.e. destroy photon)

$\xi < a \Rightarrow$  scatter

6) Scatter photon

- i) Doppler shift into rest frame of scatterer
- ii) pick new direction

$$\frac{dP}{d\Omega} = \frac{1}{\sigma_s} \left( \frac{d\sigma_s}{d\Omega} \right) = g(\hat{n}_{in}, \hat{n}_{out})$$

- iii) Doppler shift back to lab frame  
update photon weight

7) repeat  $10^6 - 10^9$  times



## Weighted (biased) Monte Carlo

- Suppose you want to send more photons to undersampled regions (e.g. interior of a dense cloud)

Answer: i) change scattering/emission probability distributions to send more photons in desired direction (i.e., bias the distribution functions).

ii) change photon weight by ratio of probabilities

$$w = \left( \frac{P_{\text{normal}}}{P_{\text{biased}}} \right) w$$

## Making Measurements

### • Sample Events:

- i) photon escape
- ii) photon interactions: absorption  
scattering
- iii) cell/boundary crossing
- iv) photon motion

### • Count Samples

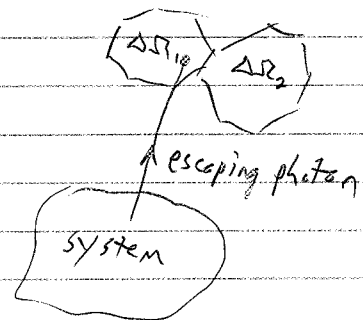
$$\text{energy} = \sum_i W_i E_i$$

### • Convert Energy / Counts to desired measurement

#### Examples

##### i) Emergent Flux

- divide emergent directions into solid angle bins,  $\Delta\Omega_i$
- count # in each bin



$$N_i = \sum_{\text{escapes into } i} W$$

- For an observer at a distance  $d$ , the flux

$$F_i = \frac{dE}{dAdt} = \frac{N_i E_i}{(d^2 \Delta\Omega_i) \Delta t} \quad \text{but } E_i = \frac{L \Delta t}{N_i}$$
$$= \left( \frac{L}{d^2 \Delta\Omega_i} \right) \frac{N_i}{N_i}$$

normalizing to the stellar flux  $F_* = \frac{L}{4\pi d^2}$

$$\left( \frac{F_i}{F} \right) = \left( \frac{4\pi}{\Delta\Omega_i} \right) \frac{N_i}{N_i}$$

- To produce an SED / spectrum, bin in frequency and direction

$$\frac{F_\nu}{F} = \left( \frac{4\pi}{\Delta\Omega_i \Delta\nu_j} \right) \frac{N_{ij}}{N_\nu}$$

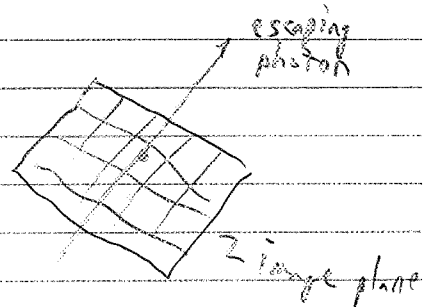
- To produce an image, project escaping photon into pixel on an image plane.

Let  $k, l$  be the pixel indices,

The specific intensity

$$I_\nu = \frac{dE/dt}{\Delta A_i \Delta\nu_j \Delta\Omega_{kl}}$$

$$= \frac{N_{ijkl} (E_\nu / \Delta t)}{d^2 \Delta\Omega_i \Delta\nu_j (A_{kl} / d^2)}$$



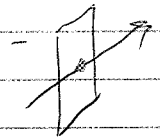
$$= \frac{L}{\Delta\Omega_i \Delta\nu_j A_{kl}} \left( \frac{N_{ijkl}}{N_\nu} \right)$$

$$= \frac{L}{4\pi d^2} \left( \frac{4\pi d^2}{\Delta\Omega_i \Delta\nu_j A_{kl}} \right) \left( \frac{N_{ijkl}}{N_\nu} \right)$$

$$\therefore \frac{I_\nu}{F_\nu} = \left( \frac{4\pi}{\Delta\Omega_i \Delta\nu_j \Delta\Omega_{kl}} \right) \frac{N_{ijkl}}{N_\nu}$$

## ii) Flux crossing a boundary

- construct a loop of area  $\Delta A_i$
- count net number of photons crossing loop



$$N_i = \sum_{\text{outward crossings}} W - \sum_{\text{inward crossings}} W$$

Flux normalized to stellar surface flux

$$\frac{F_i}{F_*} = \left( \frac{4\pi R^2}{L} \right) \frac{N_i (E_\nu / \Delta t)}{A_i} = \left( \frac{4\pi R^2}{A_i} \right) \frac{N_i}{N_\nu}$$

