# **Rendering: Monte Carlo Integration II**

#### **Bernhard Kerbl**

# **Research Division of Computer Graphics** Institute of Visual Computing & Human-Centered Technology TU Wien, Austria

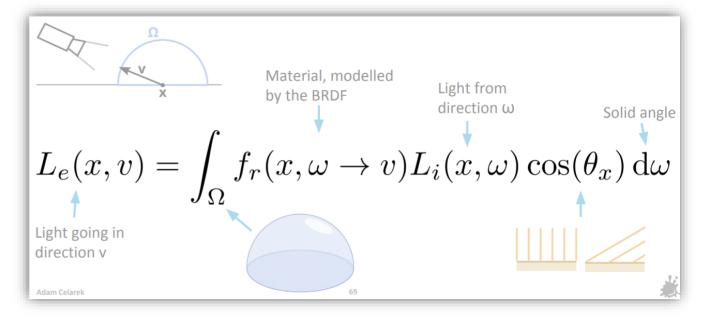
With slides based on material by Jaakko Lehtinen, used with permission

#### Today's Goal



#### Integrating the cosine-weighted radiance $L_i(x, \omega)$ at a point x

Integral of the light function
 over the hemisphere, w.r.t.
 direction/solid angle at ω



- Let's find a solution!
  - How do we integrate over the hemisphere?
  - How do we do it **smartly**?



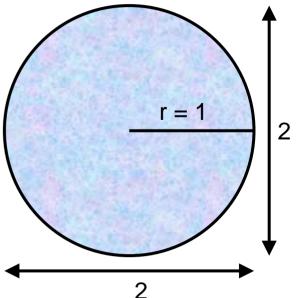


Imagine we have a disk-shaped surface with radius r = 1 that registers incoming light (color) from directional light sources

As an exercise, we want to approximate the total incoming light over the disk's surface area

• We integrate over an area of size  $\pi$ 

We will use the Monte Carlo integral for that

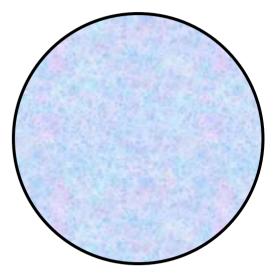




If we can manage to uniformly sample the disk, then we can compute the Monte Carlo integral as a simple average  $\times \pi$ 

By drawing uniform samples in x and y, we cannot cover the area precisely

Inscribed square: information lost



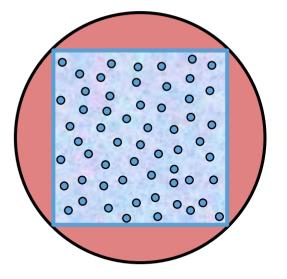
Circumscribed square: unnecessary samples



If we can manage to uniformly sample the disk, then we can compute the Monte Carlo integral as a simple average

By drawing uniform samples in x and y, we cannot cover the area precisely

Inscribed square: information lost



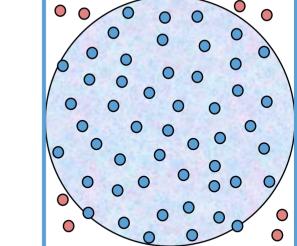
Circumscribed square: unnecessary samples



we cannot cover the area precisely

Uniformly Sampling the Unit Disk

Inscribed square: information lost



Circumscribed square: unnecessary samples

If we can manage to uniformly sample the disk, then we can

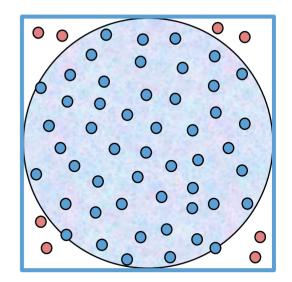
compute the Monte Carlo integral as a simple average



If we can manage to uniformly sample the disk, then we can compute the Monte Carlo integral as a simple average

By drawing uniform samples in x and y, we cannot cover the area precisely

Inscribed square: information lost



Circumscribed square: unnecessary samples
 This is actually somewhat ok!

Rendering – Monte Carlo Integration II



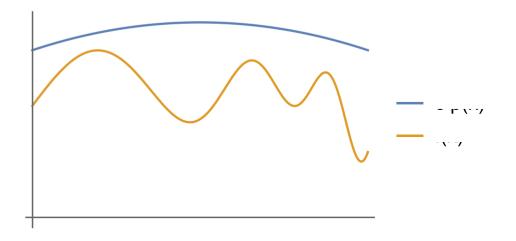




Requires a PDF  $p_X(x)$  and a constant c such that  $f(x) < cp_X(x)$ 

Draw  $\xi_i$  and  $X_i$  from their respective distributions. If the point  $(X_i, \xi_i c p_X(X_i))$  lies under f(x), then the sample is accepted

loop forever: sample X from  $p_X$ 's distribution if  $\xi_i \cdot cp_X(X_i) < f(X_i)$  then return  $X_i$ 

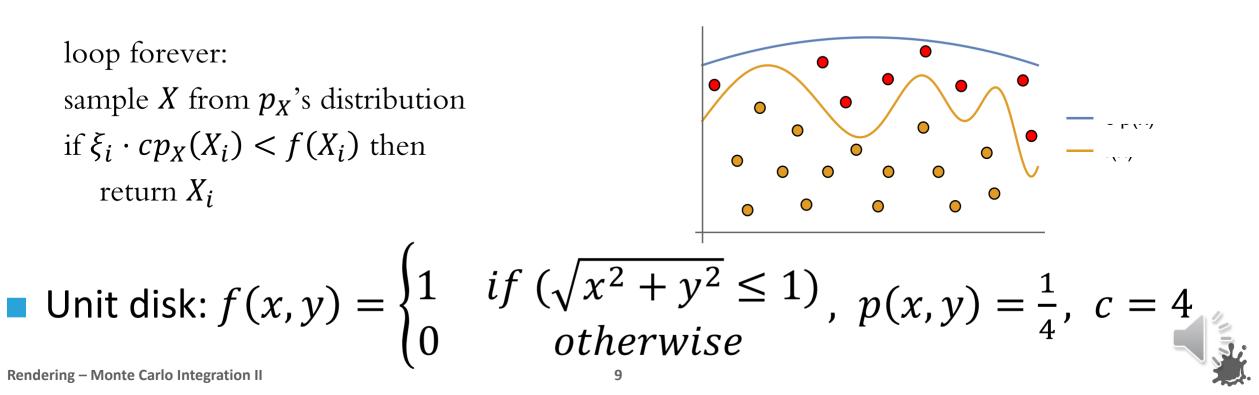






• Requires a PDF  $p_X(x)$  and a constant c such that  $f(x) < cp_X(x)$ 

Draw  $\xi_i$  and  $X_i$  from their respective distributions. If the point  $(X_i, \xi_i c p_X(X_i))$  lies under f(x), then the sample is accepted





