

Anti-aliasing and sampling

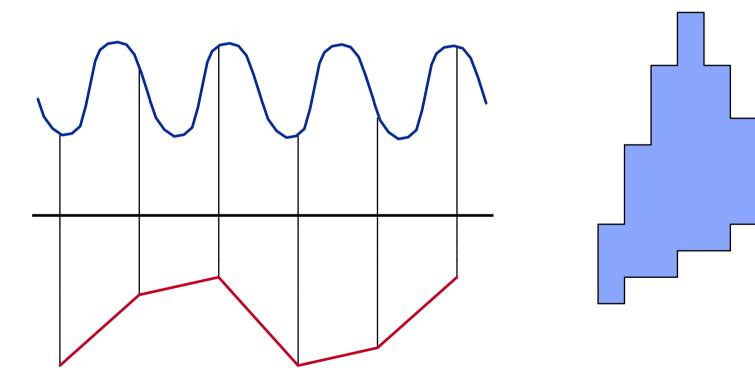
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Artefacts caused by a regular-grid or regular sampling



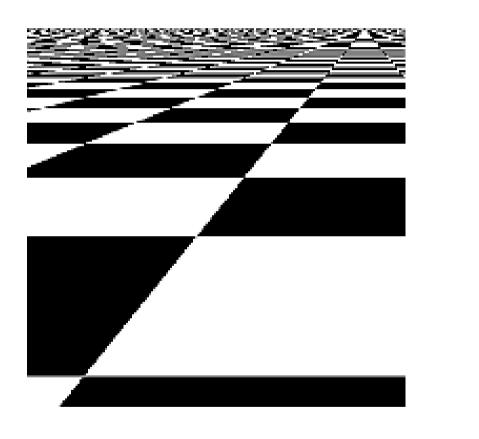
Spatial alias



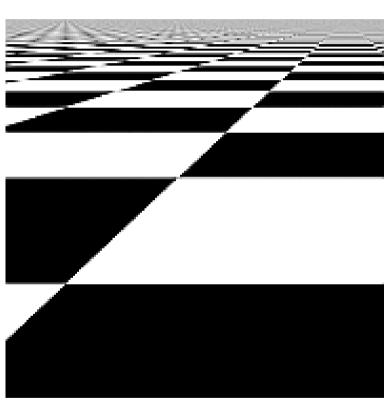
- jagged oblique lines
 - regular dense system of lines or stripes on a texture can lead to "Moiré effect"
- interference of fast periodic image changes with a pixel raster
 - example: picket fence in perspective projection
 - too fine or too distant regular texture (checkerboard viewed from distance)







1 sample per pixel



256 spp (jittering)

Temporal alias



- shows in slow motion animation
- blinking pixels on contours of moving objects
 - the whole small objects can blink
- interference of a periodic movement with a frame frequency
 - spinning wheel seems to be still or even rotating the other direction

Real world



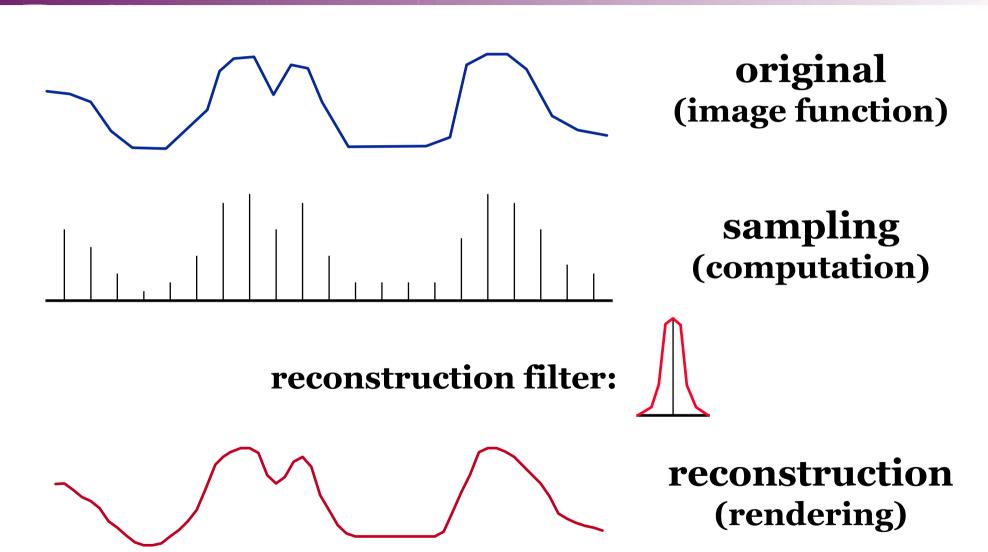
Human visual system has no alias.

