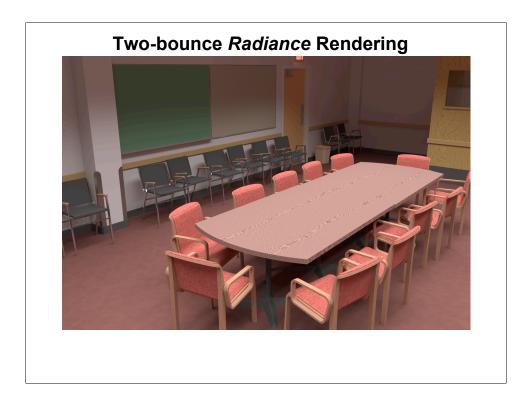


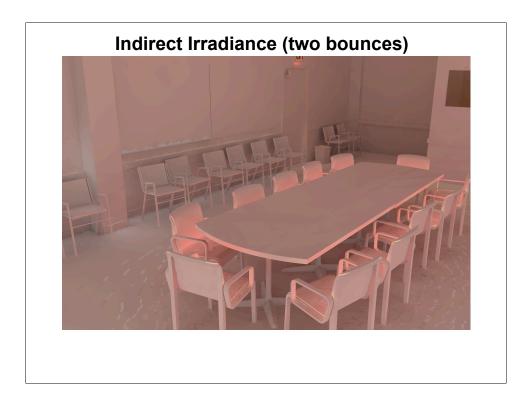
Spatial Coherence of Indirect

- In nearly all cases, indirect illumination changes slowly over surfaces
- Cost of indirect sampling may therefore be reduced through interpolation

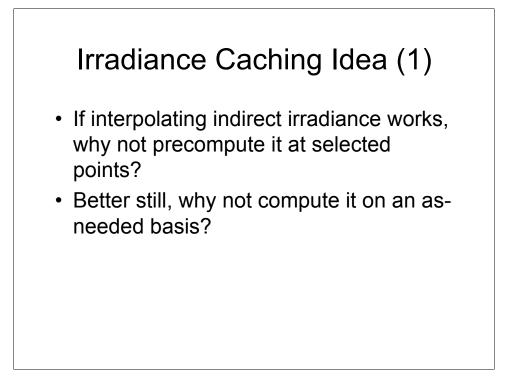
Radiosity techniques work for the same reason -- and have trouble with point light sources because direct illumination does *not* change slowly over surfaces.



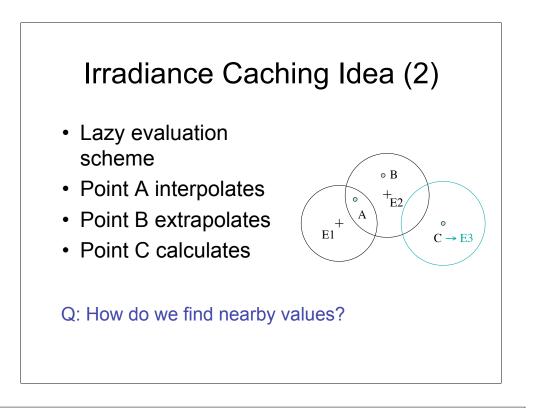
Radiance is a physically-based renderer that has been around for over 20 years, and the irradiance cache was in the first public release in 1989.



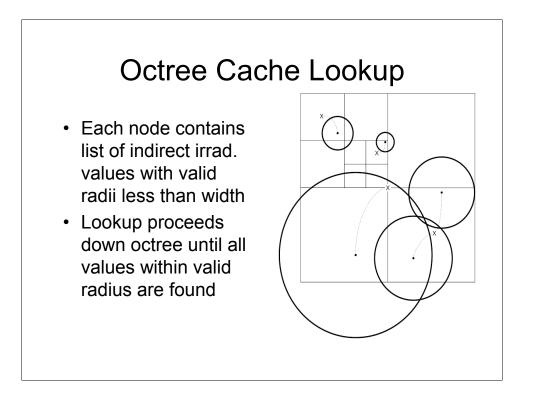
Shown here is the indirect (RGB) irradiance, which changes slowly over flat surface areas, and more rapidly over curved regions. Notice the saturation of red near the chairs, though their final color is not shown. Depicted is only the light arriving at surfaces after one or more bounces.