We do not want to waste samples if we can avoid it

Instead, find a way to generate uniform samples on the disk

- Second attempt: draw from 2D polar coordinates
  - Polar coordinates defined by radius  $r \in [0,1)$  and angle  $\theta \in [0,2\pi)$
  - Transformation to cartesian coordinates:
    - $x = r\sin\theta$
    - $\mathbf{y} = r\cos\theta$



### Uniformly Sampling the Unit Disk?

# Convert two $\xi$ to ranges [0, 1), $[0, 2\pi)$ for polar coordinates

#### Convert to cartesian coordinates

```
void sampleUnitDisk()
{
      std::default random engine r rand eng(0xdecaf);
      std::default random engine theta rand eng(0xcaffe);
      std::uniform real distribution<double> uniform dist(0.0, 1.0);
      for (int i = 0; i < NUM SAMPLES; i++)</pre>
      {
            auto r = uniform dist(r rand eng);
            auto theta = uniform dist(theta rand eng) * 2 * M PI;
            auto x = r * sin(theta);
            auto y = r * cos(theta);
            samples2D[i] = std::make pair(x, y);
```



# Clumping

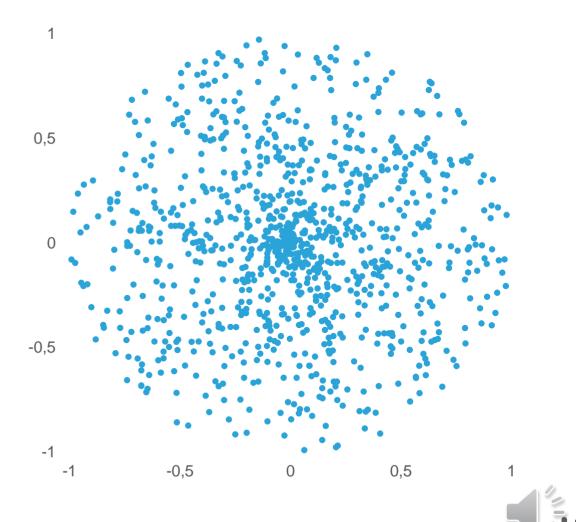


#### We successfully sampled the unit disk in the proper range

However, the distribution is not uniform with respect to the area

Samples clump together at center

Averaging those samples will give us a skewed result for the integral!





The area of a disk is proportional to  $r^2$ , times a constant factor  $\pi$ 

If we see the disk as concentric rings of width  $\Delta r$ , the *j* inner rings up to radius  $r_j = j\Delta r$  should contain  $\left(\frac{r_j}{r}\right)^2 N$  out of *N* total samples

Conversely, the  $i^{th}$  sample should lie in the ring at radius  $r_i = r \sqrt{\frac{i}{N}}$ 

Since 
$$\xi$$
 is uniform in [0, 1), we can switch  $\frac{j}{N}$  for  $\xi$  to get  $r_i = r\sqrt{\xi_i}$ 

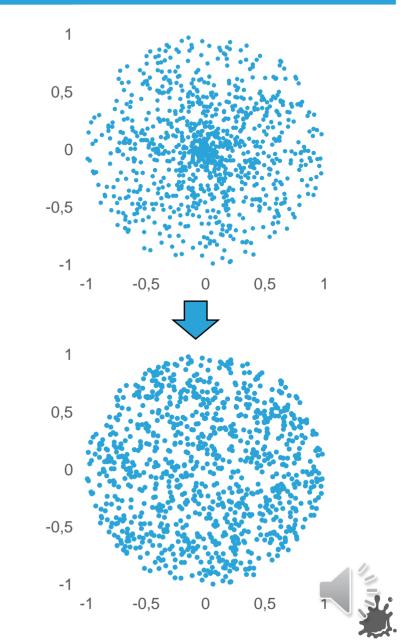
# Uniformly Sampling the Unit Disk: A Solution



It works, and it is not even a bad way to arrive at the correct solution

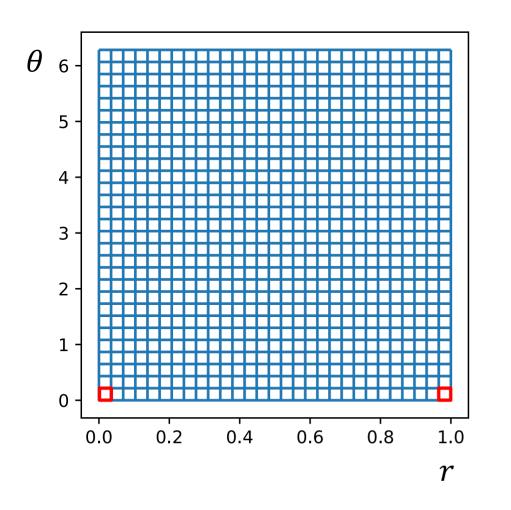
However, for more complex scenarios, we might struggle to find the solution so easily

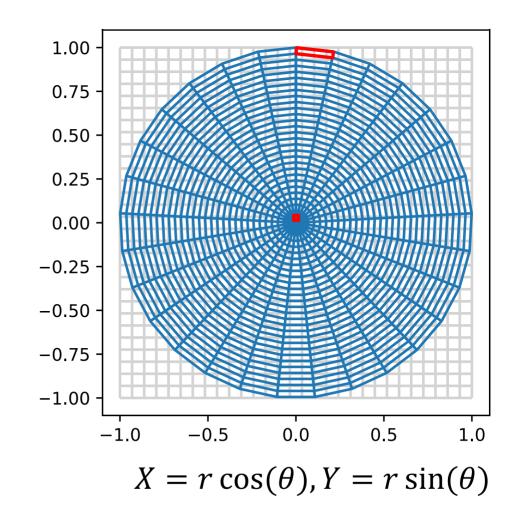
With the tools we introduced earlier, we can formalize this process for arbitrary setups