Alias manifests only mildly in photography.

- objects smaller than resolution of a sensor are without details / blurry
 - fence from large distance is percieved as an average-color area (mix of background + foreground colors)
- too fast movement generates fuzzy (blurred) perception



Reconstruction in raster context



Sampling and reconstruction



- image sampling or computing of image function
 - higher frequencies should be reduced/removed from an image before sampling
 - low pass filter (convolution window averaging)
 - image synthesis can reduce higher frequencies directly (anti-aliasing by pixel supersampling)
- reconstruction filter is defined by an output device
 - e.g. neighbour CRT monitor pixels overlap
 - LCD pixels behave differently





Image function with continuous domain and unlimited spectrum:

f(x, y)

Anti-aliasing filter (function with limited support):

h(x, y)

Pixel color [i,j]:

$$\mathbf{I}(\mathbf{i},\mathbf{j}) = \int_{-\infty-\infty}^{\infty} \int_{-\infty}^{\infty} \mathbf{f}(\mathbf{x},\mathbf{y}) \cdot \mathbf{h}(\mathbf{x} - \mathbf{i},\mathbf{y} - \mathbf{j}) \, d\mathbf{x} \, d\mathbf{y}$$





Assuming box smoothing filter and unit square pixel:

$$(i, j) = \int_{j}^{j+1} \int_{i}^{i+1} f(x, y) dx dy$$

(integral average value of the image function on the pixel area)

Quadrature



- analytic (close form solution)
 - in rare cases (simple image function)
- numeric solution using sampling
 - finite set of samples $[x_i,y_i]$
 - integral estimate by the sum

$$\mathbf{I}(\mathbf{i}, \mathbf{j}) = \frac{\sum_{k} \mathbf{f}(\mathbf{x}_{k}, \mathbf{y}_{k}) \cdot \mathbf{h}(\mathbf{x}_{k}, \mathbf{y}_{k})}{\sum_{k} \mathbf{h}(\mathbf{x}_{k}, \mathbf{y}_{k})}$$

stochastic sampling – Monte-Carlo method

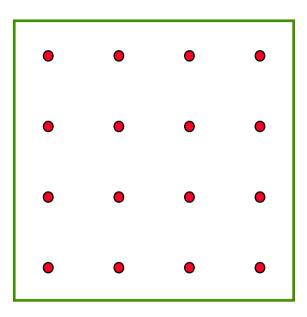
Sampling methods



- mapping: $k \rightarrow [x_k, y_k]$
 - sample is selected from the given 2D region (domain):
 usually rectangular, square or circular
 - sampling in higher dimensions (dim=10)
- required **properties** of sampling algorithms
 - uniform probability over a domain
 - high regularity is not desirable (interference)
 - efficient computation



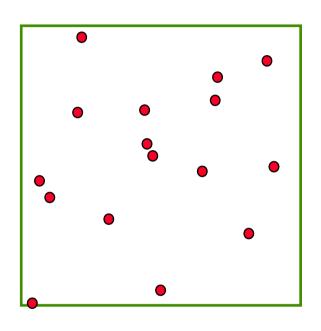




Does not reduce **Moire / interference** (interference still appears in higher frequencies)





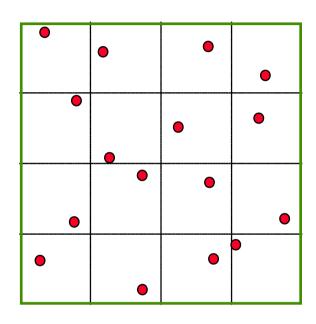


N independent random samples with uniform probability density (PDF)