• Intensity moments

$$J/F_* = \frac{R^2}{A_i} \sum (w/\mu)$$

$$H/F_* = \frac{R^2}{A_i} (\sum w^+ - \sum w^-)$$

$$K/F_* = \frac{R^2}{A_i} \sum (\mu w)$$

don't do this  
← it has bad  
statistical properties

### Problems with Simple Sampling

• The relative error for unweighted photons

$$S/N \sim \frac{1}{\sqrt{N_i}}$$

• oftentimes the number of samples is relatively small

$$N_i \sim \frac{N_{\text{bins}}}{N_{\text{y}}}$$

$$\text{for images } N_{\text{bins}} = N_{\text{freq}} \times N_{\text{directions}} \times N_{\text{pixels}}$$

$$= 10^2 \cdot 10 \cdot 10^4 = 10^7$$

$$\Rightarrow S/N = 100 \text{ require } 10^7 \text{ photons.}$$

## Advanced Sampling Techniques

• Source Function Sampling (also called peeling off)

i) sample photon interactions to measure incoming energy

ii) pick a set of observation directions

iii) calculate fraction of incoming energy escapes to each observer

a) photon emitted by star

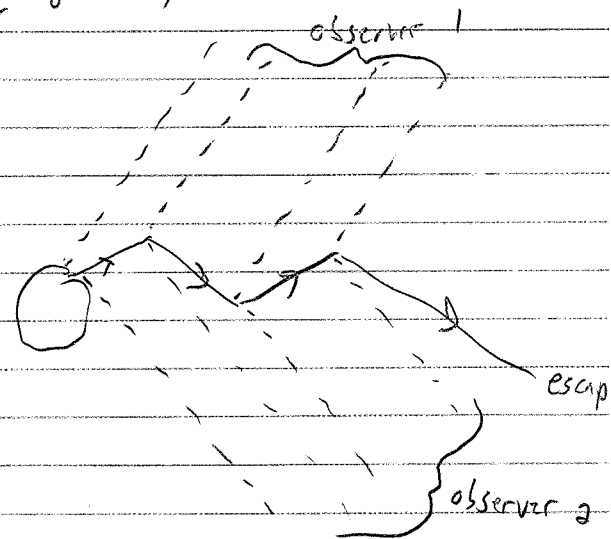
$$N_i = \int \frac{4\pi I_\nu(\mu)}{4\pi H_\nu} e^{-\tau_{esc}} \Delta\Omega_i$$

b) photons emitted in envelope

$$N_i = \int \frac{W}{4\pi} e^{-\tau_{esc}} \Delta\Omega_i$$

c) photons scattered in envelope

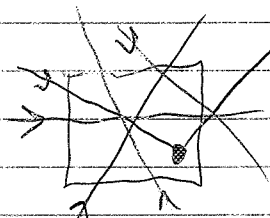
$$N_i = \int W g(\hat{n}_i, \hat{n}_{obs}) e^{-\tau_{esc}} \Delta\Omega_i$$



• Lucy Path Length Sampling

for optically thin cells very few photons interact but many photons pass through the cell

trick is to instead sample virtual absorptions using photon motion



$$\begin{aligned} N_{abs} &= \int d\tau \\ &= \int K \nu p ds = K \nu p \int ds \end{aligned}$$

## Error Estimation

### A) Equal Energy (unweighted) Photons

- For a fixed number of photons, an escaping photon is either placed in a given bin of interest or it is not.
- The probability distribution is a binomial distribution
- The relative error

$$\frac{SE}{E} = \sqrt{(1 - N_i/N_k)/N_i}$$

$$\approx 1/\sqrt{N_i} \quad (N_i \ll N_k)$$

### B) Weighted Photons

- Each entire photon track is a statistically independent event
- Let

$$W_i = \sum_{\text{track history}} W_{\text{obs}} \quad (E_{\text{obs}} = \sum_i W_i E_i)$$

be the weight contributed by the  $i^{\text{th}}$  photon track

- The relative error is approximately

$$\frac{SE}{E} \approx \sqrt{\sum_i W_i^2} / \sum_i W_i$$

Problem: Suppose one particular weight,  $W_i^{\text{max}}$ , is much larger than the rest

$$\Rightarrow \begin{cases} \sum_i W_i = W_i^{\text{max}} \\ \sum_i W_i^2 = (W_i^{\text{max}})^2 \end{cases}$$

$$\therefore \frac{SE}{E} \rightarrow 1 \quad !!!$$

and the measurement is useless

Moral: When using weighted photons, make sure all weights contributing to a measurement all are roughly equal in magnitude.