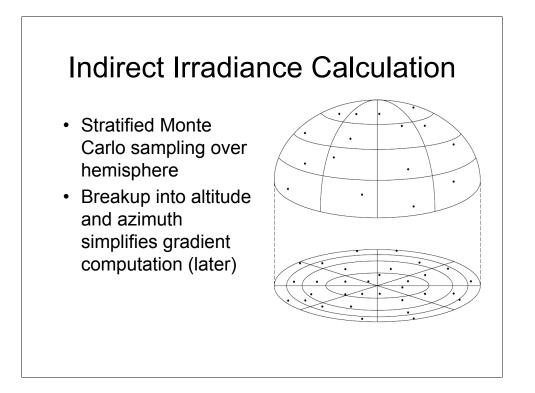
I didn't coin the term "irradiance caching" -- that was someone else's good idea. I just called it "lazy evaluation," being too lazy myself to think up a good name.



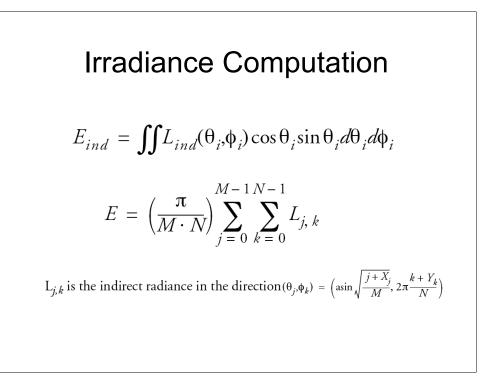
The E's are evaluation points. E1 and E2 shown here are previous evaluations, where E3 is a new evaluation triggered at query position C.



Using an octree that is independent of surfaces avoids placing restrictions on the geometric representation. It even works for volume data, sort of...



Stratified sampling reduces noise in the results at no extra cost, without adding bias. Divisions are placed to maintain constant projected areas on unit circle (Nusselt analog).



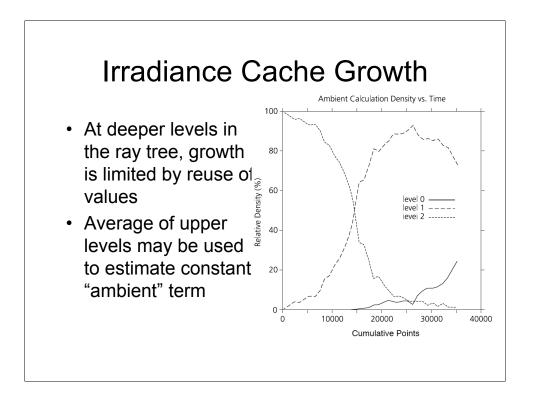
Integral equation may be converted to equal-weighted sum of selected "measures" using standard Monte Carlo integration techniques (inverting cosine projection). X_j and Y_k are uniformly distributed random variables between 0 and 1. To compute indirect irradiance, we count as zero any samples that intersect a light source.

Indirect Irradiance Example

- Send hemisphere ray samples
- Do not count direct sources
- Deeper levels use fewer samples
- Main *Radiance* controls:
 -ab number of bounces
 -ad, -as samples per hemisphere
 - -aa interpolation accuracy
 - -ar spatial resolution



The top image shows a hemispherical equal-contribution projection from the floor near one of the chairs in the conference room model. The bottom image shows the actual shape and number of samples sent for a particular indirect irradiance computation. Note that the sources are black, and the sampling areas follow altitude and azimuth divisions. Actual sample placement within each area is random.