#### Let's transform a regular grid from polar to cartesian coordinates



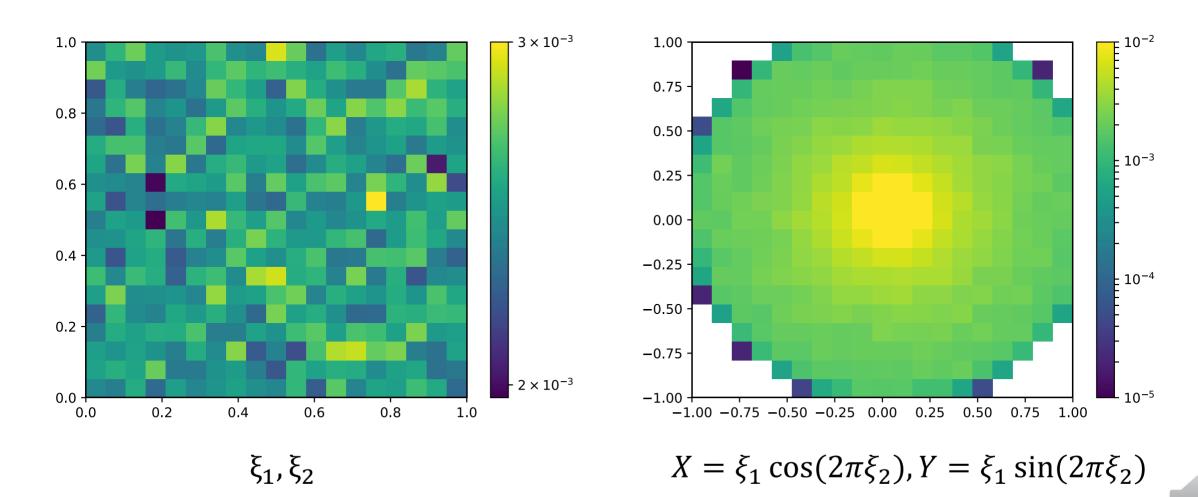




#### First Attempt to Learn the PDF



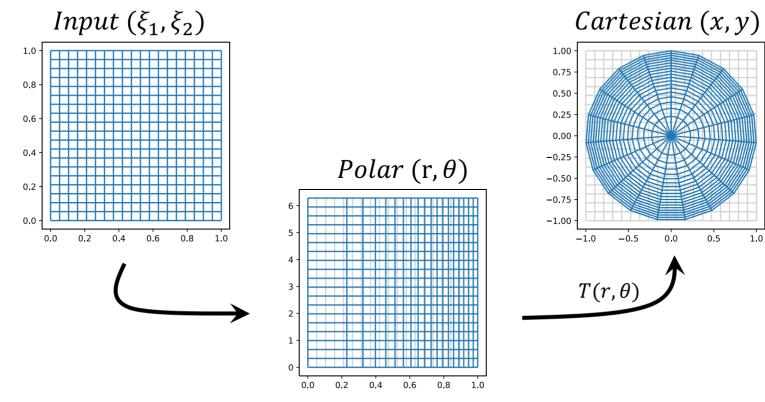
#### Take 100k samples, transform and see which box they end up in





■ If we know the effect of a transformation *T* on the PDF, we can

- Use it in the Monte Carlo integral to weight our samples, or
- Compensate to get a uniform sampling method after transformation





# Computing the PDF after a Transformation

TU

Assume a random variable A and a bijective transformation T that yields another variable B = T(A)

Bijectivity dictates that b = T(a) must be either monotonically increasing or decreasing with a

This implies that there is a unique  $B_i$  for every  $A_i$ , and vice versa

In this case, the CDFs for the two variables fulfill  $P_B(T(a)) = P_A(a)$ 



# Computing the PDF after a Transformation



If 
$$b = T(a)$$
 and  $b$  increases with  $a$ , we have:  $\frac{dP_B(b)}{da} = \frac{dP_A(a)}{da}$   
If  $b$  decreases with  $a$  (e.g.  $b = -a$ ), we have:  $-\frac{dP_B(b)}{da} = \frac{dP_A(a)}{da}$ 

Since  $p_B$  is the non-negative derivative of  $P_B$ , we can rewrite as:

$$p_B(b) \left| \frac{db}{da} \right| = p_A(a), \qquad U_{sing:} \frac{dP_X(x)}{dy} = \frac{p_X(x) dx}{dy}$$
$$p_B(b) = \left| \frac{db}{da} \right|^{-1} p_A(a)$$

# Computing the PDF after a Transformation

Let's interpret 
$$p_B(b) = \left|\frac{db}{da}\right|^{-1} p_A(a)$$

It is the probability density of A, multiplied by  $\left|\frac{db}{da}\right|^{-1}$ 

 $\left|\frac{db}{da}\right|^{-1}$  has two intuitive interpretations:

the change in sampling density at point a if we transform a by Tor, the inverse change in the volume of an infinitesimal hypercube at point a if we transform a by T



If we try to apply the above to the unit disk, we fail at  $x = r \sin \theta$ 

# • We can't evaluate $\left|\frac{dx}{dr}\right|^{-1}$ : the transformation that produces one target variable is dependent on both input variables and vice-versa

We cannot compute the change in the PDF between individual variables, we must take them all into account simultaneously



Rendering – Monte Carlo Integration II



We write the set of N values from a **multidimensional** variable  $\vec{A}$  as a vector  $\vec{a}$  and the N outputs of transformation T as a vector  $\vec{b}$ :

$$\vec{a} = \begin{pmatrix} a_1 \\ \vdots \\ a_N \end{pmatrix}, \vec{b} = \begin{pmatrix} b_1 \\ \vdots \\ b_N \end{pmatrix} = \begin{pmatrix} T_1(\vec{a}) \\ \vdots \\ T_N(\vec{a}) \end{pmatrix} = T(\vec{a})$$

Instead of quantifying the change in volume incurred by T(a),  $\left|\frac{dT(a)}{da}\right|$ , our goal is now to quantify the change incurred by  $T(\vec{a})$ 





For a transformation  $\vec{b} = T(\vec{a})$ , we can define the Jacobian matrix that contains all  $b_j$ ,  $a_i$  combinations of partial differentials

$$J_T(\vec{a}) = \begin{pmatrix} \frac{\partial b_1}{\partial a_1} & \cdots & \frac{\partial b_1}{\partial a_N} \\ \vdots & \ddots & \vdots \\ \frac{\partial b_M}{\partial a_1} & \cdots & \frac{\partial b_M}{\partial a_N} \end{pmatrix}$$

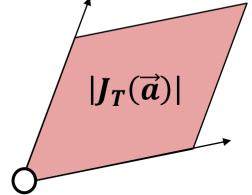
If we consider  $\vec{A}$ 's domain as a space with N axes,  $J_T(\vec{a})$  gives the change of the edges of an infinitesimal hypercube from  $\vec{a}$  to  $T(\vec{a})$ 