Samples tend to form **clusters Too much noise** in a result image

"Jittering"

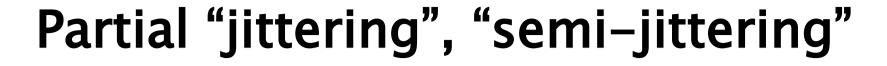




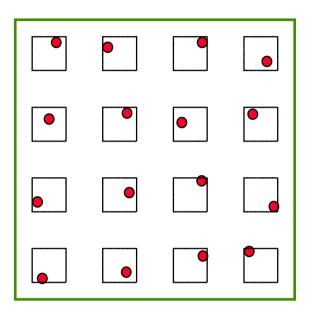
K × K <u>independent</u> random samples in K × K <u>equally sized</u> sub-regions covering the original domain completely

Big clusters are not possible

More regular coverage of a domain





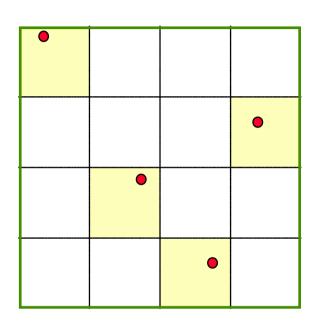


K × K <u>independent</u> random samples in K × K <u>equally sized</u> sub-regions <u>not covering</u> the original domain

Clusters are impossible **Too** regular (interference)







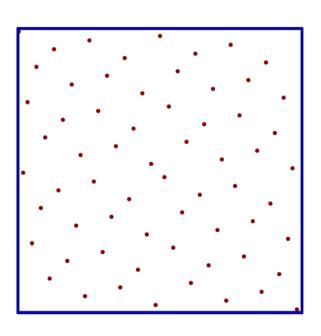
"Low-cost jittering":

there is exactly one sample in each row and in each column. Random permutation of a diagonal

Good properties of "jittering" are preserved Higher **efficiency** (especially in high dimensions)

Hammersley





- + excellent discrepancy
- + deterministic
- + very fast
- difficult adaptive refinement
- bad spatial spectrum

Famous **Halton sequence** is based on similar principles..

Deterministic sequences



- based on similar principles:
 - Halton, Hammersley, Larcher-Pillichshammer
- for a prime number **b** let **n** be positive integer expressed using b-representation:

$$n = \sum_{k=0}^{L-1} d_k(n)b^k$$

• then there is a number from [0,1) range:

$$g_b(n) = \sum_{k=0}^{L-1} d_k(n)b^{-k-1}$$

Halton, Hammersley



• famous Halton sequence (e.g. $b_1=2$, $b_2=3$):

$$x(n) = [g_{b_1}(n), g_{b_2}(n)]$$

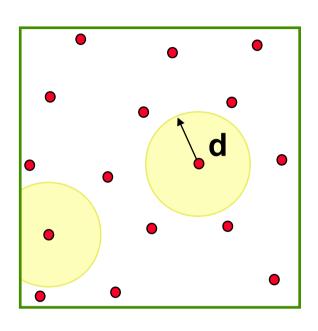
• Hammersley sequence (e.g. b=2):

$$x(n) = \left[\frac{n}{N}, g_b(n)\right]$$

Larcher-Pillichshammer sequence uses **XOR** operation instead of addition (inside the g_b(n))...







N random samples meeting condition: $|[x_k,y_k] - [x_l,y_l]| > d$ for given value d

Prevents creating **clusters**, imitates **distribution of light-perception cells** in retina of a mammal. Difficult efficient **implementation**!