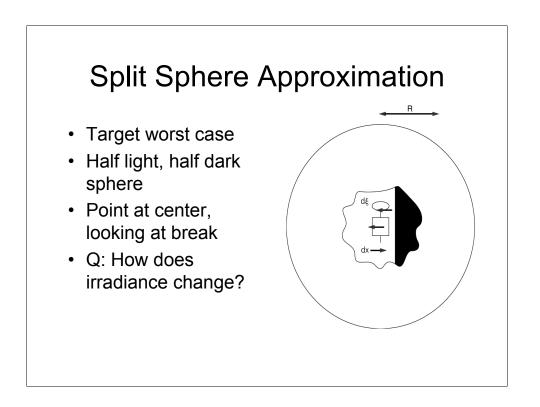


In a multi-bounce calculation, the deepest level is filled first, as the initial sample starts a hemisphere calculation, which begets another at the next level, and so on. The plot shows a three-bounce calculation, which starts by filling in the "level 2" cache, then spends time on the "level 1" cache before finally reaching the "level 0" cache -- the final indirect irradiance.

Irradiance Value Spacing & Weights

- Nearby geometry affects irradiance constancy (I.e., gradient)
- Surface curvature also influences
 gradient
- How can we quantify these to decide how closely to space our values?
- Can the same criteria be used to determine weights for interpolation?

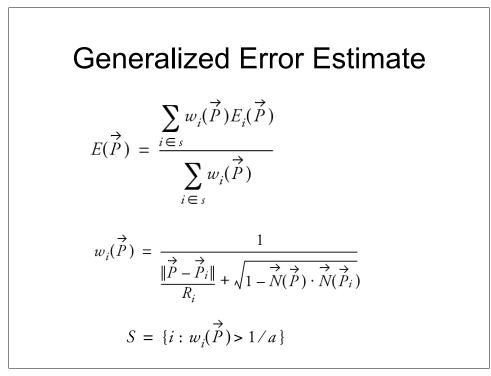
We have an intuition about the behavior of indirect irradiance -- we need to translate this into an algorithm that is robust and reliable.



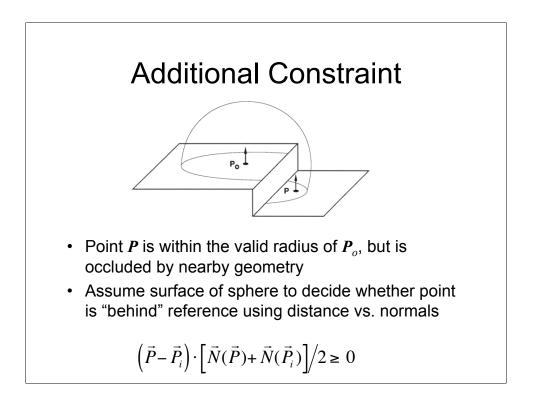
The "split sphere" is not technically a "worst case" -- it's just a "pretty bad case." It is better to qualify this as an assumption. I.e., we assume our actual environment will behave no worse than the split sphere. If it does, our technique won't necessarily break, but it won't be as accurate as we hoped, either.

 $\begin{aligned} & \mathcal{E} = \left| \frac{\partial E}{\partial x} (x - x_0) + \frac{\partial E}{\partial \xi} (\xi - \xi_0) \right| \\ & \varepsilon \leq \frac{4}{\pi} \frac{E}{R} |x - x_0| + E|\xi - \xi_0| \\ & \varepsilon(\vec{P}) \leq \frac{4}{\pi} E \frac{\|\vec{P} - \vec{P}_0\|}{R_0} + E\sqrt{2 - 2N(\vec{P}) \cdot N(\vec{P}_0)} \\ & R_0 = \text{average distance to surfaces at } \vec{P}_0 \end{aligned}$