#### ■ Change of edges of an infinitesimal hypercube with extent (1,1)



The columns of a square matrix M can be interpreted as the natural base vectors of a space  $\begin{pmatrix} 1\\0\\\vdots\\0 \end{pmatrix}$ ,  $\begin{pmatrix} 0\\1\\\vdots\\0 \end{pmatrix}$  if they were transformed by M

The determinant |M| of M computes the volume of a parallelepiped spanned by these vectors<sup>[3]</sup>



 $|J_T|$ , called the Jacobian of T, gives the volume change at  $\vec{a}$  by T



# Computing the PDF of a Transformation

• Let's try polar coordinates again: 
$$\binom{x}{y} = T\binom{r}{\theta} = \binom{r\sin\theta}{r\cos\theta}$$

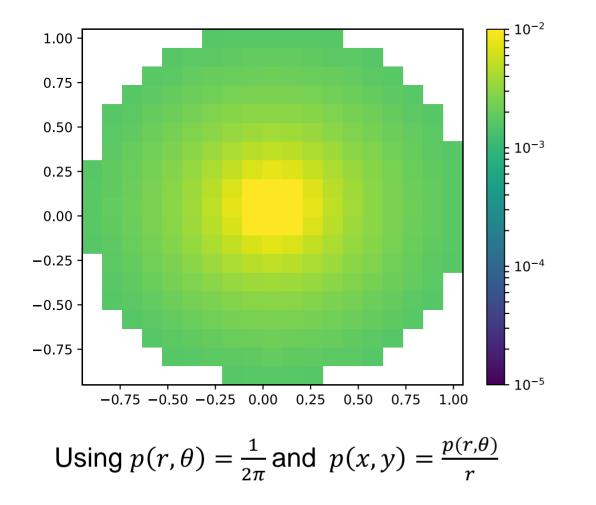
$$\left|\frac{\partial T\binom{r}{\theta}}{\partial \binom{r}{\theta}}\right| = |J_T| = \left|\begin{pmatrix}\frac{\partial x}{\partial r} & \frac{\partial x}{\partial \theta}\\ \frac{\partial y}{\partial r} & \frac{\partial y}{\partial \theta}\end{pmatrix}\right| = \left|\begin{pmatrix}\cos\theta & -r\sin\theta\\\sin\theta & r\cos\theta\end{pmatrix}\right| = r$$

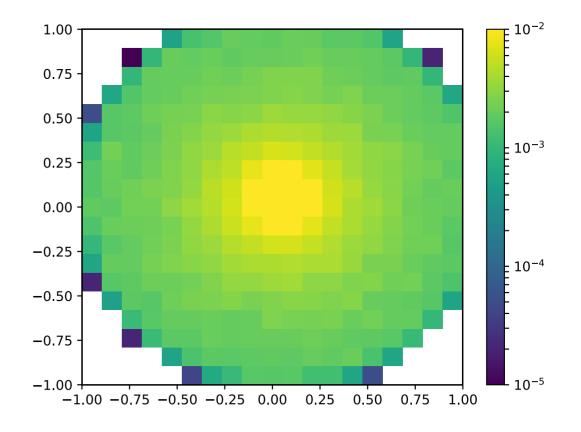
$$p(x,y) = \frac{p(r,\theta)}{r}, \text{ or } p(r,\theta) = r p(x,y):$$
  
a uniform density in  $(x,y)$  must be proportional to  $r$  in  $(r,\theta)$ 



#### **Compare PDFs After Transformation**







Measured



#### Python code, for the interested

```
num_bins = 20
```

```
xlist = np.linspace(-1.0, 1.0, num_bins)
ylist = np.linspace(-1.0, 1.0, num_bins)
bin_volume = 4 / (num_bins*num_bins)
```

```
X, Y = np.meshgrid(xlist, ylist)
X += 1/num_bins
Y += 1/num bins
```

```
r = np.sqrt(X*X+Y*Y)
uniform_dist = 1/(2 * np.pi)
pdf transform = 1/r
```

```
Z = uniform_dist * pdf_transform * bin_volume
Z[r>1] = 0
fig,ax = plt.subplots(1,1)
cp = plt.pcolor(X, Y, Z,cmap='viridis',norm=matplotlib.colors.LogNorm())
fig.colorbar(cp) # Add a colorbar to a plot
plt.show()
```



For independent variables, the joint PDF p(x, y, ...) is  $p_X(x)p_Y(y)$  ...

In general, this is an assumption that we should not rely on

Furthermore, after a transformation, only the joint PDF is known

The proper way to sample multiple variables X, Y is to compute
 the marginal density function p<sub>X</sub>(x) of one
 the conditional density function p<sub>Y</sub>(y|x) of the other



# Marginal and Conditional Density Function

- Assume we have obtained the joint PDF p(x, y) of variables X, Ywith ranges  $[a_X, b_X)$  and  $[a_Y, b_Y)$
- In a 2D domain with X, Y we can think of  $p_X(x)$  as the average density of p(x, y) at a given x over all possible values y
- We can obtain the marginal density function for one of them by integrating out all the others, e.g.:  $p_X(x) = \int_{a_y}^{b_Y} p(x, y) dy$

• We can then find 
$$p(y|x) = \frac{p(x,y)}{p_X(x)}$$

#### **Adding More Variables**



- What to do for multiple variables, e.g. *X*, *Y* and *Z*?
  - Find first marginal density  $p_X(x) = \int_{aZ}^{bZ} \int_{aY}^{bY} p(x, y, z) \, dy \, dz$
  - Find first conditional density  $p_X(y, z|x) = \frac{p(x, y, z)}{p_X(x)}$
  - Find second marginal density  $p_Y(y|x) = \int_{aZ}^{bZ} p(x, y, z) dz$
  - Find second conditional density  $p_X(z|x, y) = \frac{p(y, z|x)}{p_Y(y|x)}$
  - Integrate + invert first marginal, first and second conditional densities
  - Sample each of them
  - Extend ad libitum to even more variables





The size of the sampling domain in cartesian coordinates is  $\pi$ 

Since we want uniform sampling and sample probabilities must integrate to 1, the PDF in cartesian coordinates is  $p(x, y) = \frac{1}{\pi}$ 

We know that 
$$p(r,\theta) = r p(x,y)$$
, so we want  $p(r,\theta) = \frac{r}{\pi}$ 