Implementation



rejection sampling

- candidate sample is rejected if too close to any previous accepted sample
- less efficient for higher number of samples
- choice of value d is problematic
 - maximum number of placeable samples depends on **d**
- difficult (adaptive) refinement
 - additional samples to a existing set of samples





("best candidate" algorithm)

- generates gradually refined sample set (from Poisson sampling)
 - no problems with d
 - intrinsic refinement
- compute-intensive algorithm
 - sample set can be **precomputed and reused**
 - to reduce dependency between neighbour pixels random rotation and translation can be used

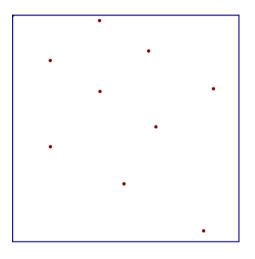
Mitchell's algorithm

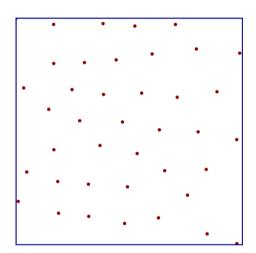


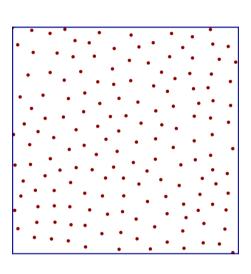
- the 1st sample is chosen randomly
- ² choice of the **(k+1)**th sample:
 - generate k· q independent candidates (q determines sample-set quality)
 - the most distant candidate (from all previous k accepted samples) is selected and accepted
- for higher **q** we get better quality set
 - choose q > 10 in demanding situations

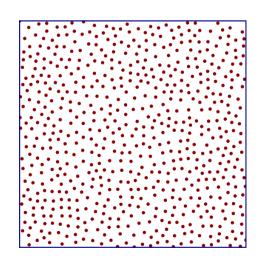
Incremental example

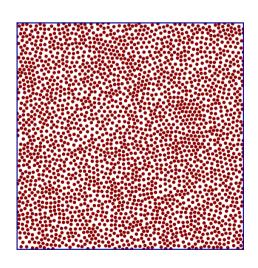












Sample numbers:

$$K = 10$$

Adaptive refinement



- sampling based on local importance (importance sampling) or interest
 - some regions have higher weight (higher probability)
 - regions with **higher variance** should be sampled more densely
- "importance" or "interest" need not be known in advance (explicitly)
 - algorithm has to adopt to intermediate results (adaptability)

Modification of static methods



• initial phase:

- compute small set of test samples (1-5)
- define refinement criterion based on previous samples

refinement phase:

- sampling is refined in regions of higher need (criterion)
- efficiency: if we can reuse all generated samples
- almost every sampling algorithm can be reformulated in that way

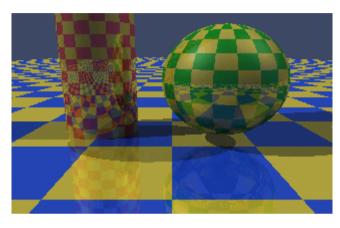
Refinement criteria



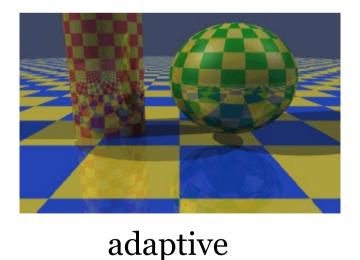
- function values (difference, variance, gradient)
 - difference between neighbour samples, ..
- Id's of hit solids (Ray-tracing specific)
 - higher priority
 - textures with repeated patterns use of <u>signatures</u>
- trace tree (recursive ray-tracing)
 - topologic comparison of complete or limited trace trees
 - tree identifier recursive hash function using solid Id's, texture signatures, shadow/light, ..