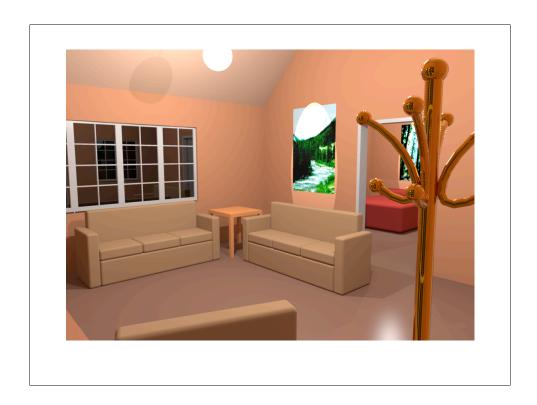
The first two inequalities bound the split sphere partial derivatives. The final equation uses these bounds to derive a "pretty bad case" error estimate for movement and rotation based on this first-order analysis. The calculation of R0 uses a harmonic mean to neighboring surfaces as measured by our ray samples. A harmonic mean is preferred since the radius appears in the denominator of our equation. Obviously, zero ray lengths are forbidden.



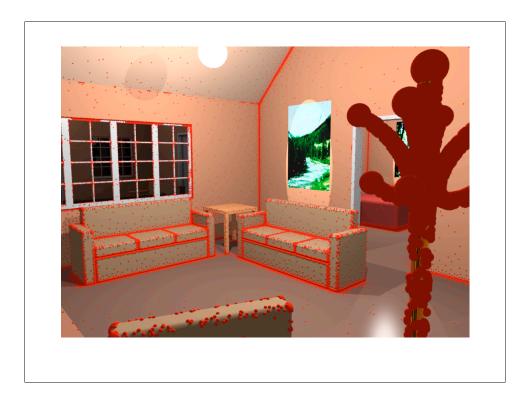
Ignoring the constants from our previous equation, we can derive a weighting function based on the split sphere approximation. Since our weights correlate to 1/error, applying them in a weighted average distributes error uniformly between the interpolated irradiance values. We hope to maintain final accuracy by restricting our set of values to ones whose estimated error is below some user tolerance, "a".



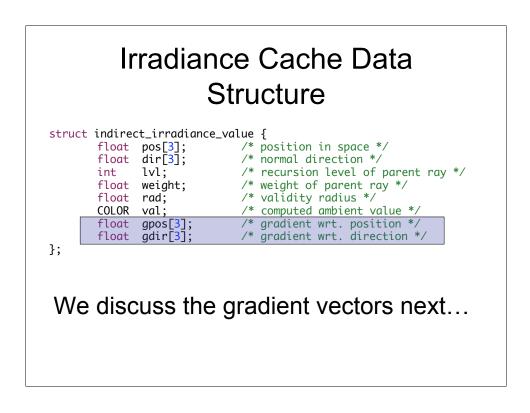
One special case we need to consider occurs when an irradiance value *thinks* it is valid over a larger region than it actually is. This case is depicted in the figure, where we are thinking about using an irradiance value at P0 at the new query position P. We can avoid this problem condition by adding a "behind test" to our criteria, which via the weighting function already incorporates a curvature test and a relative distance test.



Cabin model showing single-bounce calculation.



Indirect irradiance record placement shown as red spheres. Note how record spacing is large on flat surfaces far from neighboring geometry, then bunches up on outside curves and (especially) inside corners.

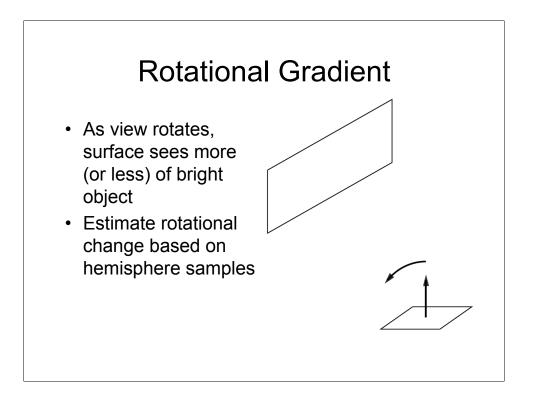


The ray level and weight are used to determine when to truncate the ray tree. The position and normal vectors are used in the weight/threshold computation, together with the "validity radius."