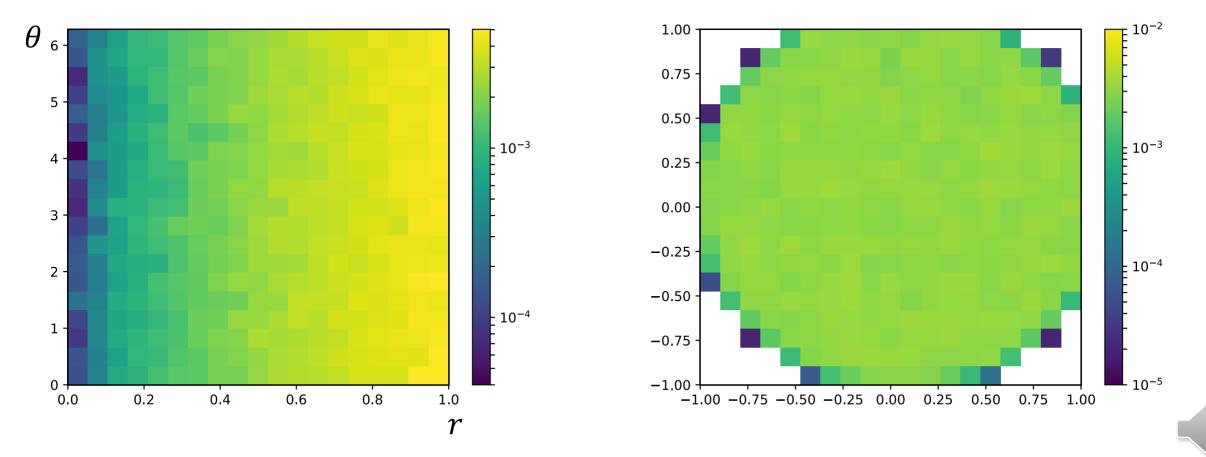
• 
$$p_R(r) = \int_0^{2\pi} p(r,\theta) d\theta = 2r$$
 and  $p(\theta|r) = \frac{p(r,\theta)}{p_R(r)} = \frac{1}{2\pi}$ 



# Sampling the Unit Disk: The Formal Solution



If we draw samples for our  $r, \theta$  with the above PDFs, we get uniform distribution in (x, y) after applying transformation T



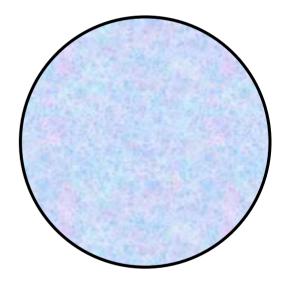
**Rendering – Monte Carlo Integration II** 

# Sampling the Unit Disk: The Formal Solution



Integrate marginal and conditional PDFs and invert: we get the same solution as before:

• 
$$r = P_R^{-1}(\xi_1) = \sqrt{\xi_1}$$
  
•  $\theta = P_{\Theta}^{-1}(\xi_2) = 2\pi\xi_2$ 



•  $p(\theta|r)$  is constant: no matter what radius we are looking at, all positions on a ring of that radius (angle) should be equally likely

Final integral: 
$$RGB_{total} = \frac{\pi}{N} \sum_{i=1}^{N} RGB(R_i \sin \Theta_i, R_i \cos \Theta_i)$$





This took as a while, but we have seen all the formal procedures

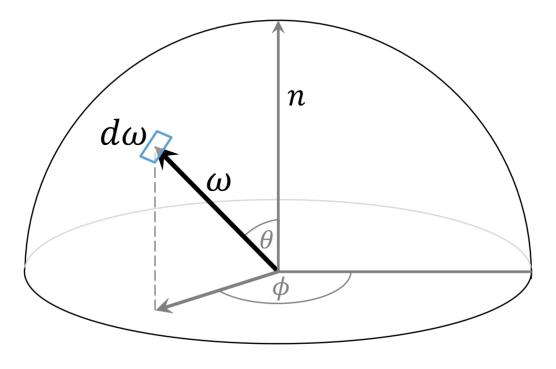
• We only need to switch from integrating planar area to points  $\omega$  on hemisphere surface (i.e., vectors (x, y, z) with length 1)

• Use spherical coordinates and bijective *T* from  $(r, \theta, \phi)$  to (x, y, z):  $x = r \sin \theta \cos \phi$   $y = r \sin \theta \sin \phi$  $z = r \cos \theta$ 



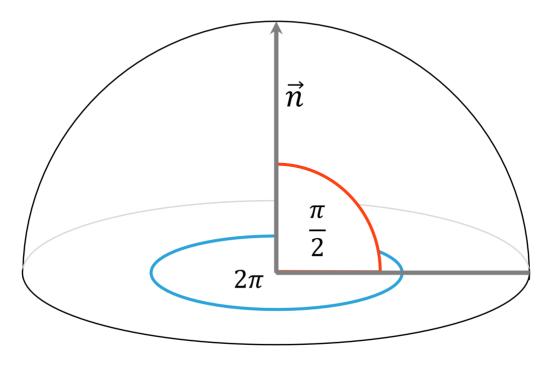
#### **Deriving Integration Over Hemisphere**

- TU
- Each direction  $\omega$  represents an infinitesimal surface area portion  $d\omega$
- How do we integrate a function  $f(\omega)$  with differential  $d\omega$ ?
- Integration over points on hemisphere surface  $\omega$ , w.r.t.  $(\theta, \phi)$





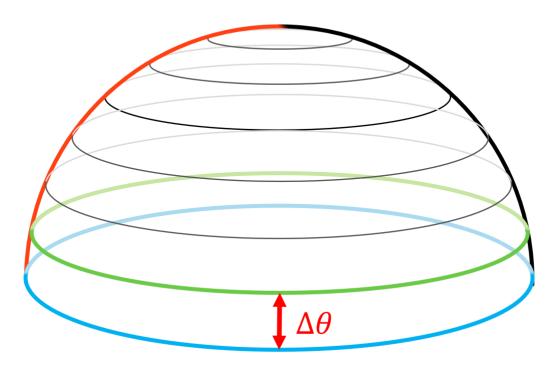
- We assume a planar surface with an upright facing normal *n*
- We use the integral intervals  $\theta \in \left[0, \frac{\pi}{2}\right)$ ,  $\phi \in \left[0, 2\pi\right)$
- I.e., a curve from perpendicular to parallel for  $\theta$ , a ring for  $\phi$