Theorem: The minimum measurement error occurs when all photons have equal weights.

Unless you have no choice, always use unweighted photons

## Radiative Equilibrium

• The circumstellar envelope is heated by the stellar radiation

⇒ Envelope emission is powered by stellar photons

Note: there may be additional energy sources, e.g. gravitational potential energy in accretion

• Radiative Equilibrium

energy absorbed = energy emitted

This is both a global and local condition

• MC Radiative Equilibrium

i) whenever a photon is absorbed reemit it locally

ii) reduce  $k_{\text{env}}$  to include only energy from other sources (e.g. accretion)

⇒ only photons generated in envelope are from other sources

iii) choose frequency of reemitted photon from local emissivity

$$\frac{dP}{d\lambda} \propto j_{\nu}$$

iv) outgoing direction is isotropic

v) doppler shift as necessary

Note: 1) Energy is automatically and explicitly conserved

2) Since no energy is lost, lambda iteration converges (lower boundary necessarily communicates with outer boundary)



• Radiative Equilibrium Temperature of Grid Cell

$$E_{abs} = E_{emit}$$

$$N_{abs} E_{\nu} / \Delta T = \int 4\pi j_{\nu} dV d\nu \quad \text{but} \quad j_{\nu} = K_{\nu} P_{\nu}(T_{cell})$$

$$= 4\pi \int \rho dV \int_0^{\infty} K_{\nu} B_{\nu}(T_{cell}) d\nu$$

$$= 4\pi m_{cell} K_p B(T_{cell})$$

$K_p =$  Planck mean opacity

note:  $K_p$  is func of  $T_{cell}$

$$= 4\pi m_{cell} K_p \left( \frac{\sigma T_{cell}^4}{\pi} \right)$$

$$\sigma T_{cell}^4 = \frac{N_{abs} (E_{\nu} / \Delta T)}{4 m_{cell} K_p(T_{cell})}$$

$$\therefore \sigma T_{cell}^4 = \frac{L}{4 m_{cell} K_p(T)} \left( \frac{N_{abs}}{W_{\nu}} \right)$$

• How to measure  $N_{abs}$

i) sample photon absorptions  $N_{abs} = \sum W$

ii) use Lucy path length sampling

$$N_{abs} = \sum W d\tau_{abs}$$

$$= \sum W K_{\nu} \rho ds$$

$$= \rho \sum W K_{\nu} ds$$

$$\sigma T_{cell}^4 = \frac{L \sum W K_{\nu} ds}{4 V_{cell} K_p(T_{cell}) W_{\nu}}$$

Note: typically path length sampling has much smaller random measurement errors than counting absorptions, especially for optically thin cells

## Temperature Correction

- In radiative equilibrium photon is emitted locally using  $j_\nu$  to pick the frequency

Problem:  $j_\nu = j_\nu(T_{\text{cell}})$   
and we do not know  $T_{\text{cell}}$  a priori

- Solution:
- 1) guess  $T$  and iterate
  - 2) use instantaneous cell temperature and choose frequency to correct the frequency distribution (Bjorkum + Wood)

- Consider a single photon absorption event

$$T \rightarrow T + \Delta T$$

$$\begin{aligned} j_\nu &\rightarrow j_\nu(T + \Delta T) = j_\nu(T) + \Delta T \frac{dj_\nu}{dT} \\ &= K_\nu B_\nu(T) + \Delta T K_\nu \frac{dB_\nu}{dT} \end{aligned}$$

The change in the frequency distribution is

$$\Delta j_\nu = \Delta T K_\nu \frac{dB_\nu}{dT}$$

- Remitt the photon choosing its frequency from

$$\frac{dP}{d\nu} = K_\nu \frac{dB_\nu}{dT} / \int_0^\infty K_\nu \frac{dB_\nu}{dT} d\nu$$

- Notes:
- 1) Iteration no longer required
  - 2) This is an example of adaptive Monte Carlo
  - 3) It works by reordering which frequencies are emitted first and by "filling in" the missing ones
  - 4) it is not easily parallelized

Generally, Lucy path length sampling + iteration (with perhaps Bjorkum + Wood for the 1<sup>st</sup> iteration) is the method of choice.