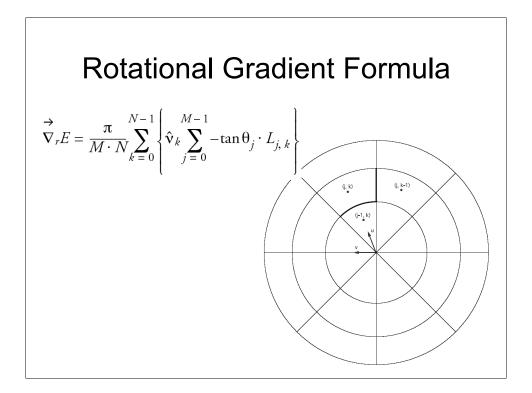
Irradiance Gradients

- From hemisphere sampling, we can also compute change w.r.t. position and direction
 - Gradient information comes essentially free
- Equivalent to higher-order interpolation method, i.e., cubic vs. linear

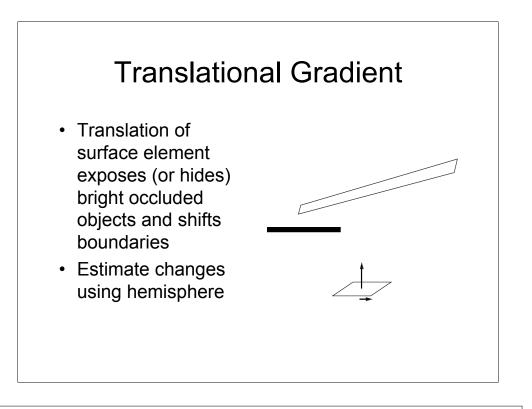
Initially, Paul Heckbert and I were thinking a gradient calculation could inform better value spacing. A year went by before we realized that it was better to apply them directly during interpolation to reduce discontinuities. This also avoids bias issues.



Caveat: we cannot know what will appear over the horizon.



The rotational gradient is zero at the zenith because the derivative of the cosine is zero, hence small rotations don't affect contribution for samples looking straight up.



The distance to our sampled geometry is important, and we have to assume that disocclusions look like what we already see of the background object. Hence, the boundary moves as the nearer edge moves rather than the further surface.

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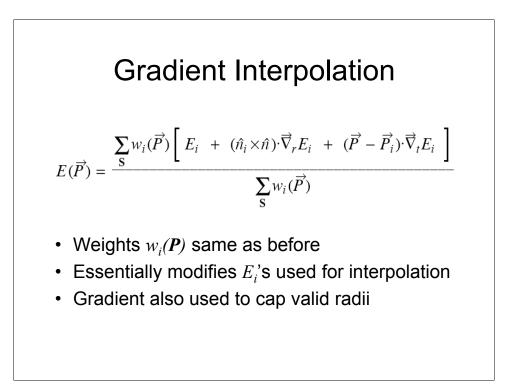
$$\vec{\nabla}_{t}E = \sum_{k=0}^{N-1} \left\{ \hat{u}_{k} \frac{2\pi}{N} \sum_{j=1}^{M-1} \frac{\sin\theta_{j_{-}} \cos^{2}\theta_{j_{-}}}{Min(r_{j,k}, r_{j-1,k})} \cdot (L_{j,k} - L_{j-1,k}) + \hat{v}_{k_{-}} \sum_{j=0}^{M-1} \frac{\sin\theta_{j_{+}} - \sin\theta_{j_{-}}}{Min(r_{j,k}, r_{j,k-1})} \cdot (L_{j,k} - L_{j,k-1}) \right\}$$

where:

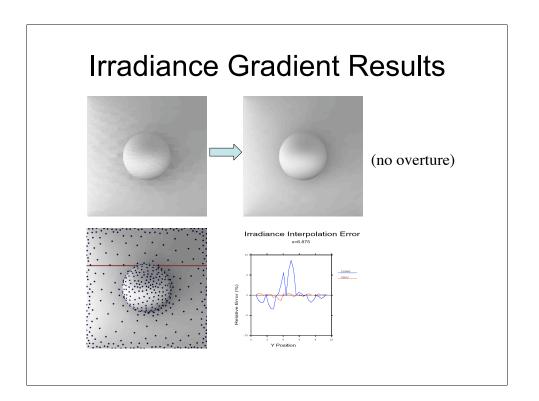
 \hat{u}_k is the unit vector in the ϕ_k direction

- $\hat{v}_{k_{-}}$ is the unit vector in the $\phi_{k_{-}} + \frac{\pi}{2}$ direction
- $\theta_{j_{-}}$ is the polar angle at the previous boundary, $\sin^{-1}\sqrt{\frac{j}{M}}$
- $\theta_{j_{+}}$ is the polar angle at the next boundary, $\sin^{-1}\sqrt{\frac{j+1}{M}}$
- $\phi_{k_{-}}$ is the azimuthal angle at the previous boundary, $2\pi \frac{k}{\lambda_{l}}$
- $r_{j,k}$ is the intersection distance for cell (j,k)

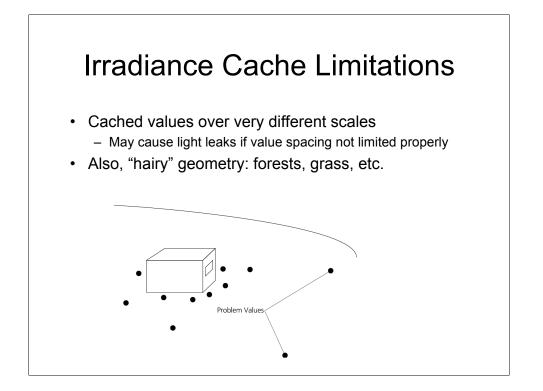
Refer to the vector diagram on the rotational gradient slide for variable definitions. Here, the '+' and '-' suffixes refer to the vectors (not drawn) corresponding to the leading and trailing edges of each cell, respectively. The Min(a,b) function is used to determine the closer of two neighboring surface samples.



Once we have our rotational and translational gradients corresponding to each irradiance value, we can apply them in a first-order, weighted interpolation as shown.



The effect of applying the irradiance gradient technique is highlighted in this extrapolated calculation, where lazy evaluation is shown in its "full glory." Despite discontinuities introduced by adding values to the cache during scanline rendering, our first-order extrapolation makes these artifacts all but disappear. Value placement is shown together with the comparative error when applying gradients to interpolation (I.e., adding an overture calculation).



Jaroslav has a solution called "neighbor clamping" for this problem. Hairy geometry is best dealt with by switching to a noisy MCPT method for busy topologies.

Sources of Bias in Irradiance Cache

- Super-sampling of hemisphere
 - Naïve adaptive sampling approach is biased
- · Truncation bias from limited bounces
 - Caching overrides Russian roulette in *Radiance*
- Placing limits on value spacing
 - Degrades fine-scale features
- Assumed average scene reflectance
 - Undermines accuracy in white-walled enclosure

To eliminate bias, don't use adaptive super-sampling, add Russian roulette to final bounce via path tracing, eliminate minimum value spacing (expensive) and use actual surface reflectances (undermines sharing). Decreasing number of hemisphere samples (by reflectance) and increasing sample spacing (by reflectance^-0.5) distributes errors evenly between levels and speeds convergence.