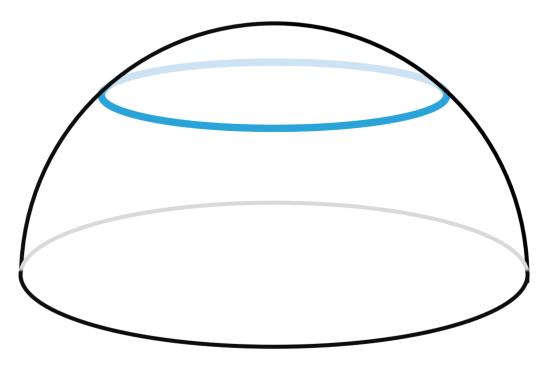
- We can split the surface along  $\theta$  into ribbons of width  $\Delta \theta \rightarrow d\theta$
- The upper edge of the ribbon is slightly shorter than the lower
- If we keep adding more and more ribbons, this difference vanishes







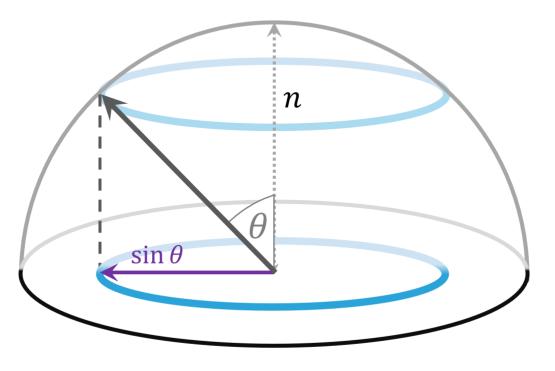
- As a ribbon's width goes to  $d\theta$ , its area becomes its length times  $d\theta$
- We can find this length by projecting the ribbon to the ground
- Using trigonometry, we find the length of a ribbon is  $2\pi \sin \theta$







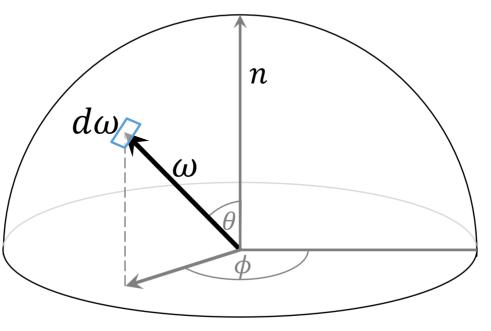
- As a ribbon's width goes to  $d\theta$ , its area becomes its length times  $d\theta$
- We can find this length by projecting the ribbon to the ground
- Using trigonometry, we find the length of a ribbon is  $2\pi \sin \theta$







- The length of a ribbon spans the entire interval  $\phi \in [0, 2\pi)$
- Convert the length to an integral over  $d\phi$ :  $2\pi \sin \theta = \int_0^{2\pi} \sin \theta \, d\phi$
- The final integral:  $\int_{\Omega} f(\omega) d\omega = \int_{0}^{\frac{\pi}{2}} \int_{0}^{2\pi} f(\omega) \sin \theta d\phi d\theta$





# **Deriving PDF for Hemisphere Sampling**



Integral of 
$$f(\omega)$$
 over area  $\Delta \omega = \int_{\Delta \omega} f(\omega) \, d\omega$ 

Integral of  $f(\omega)$  w.r.t.  $(d\theta, d\phi) = \int_{\Delta\theta} \int_{\Delta\phi} f(\omega) \sin\theta \, d\phi \, d\theta$ 

Integration domain and  $f(\omega)$  are identical, thus:  $d\omega = \sin \theta \ d\phi \ d\theta$ 

•  $\omega \to (\theta, \phi)$  is bijective, we have  $p(\theta, \phi) d\theta d\phi = p(\omega) d\omega$  and:

 $p(\theta, \phi) = \sin \theta \, p(\omega)$ 

Target distribution in  $\omega$ , which is (x, y, z) with  $\sqrt{x^2 + y^2 + z^2} = 1$ 

The transformation *T* from  $(r, \theta, \phi)$  to (x, y, z):  $x = r \sin \theta \cos \phi$   $y = r \sin \theta \sin \phi$  $z = r \cos \theta$ 

The Jacobian of the transformation T gives  $|J_T| = r^2 \sin \theta$ 

• Hence, we have  $p(r, \theta, \phi) = r^2 \sin \theta \ p(x, y, z) = 1 \sin \theta \ p(\omega)$ 



# Uniformly Sampling the Unit Hemisphere

The domain, i.e., the unit hemisphere surface area, is  $2\pi$ . Uniformly sampling the domain over  $\omega$  implies  $p(\omega) = \frac{1}{2\pi}$ 

Hence, since 
$$p(\theta, \phi) = r^2 \sin \theta \, p(\omega)$$
, we want  $p(\theta, \phi) = \frac{\sin \theta}{2\pi}$ 

• Marginal density 
$$p_{\Theta}(\theta): \int_{0}^{2\pi} p(\theta, \phi) \, d\phi = \sin \theta$$

Conditional density 
$$p(\phi|\theta): \frac{p(\theta,\phi)}{p_{\Theta}(\theta)} = \frac{1}{2\pi}$$

**Rendering – Monte Carlo Integration II** 



# Uniformly Sampling the Unit Hemisphere – Complete



• Antiderivative of  $p_{\Theta}(\theta)$ :  $\int \sin \theta \ d\theta = 1 - \cos \theta$  (added constant 1)

• Antiderivative of 
$$p(\phi|\theta)$$
:  $\int \frac{1}{2\pi} d\phi = \frac{\phi}{2\pi}$ 

Invert them to get  $\theta = \cos^{-1} \xi_1$  (cos is symmetric),  $\phi = 2\pi\xi_2$ 

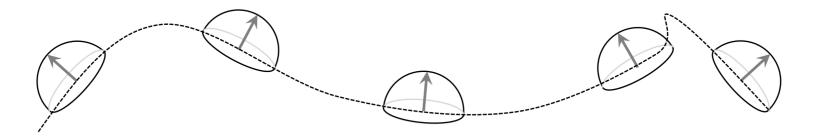
• Apply transformation T on  $(\theta, \phi)$  to obtain uniformly distributed  $\omega$ 







The orientation of sampled hemispheres depends on surface normal



Use the tangent, bitangent and normal vectors of the intersection

•  $\omega_{world} = x \cdot tangent + y \cdot bitangent + z \cdot normal$ 

#### • Or, if available, use *ToWorld* transformation methods

# **Applications for Uniform Hemisphere Sampling**



#### **Diffuse** lighting based on last lecture's insights with constant $f_r$

• What do we use for  $L_i$ ?