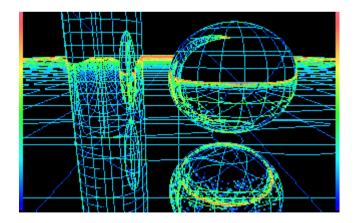




1 spp



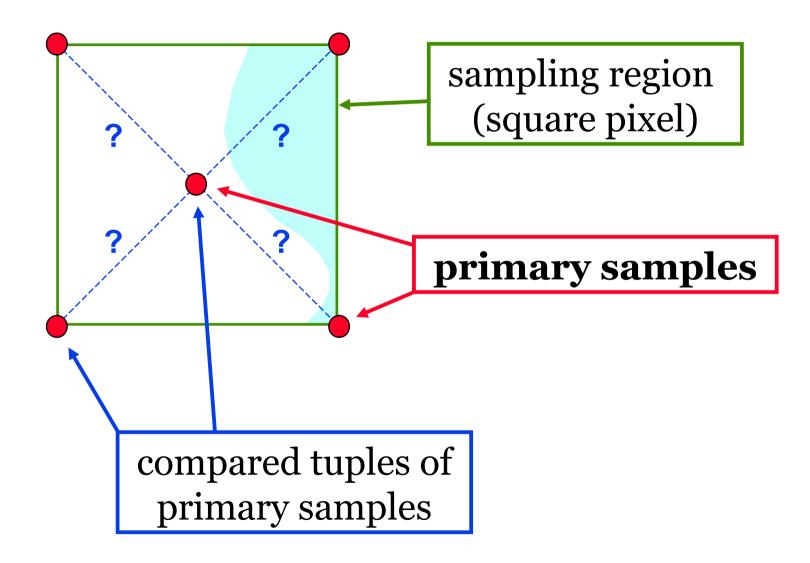
1/2 spp



refinement map

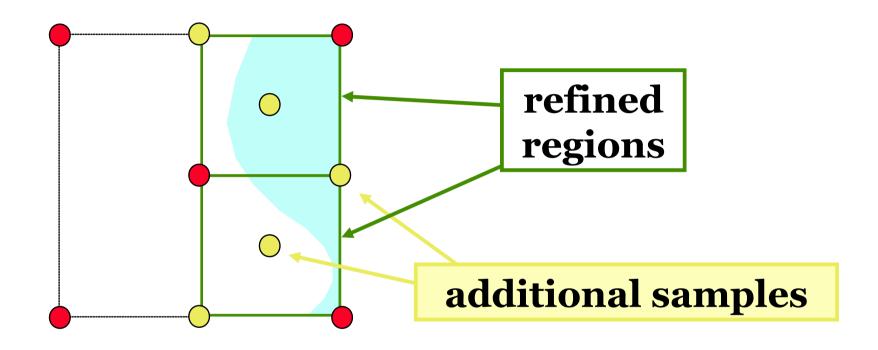
Recursive refinement (Whitted)







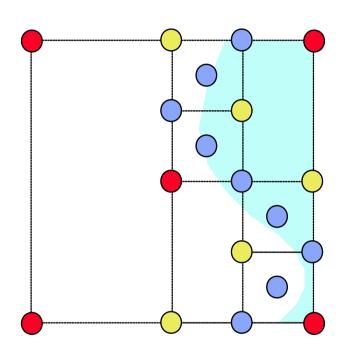




The same procedure is executed in refined regions **recursively** (up to the declared maximum level)





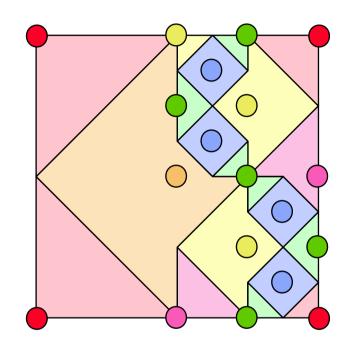


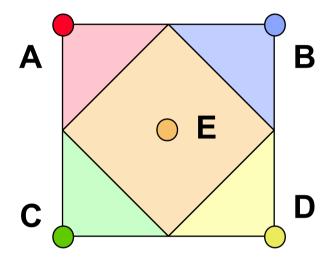
- phase I
- phase II
- phase III

Evaluated: 5+5+9 = 19 samples (from total number of 41)









$$\frac{1}{2}E + \frac{1}{8}[A + B + C + D]$$

If the refinement stops in a specific square, its area is split to two triangles (diagonal samples)

References



- A. Glassner: An Introduction to Ray Tracing, Academic Press, London 1989, 161-171
- A. Glassner: Principles of Digital Image Synthesis, Morgan Kaufmann, 1995, 299-540
- J. Pelikán: Náhodné rozmisťování bodů v rovině (Random point placement), CSGG 2014, slides & paper available online