 Full rendering equation: next time

TU Apply (from the physics chapter) Material, modelled Light from by the BRDF direction  $\omega$ Solid angle  $L_e(x,v) = \int_{\Omega} f_r(x,\omega \to v) L_i(x,\omega) \cos(\theta_x) \,\mathrm{d}\omega$ Light going in direction v

#### This time: ambient occlusion, direct lighting

#### **Ambient Occlusion**



Consider all unblocked directions around x as indirect light sources, integrate  $V(x, x + \alpha \omega) \frac{\cos \theta}{\pi}$  over directions  $\omega$  around the normal Limit ray length to  $\alpha$ , return free if no intersection closer than  $\alpha$ 

Fine geometric details on objects are accentuated by the absence of ambient light due to the shadows cast by close-by geometry

$$L_e(x) = \int_{\Omega} \mathbf{V}(x, x + \alpha \omega) \frac{\cos(\theta)}{\pi} \,\mathrm{d}\omega$$

Integrate over directions ω on the unit hemisphere defined by point x, normal n

$$V(x, x + \alpha \omega) = \begin{cases} 1 & if \ x \to (x + \alpha \omega) \ free \\ 0 & if \ x \to (x + \alpha \omega) \ blocked \end{cases}$$







Recap: variance is low if the sampling function mimics the signal

• We use  $f_r = \frac{1}{\pi}$  for ambient occlusion, therefore the contribution of signal samples varies mostly with  $\cos \theta$ 

- It would be best to apply importance sampling: use a sampling strategy for  $\omega$ , such that  $p(\omega) \propto \cos \theta$
- We have gone through all the necessary steps.
  <u>Try to solve this formally with the inversion method as an exercise</u>!

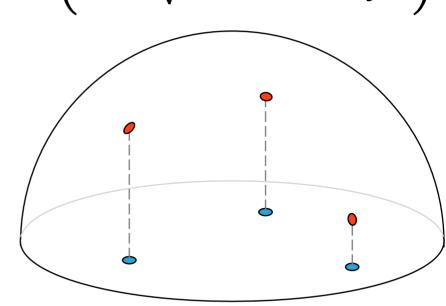


# Smart Cosine-Weighted Importance Sampling

• Malley's method: uniformly pick (x, y) samples on the unit disk

Project them to the hemisphere surface  $\left(z = \sqrt{1 - x^2 - y^2}\right)$ 

Done! Your samples are now distributed with  $p(\omega) \propto \cos \theta$ 



## Why does this work? Try to come up with your own proof!



# Finding Light

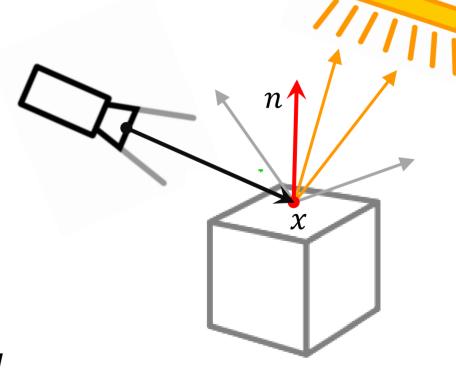


We can use Monte Carlo integration to compute the direct illumination from light sources in the scene at a point x

 Naïve version: sample unit hemisphere uniformly, hoping to hit light sources

Check **closest** hit for each direction  $\omega$ 

Use 
$$L_i(x, \omega) = \begin{cases} L_e^{[l]} & if hit a light l \\ 0 & otherwise \end{cases}$$



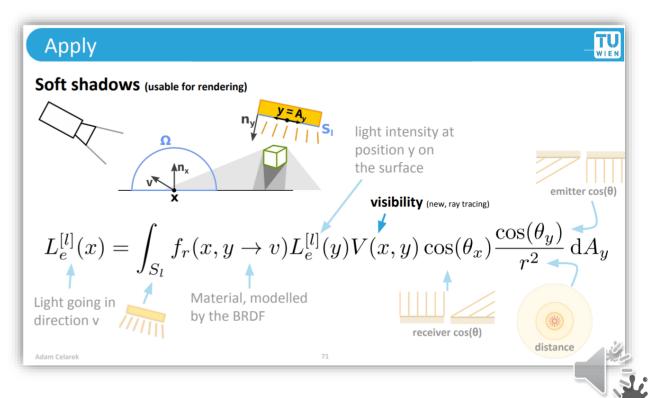


• A special kind of importance sampling: integrate over light sources!

Use  $V(x, y) = \begin{cases} 1 & if path from x to y is unblocked \\ 0 & otherwise \end{cases}$ 

Pick y on light, e.g. uniformly

 $\frac{\cos \theta_y \, dA_y}{r^2}$ : change in volume of infinitesimal 2D hypercube at y projected onto x's hemisphere



# Sampling a Light Source, Revisited



# For inclusion of simple material color and BRDF, use $f_r = \frac{\rho}{\pi}$

#### Works extremely well if the light occupies only a small solid angle



100 samples per pixel, hemisphere sampling



100 samples per pixel, light source sampling



# Hemisphere Sampling Illustrated







# Light Source Sampling Illustrated









But where does the actual integration step happen?

In the basic case, directly in the main sampling loop for each pixel!

- Static scene: Samples through pixels p always hit the same point x
- Once x has been hit, the sampling of its hemisphere follows PDF
- Return sampled values (colors), weighted by the corresponding PDF

■ Use *N* samples for *p*, sum color values weighted by PDF, average:

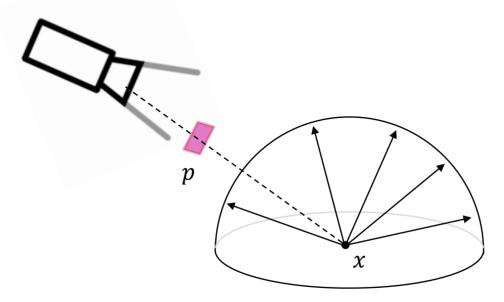
$$F_N = \frac{1}{N} \sum_{i=1}^{N} \frac{L_i(x, \omega)}{p(\omega)}$$



# Monte Carlo Integration as Loop Over Pixel Samples

- We achieve integration around x with multiple samples through p
  - A bit wasteful, but is a general, valid solution
  - We will see in a second why this is convenient

Weight returned values by PDF, sum up and divide by N





# Monte Carlo Integration as Loop Over Pixel Samples

Given: camera, pixel p, scene, pdf

 $rgb = \{0,0,0\}$ for i in [0, N) do ray = rayThroughPixel(camera, p) x = findIntersection(ray, scene)color = getIntegratorValue(x) $rgb += color \bullet$ end for rgb /= Nfunction getIntegratorValue(x) omega = getDirectionOnHemisphere(x, pdf) Li = evaluateLight(x, omega, scene)return Li / pdf(omega) 🔺

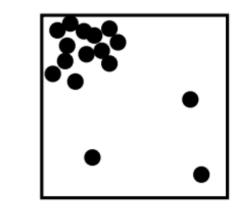
**Rendering – Monte Carlo Integration II** 

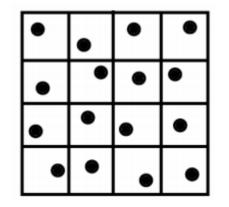


# **Stratified Sampling**

With random uniform sampling, we can get unlucky

- E.g. all samples clump in a corner
- If we don't know anything of the integrand, we want a relatively uniform sampling
  - Not regular, though, because of alias patterns!
- Subdivide domain into non-overlapping regions (e.g. a regular grid). Each region is called a *stratum*











In the first lecture, we used supersampling to fight off aliasing

Pixels are another instance where we use Monte Carlo integration

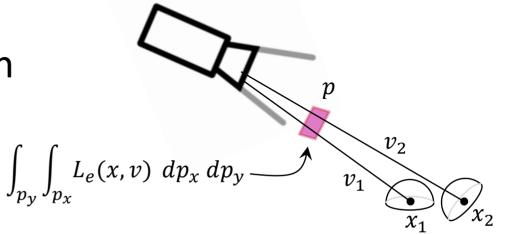
Choosing samples within pixels is an instance of stratified sampling

- Uniform 2D distribution, average over a pixel rectangle: box filter
  - We will see more advanced methods for filtering in future lectures
  - If we didn't use at least one sample per pixel, we would leave holes



# Monte Carlo Integration for Pixels

- Samples are randomly jittered in each stratum
- Ergo, we **don't** hit the same x with each pixel sample (p<sub>x</sub>, p<sub>y</sub>) inside pixel p
- We just add two random variables!



Instead of integrating over a hemisphere, we are integrating over all surface points visible to a pixel and their respective hemispheres

The box filter over pixel color samples implements uniform integral



# The Box Filter for a Pixel (Monte Carlo Integration)



Given: camera, pixel coordinates (pixel\_x, pixel\_y), scene

 $rgb = \{0,0,0\}$ 

for i in [0, N) do

 $px = pixel_x + uniform_random_sample() // \xi_i = P_{\xi}(\xi_i) = P_{\xi}^{-1}(\xi_i)$   $py = pixel_y + uniform_random_sample()$  ray = rayThroughPixelPos(camera, px, py) x = findIntersection(ray, scene) rgb += getIntegratorValue(x)Nothing else necessary: offsets in x and y are uniformly distributed, they are independent, the domain size of each pixel is 1x1=1 and so we don't have to change the integration at all.

end for

rgb /= N





Stratified sampling is a special case of low-discrepancy sampling<sup>[4]</sup>

Replace the built-in RNG with a sample generation algorithm that sacrifices randomness for good spatial distribution

Great for:

- Faster convergence (reduction of noise)
- GPUs (they don't love random numbers)
- Transparent and portable sampling behavior





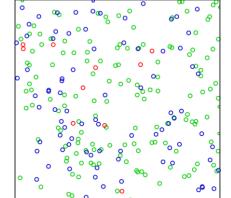
Ô

For *n*-D, pick *n* different, co-prime bases (e.g. 2,3)

Based on *radical inverse*: an integer a can be written with base b and  $d \in [0, b - 1]$  as  $a = \sum_{i=1}^{m} d_i(a)b^{i-1}$ 

For  $b = 2, d \in [0,1]$ : this is binary representation of integers:  $13 = 0 \times 2^0 + 1 \times 2^1 + 1 \times 2^2 + 1 \times 2^3$ 

• Radical inverse with *m* digits:  $\Phi(a) = \sum_{i=1}^{m} \frac{d_i(a)}{h^i}$ 



Default RNG

2,3 Halton Sequence

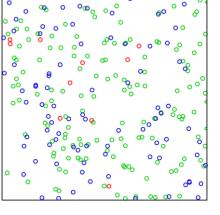
Start at any integer value for each of the *n* variables

- Need data structure to represent  $a = \sum_{i=1}^{m} d_i(a) b^{i-1}$
- Can be written for arbitrary *b* with some bit fiddling

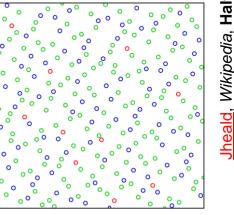
For each new *n*-D sample, increment all *n* integers

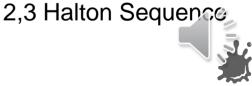
Compute their radical inverse with their proper base

Using same sequence for all pixels  $\rightarrow$  patterns  $\bigcirc$ 



**Default RNG** 







#### Halton Sequences

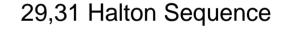
Large primes will behave similarly  $\rightarrow$  patterns (::)

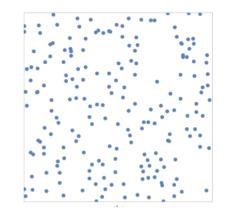
Repetitive patterns should always be avoided!

 Option 1: use different starting values for different instances (e.g. for each pixel)

Option 2: instead of incrementing  $d_i$ , cycle through random permutations of [0, b - 1] in each instance

E.g.: b = 3, {0, 2, 1} instead of {0, 1, 2}





29,31 Halton Scrambled







- Get comfortable with all approaches to integration and sampling
  - The mental image of "area under the curve" eventually collapses (infinite-dimensional integral coming up next!)
  - Transformations between sample domains may be non-trivial
  - Once you grasp the underlying concepts, applying the math is easy

- We have seen simpler explanations for the most important parts
  - Uniformly sampling a hemisphere
  - Cosine-weighted sampling of a hemisphere



## **References and Further Reading**

- Slide set based mostly on chapter 13 of *Physically Based Rendering: From Theory to Implementation*
- [1] Steven Strogatz, Infinite Powers: How Calculus Reveals the Secrets of the Universe
- [2] Video, Why "probability of 0" does not mean "impossible" | Probabilities of probabilities, part 2: <u>https://www.youtube.com/watch?v=ZA4JkHKZM50</u>
- [3] Video, The determinant | Essence of linear algebra, chapter 6: <u>https://www.youtube.com/watch?v=lp3X9LOh2dk</u>
- [4] SIGGRAPH 2012 Course: Advanced (Quasi-) Monte Carlo Methods for Image Synthesis, <u>https://sites.google.com/site/qmcrendering/</u>
- [5] Wikipedia, Van der Corput Sequence, <u>https://en.wikipedia.org/wiki/Van\_der\_Corput\_sequence</u>

